

OPTIMAL POWER FLOW STUDIES

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THESIS

In recent years a great deal of work has been done on the optimal power flow studies making use of nonlinear programming formulation with full a.c. simulation of the network. Important advances made include application of various optimization techniques in conjunction with efficient computation methods like triangular factorization, compact storage and optimal renumbering. These methods recently introduced to power system engineers by J.R. Finney and his associates of Bonneville Power Administration^{1,45} have made possible the study of very large power systems without partitioning and resulted in considerable saving of computer time. Chapter II discusses mainly these.

In Chapter III algorithm of forming the admittance matrix and simultaneously storing the non zero elements in the compact form is given. Power flow equations are discussed in Chapter IV.

In Chapter V various available load flow techniques are lumped into general categories and compared through fixed point analysis. To date not much use has been made of such analysis to power system studies. The author is familiar with only one paper³ making its use. This paper mainly discusses Newton's method through Kantrovich's theorem which makes use of Jacobian and the three dimensional Hessian matrix, computation of which is very hard. By making use of fixed point analysis, convergence properties of various

algorithms which up to now were experimentally established are analytically established.

A discussion of computer programme for load flow solution by Newton's method is given in Chapter VI. In Chapter VII criterie and techniques for identifying the optimal point of the nonlinear programming problem and convergence of sequential optimization through penalty and barrier methods are discussed.

In Chapter VIII various published techniques for optimal flow solutions are discussed and some are studied through fixed point analysis, which indicate the doubtful convergence of many of these. Dommel and Tinney's²¹ technique is found to be the most promising. In Chapter IX the computer programme for optimal power flow problem making use of a method which is a modification of the above is presented. It uses a modified Lagrangian, a different gradient and penalty method and handles the inequality constraints on voltages for the buses with reactive power generation on different line. In Chapter X algorithms developed by the author for optimal hydrothermal operations are presented.

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Definitions

Following is the list of more commonly used symbols.

In rare cases, some of these symbols represent a different quantity, explained in the text.

| | |
|------------------------------------|---|
| $v_i L_i = e_i + j f_i$ | = complex voltage of node i |
| I_i | = nodal current |
| $y_{ij} L_{ij} = g_{ij} + jb_{ij}$ | = element of admittance matrix |
| V | = column vector of bus voltages |
| I | = column vector of bus currents |
| Y | = admittance matrix |
| $P_i + j Q_i$ | = net injected power at node i |
| $P_i^G + j Q_i^G$ | = net generation at node i |
| $P_i^L + j Q_i^L$ | = net load at node i |
| \bar{x} | = upper limit of x |
| \underline{x} | = lower limit of x |
| M, A | = Jacobian matrix |
| n | = number of nodes |
| r | = step size |
| R^k | = constant with the sequential minimization |
| t_h | = off nominal turn ratio of hth transformer |
| ϕ_h | = angular shift of hth phase shifting transformer |
| x^* | = the required value of x |
| P_L | = system transmission losses |
| P_D | = system load demand |

- subscript s** - refers to the slack bus
- j** - set of nodes with active power generations
- R** - set of nodes with reactive power generations
- t** - set of nodes with thermal generations
- H** - set of nodes with hydrogenerations.
- $S_1 + S_2$ - union of sets S_1 and S_2
- $S_1 \cap S_2$ - intersection of sets S_1 and S_2

Chapter I

INTRODUCTION

Optimal power flow studies were made as early as 1940's when the transmission losses were neglected and the generation powers were scheduled on the basis of equal incremental costs. With the development of interconnected power systems, need for consideration of these losses was felt. These losses initially were approximated by a quadratic loss formula. Further work was aimed at the improvement in the method of obtaining the loss formula and generator scheduling initially on network analyser and later on digital computer. This method optimises for real powers only. Limitations of the loss formula led to search for better methods. A partial break through was first achieved in 1962 by Carpentier of Electricite' de France who made use of non-linear programming formulation through Kuhn Tucker's theorem with full a.c. representation of the network, however the algorithm needed further improvements. Further work in this direction published in English speaking literature was carried out at Bonneville Power Administration and Stanford Research Institute for the BPA system which has 100% hydrogeneration, for which transmission losses are to be minimised by reactive powers and tap adjustments. Carpentier's algorithm was not found suitable for the system. BPA had meanwhile perfected the computer programme for load flow studies and published in 1967. Making use of this programme a method for reactive

power optimization was suggested^{19, 36}. A lot of work meanwhile in this direction was also reportedly carried out in U.S.A. most of which is not yet available in English literature. Thereafter a number of papers in this direction have appeared, most claiming some advantages over other methods. An entirely different approach making use of penalty functions was developed by Sesson at Imperial College, London, and Instituto Technologico y de Estudios Superiores de Monterrey, Mexico, and published in 1969-1970^{4, 15, 16}.

The author's aim was to make a comparative study of the existing published methods and to develop an efficient and reliable computer programme suitable for large power systems. This needs that the sparsity of system matrices be fully exploited. Application of matrix inverse is out of question for larger systems. Fortunately the technique of triangular factorization recently introduced to power system engineers replaces the use of matrix inverse. This along with compact storage for matrices and optimal renumbering of nodes makes the computer logic complicated.

Performance of any technique is better judged from a bigger system. Because of availability of small digital computer IBM 1130 with 16 K, 16 bit storage, the study was restricted to 30 buses only. Findings of Dommel and Tinney indicate that conclusions drawn on a system of 25 buses or so usually hold good for larger system as well¹⁴.

After some experimentation over the existing published techniques for load flow solutions, Newton's method as

perfected by Tinney and Hart³³ with some modifications, was finally adopted. Chapter VI, discussing the author's programme for load flow studies is intended as a supplement to this reference. The accuracy of this method was found restricted only by round off errors. Though a tolerance of 10^{-4} pu was prescribed, mismatch of no t buses was found much less than this at the end of fourth iteration. A need for very accurate load flow programme was felt for the optimal power flow studies in the absence of which the relative costs with different schedule do not give a realistic comparison since the improvement in the operating cost is expected between 1 to 3% only with optimal scheduling. Improvement in transmission loss reported by Feschen et al. is about 4% with reactive power scheduling³⁴. A comparison of various published algorithms for optimal power flow studies in Chapter VIII indicates the supremacy of Domel and Tinney's method over others. While trying this method for the system described in Section 9.8, a need for further improvement was felt of which the details are given in Chapter IX. Application of the optimal flow programme apart from the routine fuel cost minimization was also made for planning power requirements of power systems under constraints on line voltages in Section 9.10.

In Chapter 10 algorithms based on Lagrangian formulation and gradient techniques for optimal hydrothermal scheduling of system consisting of multireservoirs (with and without head variation) and multithermal plants are given. Initially the algorithm is based on Bryson and Denham's technique⁴⁸ for which

the solution remains in the feasible domain of power flow equations and water storage. Application of decomposition techniques, comparative advantages and disadvantages are also given. It was found that application of maximum principle and Lagrangian formulation essentially give the same mathematical equations, on which the algorithm is to be based.



CHAPTER XI

SOLUTION OF LINEAR EQUATIONS

This chapter presents some of the recently introduced techniques namely triangular factorization, compact storage scheme and optimal renumbering scheme, for solving a set of linear equations. Section 2.1 presents triangular factorization technique, in which an array of numbers, is obtained from a non singular matrix A ; that can be used to obtain the effect of A , A^{-1} , A^t , $(A^{-1})^t$ over a column vector. This technique replaces the use of matrix inverse, which is very inefficient for large sparse system of equations such as occur in power network problems, since while the system matrix is sparse, its inverse is a full matrix. Most material of this section is from reference (31). However in addition a few computer logics and few additional proofs which include 2.42, 2.51 and 2.54 have been added. Few of the compact storage schemes are described in section 2.2. Near optimal renumbering schemes are discussed in section 2.3. Description of various numbering schemes is from reference (1). In addition description and computer logic of authors sub-routine is provided. These two techniques greatly improve the efficiency of traingular factorization.

2.1 Triangular Factorization

A matrix is usually triangularized by eliminating

elements of successive columns below the diagonal. From the point of computer speed and storage, it is more efficient to eliminate elements of each row up to the diagonal before proceeding to the next row, since only one row need be formed and operated upon at one time.

Further discussion is based on the following equation

$$Ax = b \quad (2.1)$$

where A is $n \times n$ non singular matrix, x is a column vector of unknowns and b is the known vector.

This equation written explicitly is as follows -

$$\begin{bmatrix} a_{11} & a_{21} & \dots & a_{1n} \\ a_{21} & a_{22} & & a_{2n} \\ \vdots & \vdots & \ddots & \vdots \\ a_{n1} & a_{n2} & \dots & a_{nn} \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ \vdots \\ \vdots \\ x_n \end{bmatrix} = \begin{bmatrix} b_1 \\ b_2 \\ \vdots \\ \vdots \\ \vdots \\ b_n \end{bmatrix} \quad (2.2)$$

The first step in Gaussian elimination is to divide the elements of the first row by a_{11} as follows:

$$d_{11} = 1/a_{11} \quad (2.3)$$

$$(1) \quad a_{1j} = d_{11} a_{1j}; \quad j = 1, \dots, n \quad (2.4)$$

$$u_{1j} = a_{1j}^{(1)}; \quad j = 2, \dots, n \quad (2.5)$$

$$b_1^{(1)} = d_{11} b_1 \quad (2.6)$$

a_{21} is eliminated from the second row by linear combination with the derived first row, and then to divide the remaining derived elements of the second row by its derived diagonal element as follows:

$$l_{21} = a_{21} \quad (2.7)$$

$$a_{2j}^{(1)} = a_{2j} - l_{21} a_{1j}^{(1)} ; \quad j = 1, \dots, n \quad (2.8)$$

$$b_2^{(1)} = b_2 - l_{21} b_1^{(1)} \quad (2.9)$$

$$d_{22} = 1/a_{22}^{(1)} \quad (2.10)$$

$$a_{2j}^{(2)} = d_{22} a_{2j}^{(1)} ; \quad j = 2, \dots, n \quad (2.11)$$

$$u_{2j} = a_{2j}^{(2)} ; \quad j = 3, \dots, n \quad (2.12)$$

$$b_2^{(2)} = d_{22} b_2^{(1)} \quad (2.13)$$

a_{31} and a_{32} are eliminated from third row and remaining elements of this row are now divided by the diagonal element.

Thus elements to the left of diagonal are eliminated and the diagonal element is made unity. This process illustrated for k th row is as follows.

$$l_{k1} = a_{k1} \quad (2.14)$$

$$a_{kj}^{(1)} = a_{kj} - l_{k1} a_{1j}^{(1)} ; \quad j = 1, \dots, n \quad (2.15)$$

$$b_k^{(1)} = b_k - l_{k1} b_1^{(1)} \quad (2.16)$$

$$l_{k2} = a_{k2}^{(1)} \quad (2.17)$$

$$a_{kj}^{(2)} = a_{kj}^{(1)} - l_{k2} a_{2j}^{(1)} ; \quad j = 2, \dots, n \quad (2.18)$$

$$b_k^{(2)} = b_k^{(1)} - l_{k2} b_2^{(1)} \quad (2.19)$$

$$\begin{matrix} \cdot \\ \vdots \\ \cdot \end{matrix}$$

$$l_{k,k-1} = a_{k,k-1}^{(1,-2)} \quad (2.20)$$

$$a_{kj}^{(k-1)} = a_{kj}^{(k-2)} - l_{k,k-1} a_{k-1,j}^{(k-1)} ; \quad j = k-1, \dots, n \quad (2.21)$$

$$b_k^{(k-1)} = b_k^{(k-2)} - l_{k,k-1} b_{k-1}^{(k-1)} \quad (2.22)$$

$$d_{kk} = 1/a_{kk}^{(k-1)} \quad (2.23)$$

$$a_{kj}^{(k)} = d_{kk} a_{kj}^{(k-1)} ; \quad j \geq k \quad (2.24)$$

$$b_k^{(k)} = d_{kk} b_k^{(k-1)} \quad (2.25)$$

$$a_{kj} = a_{kj}^{(k)} ; \quad j > k \quad (2.26)$$

This process represented by equations 2.15 to 2.22 amounts to premultiplying the derived matrix A' and b' by a $n \times n$ non-singular matrix L_k which differ from the unit matrix only in the kth row as follows:

$$\text{matrix } L_k ; \text{ row } k = [-1_{k,1}, -1_{k,2}, \dots, -1_{k,k-1}, 1, 0, \dots, 0] \quad (2.27)$$

The process represented by equations 2.24 and 2.25 amounts to premultiplying the derived matrix A' and b' by an elementary matrix D_k which differs from the unit matrix in the k th diagonal element as follows:

$$\text{matrix } D_k ; \text{ row } k = [0, \dots, 0, d_{kk}, 0, \dots, 0] \quad (2.28)$$

After all the elements on the left side of the diagonal are eliminated and all the diagonal element made unity; the derived system of equation is as follows:

$$\left[\begin{array}{cccc|c} 1 & u_{12} & \dots & u_{1n} & x_1 \\ 0 & 1 & \dots & u_{2n} & x_2 \\ \vdots & \vdots & \ddots & \vdots & \vdots \\ 0 & \dots & 1 & u_{nn} & x_n \end{array} \right] = \left[\begin{array}{c} b_1^{(1)} \\ b_2^{(2)} \\ \vdots \\ b_n^{(n)} \end{array} \right] \quad (2.29)$$

Solution is now obtained as follows by back substitution.

$$x_n = b_n^{(n)}$$

$$x_{n-1} = b_n^{(n-1)} - u_{n-1,n}^{(n-1)} x_n$$

$$x_1 = b_1^{(1)} - \sum_{j>1} u_{1j} x_j \quad (2.30)$$

Back substitution amounts to premultiplying the derived matrix b' successively by U_{n-1} , U_{n-2} , ..., U_2 , U_1 , where U_k is a $n \times n$ non-singular matrix differing from the matrix as follows.

$$U_k ; \text{ row } k = [0, \dots, 0, 1, -u_{k,k+1}, \dots, -u_{kn}] \quad (2.31)$$

Thus

$$U_1 U_2 \cdots U_{n-1} D_n L_n \cdots L_2 L_2 D_1 = b^{-1} \quad (2.32)$$

The non trivial element of the matrices U , L and D are stored as follows:

$$\begin{bmatrix} d_{11} & u_{12} & u_{13} & \cdots & u_{1n} \\ l_{21} & d_{22} & u_{23} & \cdots & u_{2n} \\ l_{31} & l_{32} & d_{33} & \cdots & u_{3n} \\ \vdots & \vdots & \vdots & & \vdots \\ \vdots & \vdots & \vdots & & \vdots \\ \vdots & \vdots & \vdots & & \vdots \\ l_{n1} & l_{n2} & & & d_{nn} \end{bmatrix} \quad (2.33)$$

This table will henceforth be called as table of L. U. factors.

Inverse of these matrices are trivial. Inverse of D_1

matrix involves only the reciprocal of the element a_{11} . The inverses of matrix L_1 and U_1 involve only reversal of algebraic sign of the off diagonal elements.

Operations A^{-1} , \therefore , $(At)^{-1}$, (A^t) on a column vector are represented as follows:

$$A^{-1} b = U_1 U_2 \cdots U_{n-1} D_n L_n \cdots L_2 L_1 b \quad (2.34)$$

$$A b = L_1^{-1} U_2^{-1} D_2 \cdots L_n^{-1} U_n^{-1} U_{n-1}^{-1} \cdots U_2^{-1} U_1^{-1} b \quad (2.35)$$

$$(At)^{-1} b = L_1^t L_2^t D_2^t \cdots L_n^t U_n^t U_{n-1}^t \cdots U_2^t U_1^t b \quad (2.36)$$

$$(A^t) b = (U_1^{-1})^t (U_2^{-1})^t \cdots (U_{n-1}^{-1})^t (D_n^{-1})^t (L_n^{-1})^t \\ \cdots (D_2^{-1})^t (L_2^{-1})^t (L_1^{-1})^t b \quad (2.37)$$

2.1.1 Numerical example

The following example is based on the non symmetrical matrix A

$$A = \begin{bmatrix} 2 & 4 & 6 \\ 3 & 8 & 13 \\ 4 & 10 & 19 \end{bmatrix}$$

Table of L.U. factors for A is obtained as follows.

$$\left[\begin{array}{ccc} 2 & 4 & 6 \\ 3 & 8 & 13 \\ 4 & 10 & 19 \end{array} \right] \xrightarrow{\quad} \left[\begin{array}{ccc} \frac{1}{2} & 2 & 3 \\ 3 & 8 & 13 \\ 4 & 10 & 19 \end{array} \right] \xrightarrow{\quad} \left[\begin{array}{ccc} \frac{1}{2} & 2 & 3 \\ 3 & 2 & 4 \\ 4 & 10 & 19 \end{array} \right] \xrightarrow{\quad} \left[\begin{array}{ccc} \frac{1}{2} & 2 & 3 \\ 3 & \frac{1}{2} & 2 \\ 4 & 10 & 19 \end{array} \right]$$

$$\left[\begin{array}{ccc} \frac{1}{2} & 2 & 3 \\ 3 & \frac{1}{2} & 2 \\ 4 & 2 & 7 \end{array} \right] \xrightarrow{\quad} \left[\begin{array}{ccc} \frac{1}{2} & 2 & 3 \\ 3 & \frac{1}{2} & 2 \\ 4 & 2 & 3 \end{array} \right] \xrightarrow{\quad} \left[\begin{array}{ccc} \frac{1}{2} & 2 & 3 \\ 3 & \frac{1}{2} & 2 \\ 4 & 2 & \frac{1}{3} \end{array} \right]$$

With $b = [20 \quad 38 \quad 51]^t$; x is obtained from table of factors as follows:

$$\left[\begin{array}{c} 20 \\ 38 \\ 51 \end{array} \right] \xrightarrow{\quad} \left[\begin{array}{c} 10 \\ 38 \\ 51 \end{array} \right] \xrightarrow{\quad} \left[\begin{array}{c} 10 \\ 8 \\ 51 \end{array} \right] \xrightarrow{\quad} \left[\begin{array}{c} 10 \\ 4 \\ 51 \end{array} \right] \xrightarrow{\quad} \left[\begin{array}{c} 10 \\ 4 \\ 3 \end{array} \right] \xrightarrow{\quad} \left[\begin{array}{c} 10 \\ 4 \\ 1 \end{array} \right] \xrightarrow{\quad} \left[\begin{array}{c} 10 \\ 2 \\ 1 \end{array} \right] \xrightarrow{\quad} \left[\begin{array}{c} 3 \\ 2 \\ 1 \end{array} \right]$$

With $x = [3 \quad 2 \quad 1]^t$; $b = Ax$ can be obtained from table of factors as follows:

$$\left[\begin{array}{c} 3 \\ 2 \\ 1 \end{array} \right] \xrightarrow{\quad} \left[\begin{array}{c} 10 \\ 2 \\ 1 \end{array} \right] \xrightarrow{\quad} \left[\begin{array}{c} 10 \\ 4 \\ 1 \end{array} \right] \xrightarrow{\quad} \left[\begin{array}{c} 10 \\ 4 \\ 3 \end{array} \right] \xrightarrow{\quad} \left[\begin{array}{c} 10 \\ 4 \\ 51 \end{array} \right] \xrightarrow{\quad} \left[\begin{array}{c} 10 \\ 8 \\ 51 \end{array} \right] \xrightarrow{\quad} \left[\begin{array}{c} 10 \\ 38 \\ 51 \end{array} \right] \xrightarrow{\quad} \left[\begin{array}{c} 10 \\ 38 \\ 51 \end{array} \right]$$

With $b = \begin{bmatrix} 16 \\ 38 \\ 63 \end{bmatrix}^T$; $(A^{-1})^T b$ is obtained as follows:

$$\begin{bmatrix} 16 \\ 38 \\ 63 \end{bmatrix} \xrightarrow{\quad} \begin{bmatrix} 16 \\ 6 \\ 15 \end{bmatrix} \xrightarrow{\quad} \begin{bmatrix} 16 \\ 6 \\ 3 \end{bmatrix} \xrightarrow{\quad} \begin{bmatrix} 16 \\ 6 \\ 1 \end{bmatrix} \xrightarrow{\quad} \begin{bmatrix} 12 \\ 6 \\ 1 \end{bmatrix} \xrightarrow{\quad} \begin{bmatrix} 12 \\ 2 \\ 1 \end{bmatrix} \xrightarrow{\quad} \begin{bmatrix} 6 \\ 2 \\ 1 \end{bmatrix} \xrightarrow{\quad} \begin{bmatrix} 3 \\ 2 \\ 1 \end{bmatrix}$$

With $b = \begin{bmatrix} 3 \\ 2 \\ 1 \end{bmatrix}^T$; $A^T b$ is obtained as follows:

$$\begin{bmatrix} 3 \\ 2 \\ 1 \end{bmatrix} \xrightarrow{\quad} \begin{bmatrix} 6 \\ 2 \\ 1 \end{bmatrix} \xrightarrow{\quad} \begin{bmatrix} 12 \\ 2 \\ 1 \end{bmatrix} \xrightarrow{\quad} \begin{bmatrix} 12 \\ 4 \\ 1 \end{bmatrix} \xrightarrow{\quad} \begin{bmatrix} 16 \\ 6 \\ 1 \end{bmatrix} \xrightarrow{\quad} \begin{bmatrix} 16 \\ 6 \\ 3 \end{bmatrix} \xrightarrow{\quad} \begin{bmatrix} 16 \\ 6 \\ 15 \end{bmatrix} \xrightarrow{\quad} \begin{bmatrix} 16 \\ 38 \\ 63 \end{bmatrix}$$

2.1.2 Hybrid Operations

Method of triangular factorization can be extended for the hybrid operations as follows.

Let the hybrid column g is defined as follows -

$$g^T = (b_1, b_2, \dots, b_k, x_{k+1}, \dots, x_n) \quad (2.38)$$

An intermediate column vector b representing forward substitution for the first k rows is as follows:

$$b^T = (b_1^{(1)}, b_2^{(2)}, \dots, b_k^{(k)}, x_{k+1}, \dots, x_n)$$

$$= D_k L_k D_{k-1} L_{k-1} \dots, D_2 L_2 L_1 g \quad (2.39)$$

Backward substitution for the first k rows will give

$$x_k, x_{k-1}, \dots, x_2, x_1$$

$$x = U_1 U_2 \dots U_{k-1} U_k b$$

$$x = U_1 U_2 \dots U_{k-1} D_k L_k \dots D_3 L_3 D_2 L_2 D_1 b \quad (2.40)$$

$$b = D_1^{-1} L_1^{-1} D_2^{-1} L_2^{-1} \dots D_n^{-1} L_n^{-1} U_{n-1}^{-1} \dots U_{k+1}^{-1} D_k L_k \dots \\ D_3 L_3 D_2 L_2 D_1 b \quad (2.41)$$

In order to obtain the complete vector; complete operation represented by (2.41) is not necessary since first k elements are known. Last $n-k$ elements may be obtained from a matrix b' defined as follows:

$$b' = L_{k+1}^{-1} D_{k+1}^{-1} \dots L_n^{-1} D_n^{-1} U_{n-1}^{-1} \dots U_{k+1}^{-1} D_k L_k \dots \\ D_3 L_3 D_2 L_2 D_1 b \quad (2.42)$$

2.1.3 Numerical Example

Given $g^t = (b_1, b_2, x_3) = (20, 38, 1)$ to obtain
 $x = (x_1, x_2, x_3)$ for the matrix of example 2.1.1

$$\begin{array}{c|c|c|c|c|c} 20 & 10 & 10 & 10 & 10 & 3 \\ \downarrow & \downarrow & \downarrow & \downarrow & \downarrow & \\ 38 & 38 & 8 & 4 & 2 & 2 \\ \downarrow & \downarrow & \downarrow & \downarrow & \downarrow & \\ 1 & 1 & 1 & 1 & 1 & 1 \end{array}$$

To obtain (b_1, b_2, b_3) alone

$$\left[\begin{array}{c|c|c|c|c|c} 20 & 10 & 10 & 10 & 10 & 10 \\ 38 & 38 & 8 & 4 & 4 & 4 \\ 1 & 1 & 1 & 1 & 3 & 51 \end{array} \right]$$

$$b = (20, 38, 51)^t$$

Computer Logic

2.1.4 To obtain the L.U. factors of $n \times n$ matrix A

In this programme, elements of $n \times n$ matrix A are replaced by the L.U. factors of matrix A.

(i) Perform up to (vi) for $i = 1, \dots, n$.

(ii) Perform (iii) for $k = 1, \dots, i-1$.

(iii) $a_{ij} = a_{ij} - a_{ik} a_{kj}; \quad j = k+1, \dots, n$

(iv) $d = 1/a_{11}$

(v) $a_{ij} = d a_{ij}; \quad j = i+1, \dots, n$.

(vi) $a_{11} = d$

2.1.5 To simulate $A^{-1} b$

(i) Set $x_i = b_i; \quad i = 1, \dots, n$.

Forward substitution

(ii) Perform up to (iv) for $i = 1, \dots, n$.

$$(iii) x_i = x_i - \sum_{k=1}^{i-1} a_{ik} x_k$$

$$(iv) x_i = b_{ii} x_i$$

Backward substitution

(v) Perform up to (vii) for $i = 1, \dots, n-1$.

$$(vi) j = n - i$$

$$(viii) x_j = x_j - \sum_{k=j+1}^n a_{jk} x_k$$

2.1.6 To simulate Ax

(i) Set $b_i = x_i$ for $i = 1, \dots, n$.

$$(ii) b_i = b_i + \sum_{j=i+1}^n a_{ij} b_j ; i = 1, \dots, n-1$$

(iii) Perform up to (vi) for $i = 0, \dots, n-1$.

$$(iv) j = n - i.$$

$$(v) b_j = b_j / a_{jj}$$

$$(vi) b_j = b_j + \sum_{k=1}^{j-1} a_{jk} b_k$$

2.1.7 To simulate $(A^{-1})^T b$

- (i) Set $x_1 = b_1 ; i = 1, \dots, n.$
- (ii) Perform (iii) for $i = 1, \dots, n-1.$
- (iii) $x_j = x_j - x_i a_{ij} ; j = 1+1, \dots, n.$
- (iv) Perform up to (vii) for $i = 0, \dots, n-1.$
- (v) $j = n - 1.$
- (vi) $x_j = x_j / a_{jj}$
- (vii) $x_k = x_k - a_{jk} x_j ; k = 1, \dots, j-1.$

2.1.8 To simulate $A^T x$

- (i) Set $b_i = x_i ; i = 1, \dots, n.$
- (ii) Perform up to (iv) for $i = 1, \dots, n.$
- (iii) $x_k = x_k + a_{ik} x_i ; k = 1, \dots, i-1.$
- (iv) $x_i = x_i / a_{ii}$
- (v) Perform up to (vii) for $i = 1, \dots, n-1.$
- (vi) $k = n - 1.$
- (vii) $x_j = x_j + x_k a_{kj} ; j = k+1, \dots, n.$

2.1.9 Simulation of hybrid matrix operations

(a) Given vector $\mathbf{g}^T = (b_1, \dots, b_k, x_{k+1}, \dots, x_n)$ to obtain (x) .

(i) Set $x_i = b_i$; $i = 1, \dots, k$.

(ii) Perform up to (iv) for $i = 1, \dots, k$.

$$(iii) x_i = x_i - \sum_{j=1}^{i-1} a_{ij} x_j$$

$$(iv) x_i = a_{ii} x_i$$

(v) Perform (vi) for $i = 0, \dots, k-1$.

$$(vi) x_{(k-i)} = x_{(k-i)} - \sum_{j=k-i+1}^n a_{k-i,j} x_j$$

(b) To obtain vector (b) alone.

(i) Set $x_i = b_i$; $i = 1, \dots, k$.

(ii) Perform up to (iv) for $i = 1, \dots, k$.

$$(iii) x_i = x_i - \sum_{j=1}^{i-1} a_{ij} x_j$$

$$(iv) x_i = a_{ii} x_i$$

(v) Perform (vi) for $i = k+1, \dots, n-1$.

$$(vi) x_i = x_i + \sum_{j=i+1}^n a_{ij} x_j$$

(vii) Perform up to (x) for $i = 0, \dots, n-1-k$

$$(viii) j = n - i.$$

$$(ix) \quad x_j = x_j/a_{jj}$$

$$(x) \quad x_j = x_j + \sum_{k=1}^{j-1} a_{jk} x_k$$

$$(xi) \quad b_i = x_i \quad \text{for } i > k.$$

2.1.10 Symmetrical matrices

If the matrix A is symmetrical; its inverse is also symmetrical for which the elements on one side of the diagonal do not have to be stored. In this section, it is shown how symmetry could be exploited in the table of factors.

$$u_{ji} = a_{ji}^{(j)} = d_{jj} a_{ji}^{(j-1)} ; \quad i > j \quad (2.43)$$

$$l_{ij} = a_{ij}^{(j-1)} \quad (2.44)$$

For a symmetrical matrix;

$$\begin{aligned} a_{ji} &= a_{ij} \\ a_{ji}^{(1)} &= a_{ji} - a_{ji} a_{ii}^{(1)} ; \\ &= a_{ji} - a_{ji} d_{ii} a_{ii} ; \quad \text{for } i > j \geq 2 \end{aligned} \quad (2.45)$$

$$\begin{aligned} a_{ij}^{(1)} &= a_{ij} - a_{ii} a_{ij} \\ &= a_{ij} - a_{ii} d_{ii} a_{ij} \end{aligned} \quad (2.46)$$

Hence

$$a_{ji}^{(1)} = a_{ij}^{(1)} ; \quad i > j \geq 2 \quad (2.47)$$

$$\begin{aligned} a_{ji}^{(2)} &= a_{ji}^{(1)} - a_{j2}^{(1)} a_{21}^{(2)} \\ &= a_{ji}^{(1)} - a_{j2}^{(1)} d_{22} a_{21}^{(1)} ; \text{ for } i > j \geq 3 \end{aligned} \quad (2.48)$$

$$\begin{aligned} a_{ij}^{(2)} &= a_{ij}^{(1)} - a_{i2}^{(1)} a_{2j}^{(2)} \\ &= a_{ij}^{(1)} - a_{i2}^{(1)} d_{22} a_{2j}^{(1)} \end{aligned}$$

Hence

$$a_{ij}^{(2)} = a_{ji}^{(2)} \quad i > j \geq 3$$

$$\begin{aligned} a_{ij}^{(k)} &= a_{ij}^{(k-1)} - a_{ik}^{(k-1)} a_{kj}^{(k)} \\ &= a_{ij}^{(k-1)} - a_{ik}^{(k-1)} d_{kk} a_{kj}^{(k-1)} \end{aligned} \quad (2.49)$$

for $i > j > k$

$$\begin{aligned} a_{ji}^{(k)} &= a_{ji}^{(k-1)} - a_{jk}^{(k-1)} a_{ki}^{(k)} \\ &= a_{ji}^{(k-1)} - a_{jk}^{(k-1)} d_{kk} a_{ki}^{(k-1)} \end{aligned}$$

Hence, if

$$\begin{aligned} a_{ij}^{(k-1)} &= a_{ji}^{(k-1)} ; \quad i > j > k \\ a_{ij}^{(k)} &= a_{ji}^{(k)} \end{aligned} \quad (2.50)$$

Since (2.50) is valid for $k = 2$; it is valid for all $k < j$.

$$a_{ij}^{(j-1)} = a_{ji}^{(j-1)} \quad ; \quad i > j$$

$$u_{ji} = d_{jj} a_{ji}^{(j-1)} - d_{jj} l_{ij} \quad (2.51)$$

Equation (2.51) indicates that it is not necessary to store both l and u terms separately. l terms can be obtained from d and u terms, thus advantages of symmetry are reflected in the table of factors. If u terms are stored row by row; l terms can be obtained column by column; by dividing the corresponding row of u by 'd' term. The formulation of A^{-1} of equation (2.34) need complete row of l terms. It appears desirable to modify the formulation so that complete column of l terms is needed instead.

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Matrix L_k defined in equation (2.27) differs from the unit matrix only in k th row. This is broken in $(k-1)$ 'elementary matrices' E_{kj}^1 where $E_{kj}; j < k$ has its k jth element as an essential element with the value ' $-l_{kj}$ '.

$$L_k = E_{k1}^1 E_{k2}^1 E_{k3}^1 \dots E_{k,k-1}^1 \quad (2.52)$$

$$\begin{aligned} \text{Since } A^{-1} &= U_1 U_2 \dots U_{n-1} D_n L_n \dots L_2 L_2 D_1 \\ &= U_1 U_2 \dots U_{n-1} D_n E_{n1}^1 E_{n2}^1 \dots E_{n,n-1}^1 \\ &\quad \dots E_{n-1,1}^1 E_{n-1,2}^1 \dots E_{n-1,n-2}^1 E_{n-1,n-2}^1 \\ &\quad \dots \dots \dots \dots \dots \dots \dots \dots \dots \\ &\quad \dots \dots D_3 E_{31}^1 E_{32}^1 D_2 E_{21}^1 D_1 \end{aligned} \quad (2.53)$$

Two elementary matrices E_{ij} and E_{kl} will commute if $i \neq l$ and also $j \neq k$; and E_{ij}, E_k will commute if $i \neq k$ or $j \neq k$. Equation (2.54) is regrouped as follows:

$$\begin{aligned} A^{-1} &= U_1 U_2 \dots U_{n-1} D_n E_{n,n-1}^1 E_{n,n-2}^1 E_{n-1,n-2}^1 \\ &\quad D_{n-2} E_{n,n-3}^1 E_{n-1,n-3}^1 E_{n-2,n-3}^1 \dots \\ &\quad D_k E_{n,k-1}^1 E_{n-1,k-1}^1 \dots E_{k,k-1}^1 \dots \\ &\quad D_2 E_{n1}^1 E_{n-1,1}^1 \dots E_{21}^1 D_1 \end{aligned}$$

Let $L'_k = E_{nk}^1 E_{n-1,k}^1 \dots E_{k+1,k}^1$

Matrix L'_k differ from the unit matrix in k th column as follows:

$$L'_k : \text{ col } k = [0, \dots, 0, 1, -1_{k+1,k}, -1_{k+2,k}, \dots, -1_{nk}]$$

$$\begin{aligned} A^{-1} &= U_1 U_2 \dots U_{n-1} D_n L'_{n-1} D'_{n-1} L'_{n-2} \dots \\ &\quad D_2 L'_1 D_1 \end{aligned} \tag{2.54}$$

This equation, true for symmetrical and non-symmetrical matrices amounts to elimination by columns.

From relation (2.51);

$$L'_1 = D_1 U_1^t D_1^{-1} \text{ for symmetrical matrices;}$$

$$\mathbf{A}^{-1} = U_1 U_2 \dots U_{n-1} D_n D_{n-1} U_{n-1}^T D_{n-1}^{-1} D_{n-1}$$

$$D_{n-2} U_{n-2}^T D_{n-2}^{-1} \dots \dots \dots \dots \dots \dots$$

$$\dots \dots \dots \dots \dots \dots \dots \dots \dots \dots$$

$$\dots \dots \dots D_2 U_2^T D_2^{-1} U_2 D_1 U_1^T D_1^{-1} D_1$$

$$= U_1 U_2 \dots U_{n-1} D_n D_{n-1} U_{n-1}^T D_{n-2} U_{n-2}^T \dots$$

$$\dots D_1 U_1^T$$

Since $U_i^T D_j$ commute for $i \neq j$;

$$\mathbf{A}^{-1} = U_1 U_2 \dots U_{n-1} D_n D_{n-1} \dots D_1 U_{n-1}^T \dots U_2^T U_1^T$$

for symmetrical matrix (2.55)

2.1.11 Numerical example

Let the symmetrical matrix A is the following:

$$\begin{bmatrix} 2 & -1 & -1 & 0 \\ -1 & 4 & -2 & -1 \\ -1 & -2 & 5 & -2 \\ 0 & -1 & -2 & 4 \end{bmatrix}$$

Arithmetic steps for factorisation are as follows:

$$a_{11} = 1/a_{11} = 1/2$$

$$B = a_{11} \cdot a_{12} = (1/2) \cdot (-1) = -1/2$$

$$a_{22} = a_{22} - B \cdot a_{12} = 4 - (-1/2) \cdot (-1) = 7/2 \quad ①$$

$$a_{12} = B = -1/2$$

$$a_{23} = a_{23} - B \cdot a_{13} = 5 - (-1/2) \cdot (-1) = -5/2$$

$$a_{24} = a_{24} - B \cdot a_{14} = -1 - (-1/2) \cdot (0) = -1$$

$$a_{22} = 1/a_{22} = 2/7$$

$$B = a_{11} \cdot a_{13} = (1/2) \cdot (-1) = -1/2$$

$$a_{33} = a_{33} - B(a_{13}) = 5 - (-1/2) \cdot (-1) = 9/2$$

$$a_{34} = a_{34} - B(a_{14}) = -2 - (-1/2)(0) = -2$$

$$a_{13} = B = -1/2$$

$$B = a_{22} \cdot a_{23} = (2/7) \cdot (-5/2) = -5/7$$

$$a_{33} = a_{33} - B \cdot a_{23} = 9/2 - (-5/7) \cdot (-5/2) = 19/7$$

$$a_{34} = a_{34} - B \cdot a_{24} = -2 - (-5/7) \cdot (-1) = -19/7$$

$$a_{23} = B = -5/7$$

$$a_{33} = 1/a_{33} = 7/19$$

$$B = a_{22} a_{24} = (2/7)(-1) = -2/7$$

$$a_{44} = a_{44} - B a_{24} = 4 - (-2/7)(-1) = 26/7$$

$$a_{24} = B = -2/7$$

$$B = a_{33} a_{34} = (7/19)(-19/7) = -1$$

$$a_{44} = a_{44} - B a_{34} = 26/7 - (-1)(-19/7) = 1$$

$$a_{34} = B = -1$$

$$a_{44} = 1/a_{44} = 1.$$

Omitting the terms on the left of the diagonal, table of factors for the matrix is as follows:

| | | | |
|--------|--------|--------|-----|
| $1/2$ | $-1/2$ | $-1/2$ | 0 |
| $2/7$ | $-5/7$ | $-2/7$ | |
| $7/19$ | -1 | | |
| | | 1 | |

For a given vector $b^t = (1, 4, -1, -4)$; x may be obtained as follows:

$$\left[\begin{array}{c|ccccc|c} 1 & 1 & 1 & 1 & 1/2 & 1/2 & 2 \\ 4 & 9/2 & 9/2 & 9/2 & 9/7 & 2 & 2 \\ -1 & -1/2 & 19/7 & 19/7 & 1 & 1 & 1 \\ -4 & -4 & -19/7 & 0 & 0 & 0 & 0 \end{array} \right]$$

2.1.12 Computer logic for symmetrical matrix

(a) Computation of table of factors for the symmetrical matrix:

(i) Perform up to (viii) for $i = 1, \dots, n$.

(ii) Perform up to (vii) for $j = 1, \dots, i-1$.

$$(iii) B = a_{ji} a_{jj}$$

$$(iv) a_{ii} = a_{ii} - B a_{ji}$$

$$(v) a_{ji} = B$$

(vi) Perform up to (vii) for $k = i+1, \dots, n$.

$$(vii) a_{ik} = a_{ik} - B a_{jk}$$

$$(viii) a_{ii} = 1/a_{ii}$$

(b) Computation of $x = A^{-1}b$ from the table of factors.
(Equation 2.55):

(i) Set $x_1 = b_1$; $i = 1, \dots, n$.

(ii) Perform up to (iv) for $i = 1, \dots, n-1$.

(iii) Perform up to (iv) for $j = i+1, \dots, n$

$$(iv) x_j = x_j - x_i a_{ij}$$

$$(v) x_i = a_{ii} x_i ; i = 1, \dots, n$$

(vi) Perform up to (viii) for $i = 1, \dots, n-1$

(vii) $j = n - i$

$$(viii) x_j = x_j - \sum_{k=j+1}^n a_{jk} x_k$$

2.1.13 Comparison between the method of triangular factorisation and inverse for a full matrix

The two methods are comparable in the following respects:

- (i) Once the inverse matrix and table of factors are obtained, computation of $A^{-1}b$ and $(A^t)^{-1}b$ needs the same number of multiplication and addition (n^2).
- (ii) Elements on one side of the diagonal need not be stored for a symmetrical matrix in both the methods. Thus storage requirement is the same.

The method of triangular factorisation has the following advantages over the method of inverse:

- (i) The table of factors can be obtained in one-third the operations of the inverse.
- (ii) The method of triangular factorisation gives the effect of (A) ; A^t ; and certain hybrid operations; whereas the method of inverse does not.

The advantages of inverse are:

- (i) Complete solution needs $k \times n$ multiplications and additions if the vector 'b' has k non zero elements.
- (ii) Under some circumstances, the inverse may be more easily modified to reflect changes in the original matrix.

2.2 Compact storage scheme for sparse matrix

Although explicit matrix storage can be justified in some cases of relatively small sparse matrices, it is not practical for the matrices of most power network problems. A compact matrix storage scheme in which only the non-zero elements are retained is essential. A properly programmed compact storage scheme also results in considerable saving of computer time during matrix operations.

One possible scheme for a general matrix stores the non-zero elements of successive rows in a linear array. The column location of these non-zero elements and the location where the next row starts (row index) is stored separately. Following example explains the scheme.

2.2.1 Example

If the matrix A expressed explicitly is as follows:

| | | | | |
|-------|-------|-------|-------|-------|
| d_1 | | | u_1 | |
| | c_2 | | | u_2 |
| l_1 | | d_3 | | |
| | | | d_4 | u_3 |
| | | l_2 | | d_5 |

One possible compact storage scheme is as follows:

Table 2.2.1

| Col. No. | 1 | 2 | 3 | 4 | 5 | 6 | | | | |
|-----------------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|
| Row Index | 1 | 3 | 5 | 7 | 9 | 11 | | | | |
| Col. No. | 1 | 2 | 3 | 4 | 5 | 6 | 7 | 8 | 9 | 10 |
| Nonzero element | d_1 | u_1 | d_2 | u_2 | l_1 | d_3 | d_4 | u_3 | l_2 | d_5 |
| Column location | 1 | 4 | 2 | 5 | 1 | 3 | 4 | 5 | 3 | 5 |

Many variation of this scheme suitable for different situations, depending upon the nature of matrix or the nature of matrix operation expected are possible. An efficient storage scheme is again a trade off between the storage requirements

and the computer time during the matrix operations. In this project the sparse matrix encountered are (a) Admittance matrix (b) L.U. factors of the jacobian matrix. Methods of storing these matrices are discussed below.

2.2.2 Admittance matrix

This matrix is complex and symmetrical and can be stored in polar or rectangular form. Elements on the left side of the diagonal need not be stored. A feasible scheme is as follows.

If the symmetrical matrix Y expressed explicitly is as follows:

| | | | | |
|-------|-------|-------|-------|-------|
| d_1 | a_1 | | a_2 | |
| a_1 | d_2 | a_3 | | |
| | a_3 | d_3 | | a_4 |
| a_2 | | | d_4 | |
| | | a_4 | | d_5 |

Compact storage scheme is as follows:

Table 2.2.2 a

| No. | 1 | 2 | 3 | 4 | 5 |
|---|-------|-------|-------|-------|-------|
| Diagonal element | d_1 | d_2 | d_3 | d_4 | d_5 |
| Row index for right off sub-diagonal elements | 1 | 3 | 4 | 5 | 5 |
| No. | 1 | 2 | 3 | 4 | 5 |
| Nonzero right sub-diagonal element | a_1 | a_2 | a_3 | a_4 | - |
| Column location | 1 | 4 | 3 | 5 | - |

It is sometimes desired to obtain all the non-zero elements of one row of the admittance matrix (y_{ij} ; $j=1, \dots, n$). For $j > i$, non-zero elements can be obtained by obtaining the row index for i and $i+1$. For $j < i$, a search is needed from all the row indices from 1 to i . This may be quite time consuming and due to the fact that many of these elements are likely to be zero; this approach has not been considered very efficient.

In the computer programme; additional array of row index, column location and location of the element in the array of non-zero right sub-diagonal elements are formulated.

Additional table may be as follows.

Table 2.2.2 b

| S.No. | 1 | 2 | 3 | 4 | 5 | 6 |
|--|---|---|---|---|---|---|
| Row index for left sub-diagonal elements | 1 | 1 | 2 | 3 | 4 | 5 |
| S.No. | 1 | 2 | 3 | 4 | 5 | |
| Column location for left sub-diagonal element | 1 | 2 | 1 | 3 | - | |
| Location in the array of non-zero right sub-diagonal element | 1 | 3 | 2 | 4 | - | |

The last array of this table represents the location of the non-zero element in the main array (Table 2.2.2 a). If instead the non-zero element itself is represented, additional storage is needed since these elements are complex and are in real mode.

2.2.3 Table of L.U. factors

Equations (2.34) to (2.37) indicate that a complete row of elements on the left or right of diagonal is needed simultaneously. Hence the scheme similar to that of storing the

symmetrical matrix has been adopted. The last array of table (2.2.2 b) however has the actual non-zero element on the left of diagonal. This is explained through Table(2.2.3) for example of (2.2.1).

Table 2.2.3:

Compact storage scheme of for example 2.2.1:

| S.No. | 1 | 2 | 3 | 4 | 5 | 6 |
|---|-------|-------|-------|-------|-------|-------|
| Diagonal element | d_1 | d_2 | d_3 | d_L | d_5 | d_6 |
| Row index for right sub-diagonal elements | 1 | 2 | 3 | 3 | 4 | 4 |
| Row index for left sub-diagonal elements | 1 | 1 | 1 | 2 | 2 | 3 |
| S.No. | 1 | 2 | 3 | 4 | | |
| Non-zero right sub-diagonal element | u_1 | u_2 | u_3 | - | | |
| Non-zero left sub-diagonal element | l_1 | l_2 | - | - | | |
| Column of non-zero right sub-diagonal element | 4 | 5 | 5 | - | | |
| Column of non-zero left sub-diagonal element | 1 | 3 | - | - | | |

In case the table of factors has a symmetrical pattern of non-zero elements; as is the case with the conventional load flow solution through Newton's method, a more efficient storage scheme in terms of storage stores the complete row on the right of diagonal and complete column below the diagonal, through common index and integer array listing. Equation(2.54) may be used for A^{-1} . This needs the complete column below the diagonal at one time, instead of complete row on the left of the diagonal. This is explained through the following example.

2.2.4 Example

Let the matrix (A) with symmetrical pattern of non-zero elements written explicitly be the following:

| | | | | |
|-------|-------|-------|-------|-------|
| d_1 | | u_1 | | u_2 |
| | d_2 | | u_3 | |
| l_1 | | d_3 | | u_4 |
| | l_3 | | d_4 | |
| l_2 | | l_4 | | d_5 |

The compact storage scheme may be as follows.

Table 2.2.4

| | | | | | | |
|---|-------|-------|-------|-------|-------|---|
| S.No. | 1 | 2 | 3 | 4 | 5 | 6 |
| Diagonal element d_1 | d_1 | d_2 | d_3 | d_4 | d_5 | - |
| Index | 1 | 3 | 4 | 5 | 5 | 5 |
| S.No. | 1 | 2 | 3 | 4 | 5 | |
| Integer listing | 3 | 5 | 4 | 5 | - | |
| Non-zero right sub-diagonal element | u_1 | u_2 | u_3 | u_4 | - | |
| Non-zero sub- diagonal element below diagonal | l_1 | l_2 | l_3 | l_4 | - | |

Most versions of Fortran IV compilers store a rectangular matrix 'A' of dimension $M \times N$ by column in block of memory named A. The first column is followed by the next column and so on until the N columns are exhausted. Thus the double subscripted variable A is stored as a single subscripted variable. The element A(I, J) is interpreted as A($M \cdot J + I$). This rule is built into the Fortran compiler and the translation is automatic whenever the ordered pair is encountered in the coding. In the compact storage scheme non-zero elements

are not stored. If a sub-programme could be written such that the programmes written in conventional way could with little modification use the compact storage; this should be of great value.

2.3 Optimal renumbering of nodes

There are two primary objectives¹ for an ordered Gaussian elimination. The first and the oldest seeks to control numerical accuracy through a pivoting scheme. The second aims at conservation of matrix sparsity, since the order in which the Gaussian elimination is performed on sparse matrices affects the total number of new non-zero elements generated in the course of elimination and hence the total computation time. The physical nature of large scale electrical networks and the numerical accuracy of modern computers tend to eliminate the need for pivoting and to enhance the need for the exploitation of sparsity. Survey of this section is restricted to sparsity conserving ordering scheme.

Need for a reasonable renumbering scheme can be explained by the help of Fig. 2.3.1a. With the numbering schemes the non-zeros of Y-matrix and table of L.U. factors are shown respectively in Figures 2.3.1b and 2.3.1c.

If the same system is renumbered as shown in Fig. 2.3.2a; the total number of non-zero terms of the Y-matrix shown in



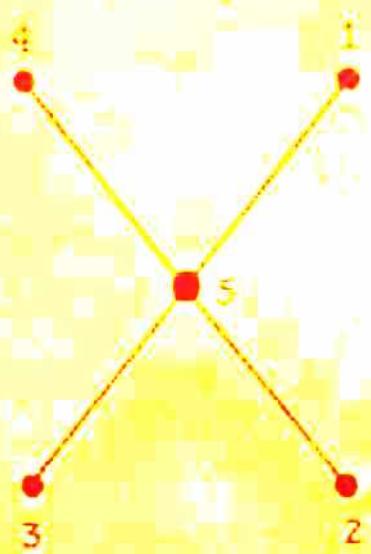
(a) LINEAR GRAPH

| | | | | |
|---|---|---|---|---|
| × | × | × | × | × |
| × | × | | | |
| × | | | × | |
| × | | | | × |
| × | | | | × |

(b) NON ZERO ELEMENTS OF Y-MATRIX.

| | | | | |
|---|---|---|---|---|
| × | × | × | × | × |
| × | × | × | × | × |
| × | × | × | × | × |
| × | × | × | × | × |
| × | × | × | × | × |

(c) NON ZERO ELEMENTS OF TABLE OF FACTORS.



LINEAR GRAPH.

| | | | | | |
|---|---|---|---|---|---|
| x | | | | | x |
| | | x | | | x |
| | | | x | | x |
| | | | | x | x |
| x | x | x | x | x | x |

(b) NON ZERO ELEMENTS OF Y-MATRIX.

| | | | | | |
|---|---|---|---|---|---|
| x | | | | | x |
| | | x | | | x |
| | | | x | | x |
| | | | | x | x |
| x | x | x | x | x | x |

(c) NON ZERO ELEMENTS OF TABLE OF FACTORS.

Fig. 2.3.2b, remain unchanged but the number of non-zero elements of the table of factors shown in Fig. 2.3.2c is very much less compared to that shown for the same system in Fig. 2.3.1c.

For a n node system; there are of course $n!$ permutations for numbering available. The absolute optimal order for elimination would result in the least possible non-zero elements in the table of factors. The only known method for determining the absolute optimal order is to examine all $n!$ possibilities. This seems impractical particularly if n is large. However, several satisfactory schemes have been developed for obtaining near optimal orders.

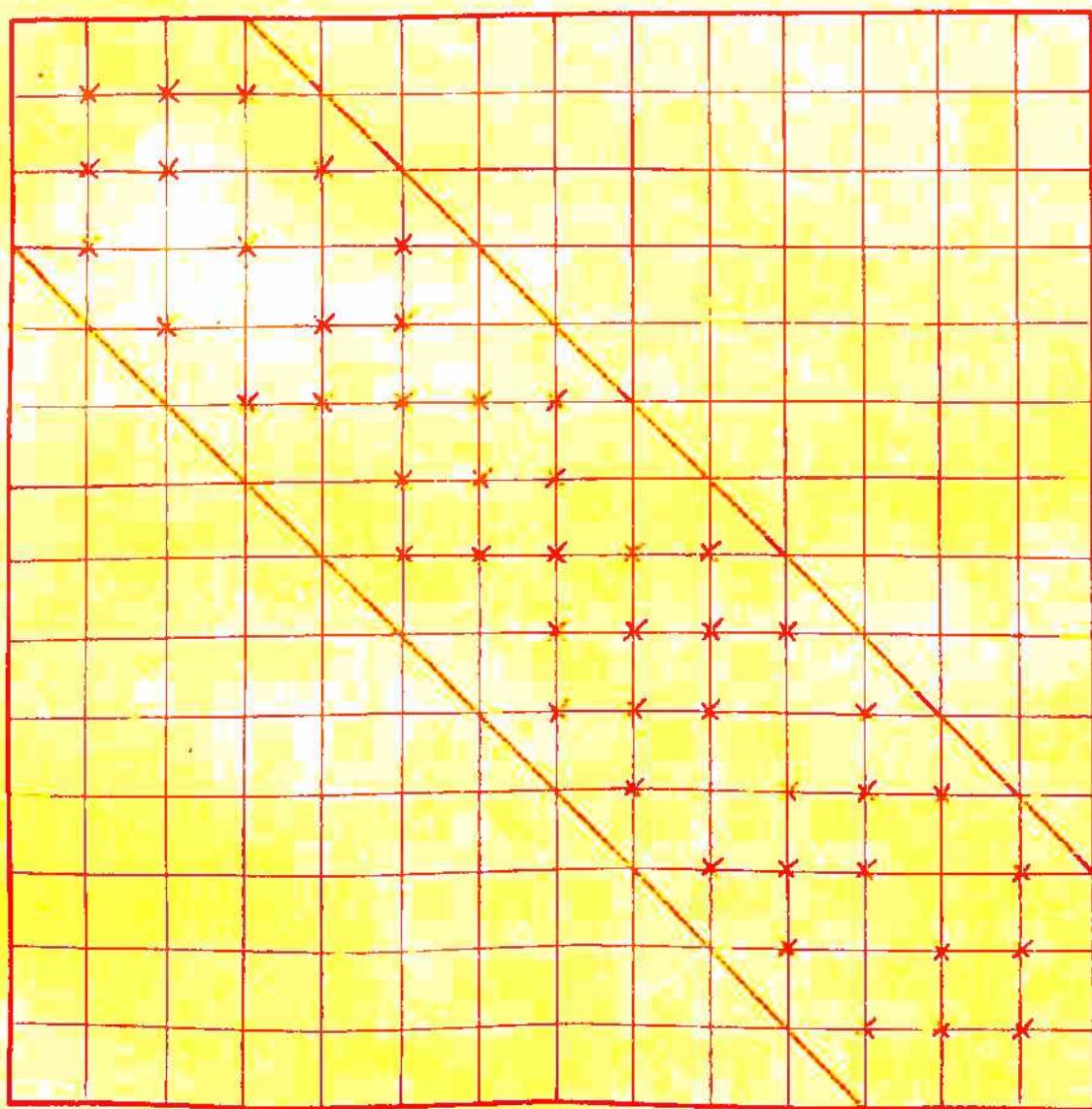
2.3.1 Matrix banding schemes

The objective of the banding schemes is to restrict the elements of a matrix to a narrow band about the major diagonal (Fig. 2.3.3.a) or about the minor diagonal (Fig. 2.3.3b).

2.3.2 Minimally adjacent cut sets (MACS)

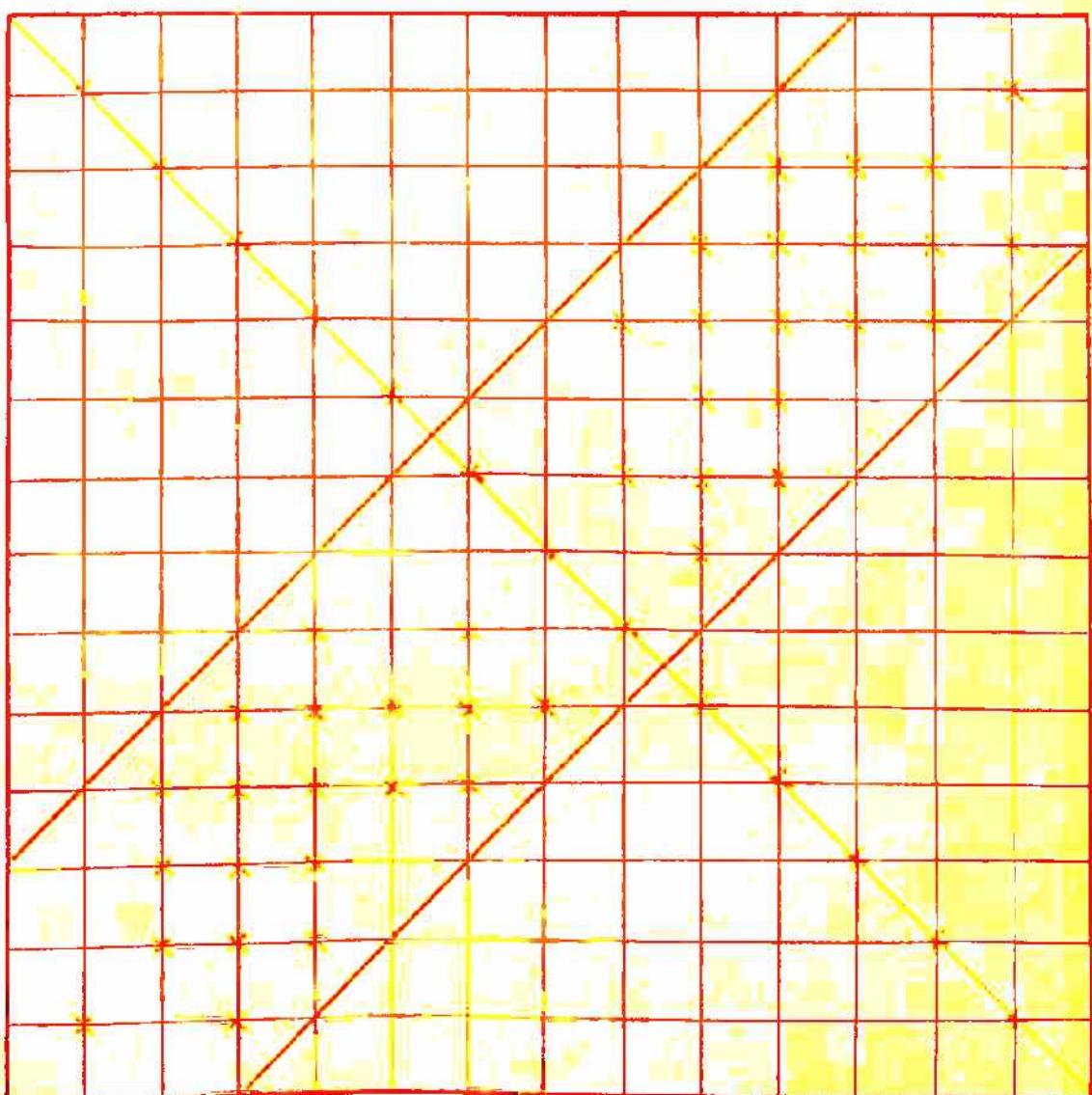
Given a set S of nodes of a connected graph G and a subset S_1 of nodes; the minimally adjacent cut sets (MACS) of S_1 are nodes from S not in S_1 but which are neighbours of nodes in S_1 .

2.3.3 An algorithm for major diagonal banding is as follows:



MAJOR DIAGONAL BANDING.

FIG. - 2.3.3 (a).



MINOR DIAGONAL BANDING .

FIG - 2.3.3 (b).

'Starting with an arbitrary node as number 1, the MACS of node 1 are assigned the next sequence of numbers; the MACS of the numbered nodes are subsequently assigned the next sequence of numbers and the process is continued until all the nodes of the network are numbered'.

Linear graph of AEP 14 bus test system, numbered as above and the non-zero elements of 'Y' matrix and table of factors are shown respectively in Fig. 2.3.4 and Fig. 2.3.5.

The algorithm is weak because the choice of initial guess is critical.

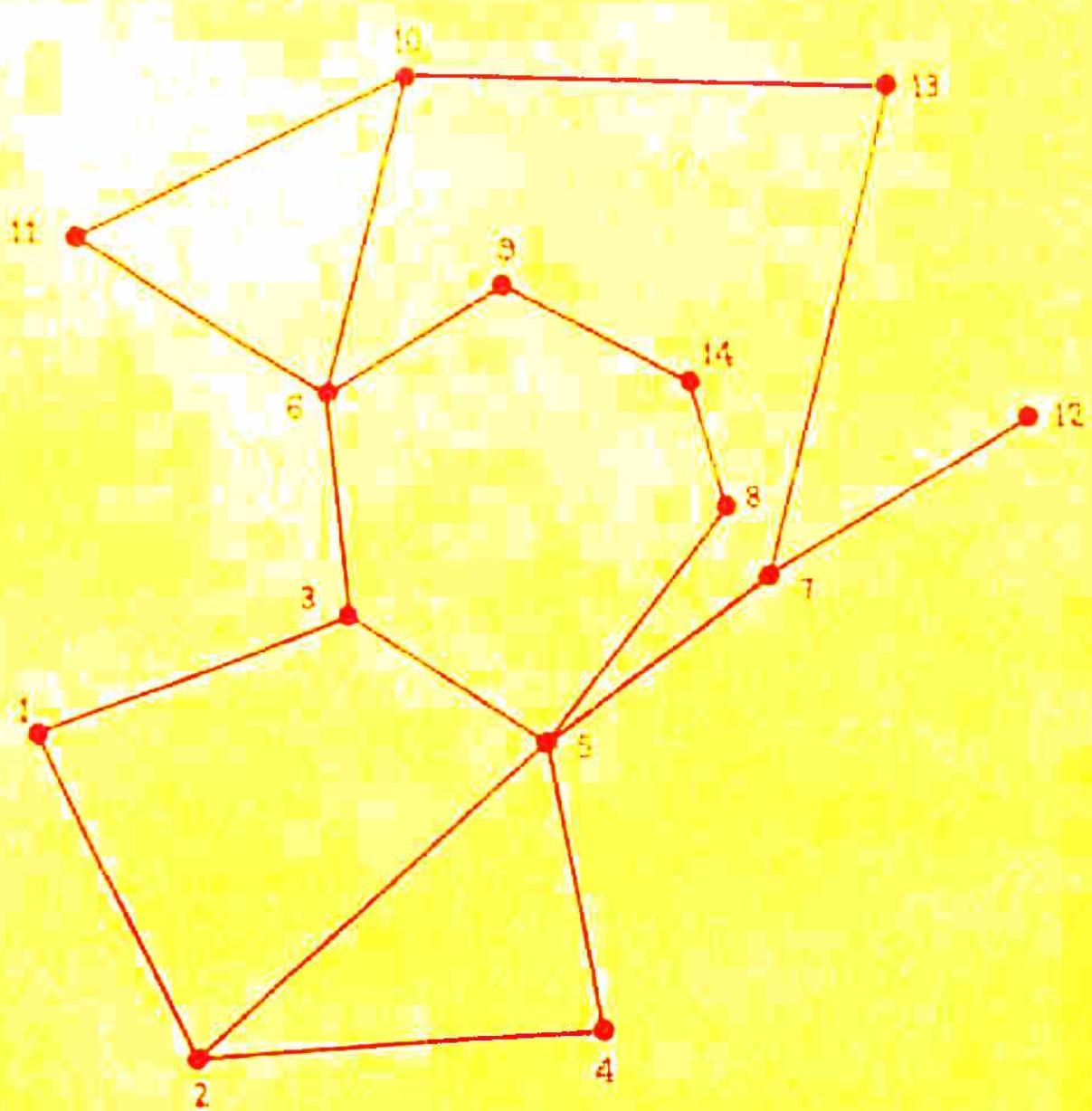
Carre²³ has suggested the following technique:

'First one node is arbitrarily chosen as last point and numbered n. Its MACS are assigned the next high valued numbers; the MACS of these numbered nodes are assigned, the next high valued numbers'.

The author does not understand how this method of numbering is effectively the same as the 'Scheme II' described by Tinney⁴⁵ (2.3.5), as claimed in the paper²³. This is essentially the major diagonal banding scheme.

2.3.4 An algorithm for minor diagonal banding is as follows:

- (1) Start with the bus having only one line and list it as node 1. In the absence of a node with only one incident branch; the node with least number of incident branches may be numbered as 1.



LINEAR GRAPH OF AEP 14 BUS TEST SYSTEM; NUMBERED THROUGH
MAJOR DIAGONAL BANDING SCHEME.

FIG. - 2·3·4

| | 1 | 2 | 3 | 4 | 5 | 6 | 7 | 8 | 9 | 10 | 11 | 12 | 13 | 14 |
|----|---|---|---|---|---|---|---|---|---|----|----|----|----|----|
| 1 | x | x | x | | | | | | | | | | | |
| 2 | x | x | o | x | x | | | | | | | | | |
| 3 | x | o | x | o | x | x | | | | | | | | |
| 4 | | x | o | x | x | o | | | | | | | | |
| 5 | | x | x | x | x | o | x | x | | | | | | |
| 6 | | | x | o | o | x | o | o | x | x | x | | | |
| 7 | | | | x | o | x | o | o | o | o | o | x | x | |
| 8 | | | | | x | o | o | x | o | o | o | o | o | x |
| 9 | | | | | | x | o | o | x | o | o | o | o | x |
| 10 | | | | | | x | o | o | o | x | x | o | x | o |
| 11 | | | | | | | x | o | o | o | x | x | o | o |
| 12 | | | | | | | x | o | o | o | o | x | o | o |
| 13 | | | | | | | | x | o | o | x | o | x | o |
| 14 | | | | | | | | | x | x | o | o | o | x |

x NON ZERO ELEMENTS OF Y-MATRIX

o ADDITIONAL NON ZERO ELEMENTS OF TABLE OF FACTORS

NON ZERO ELEMENTS OF AEP 14 BUS TEST SYSTEM, NUMBERED THROUGH
MAJOR DIAGONAL BANDING SCHEME.

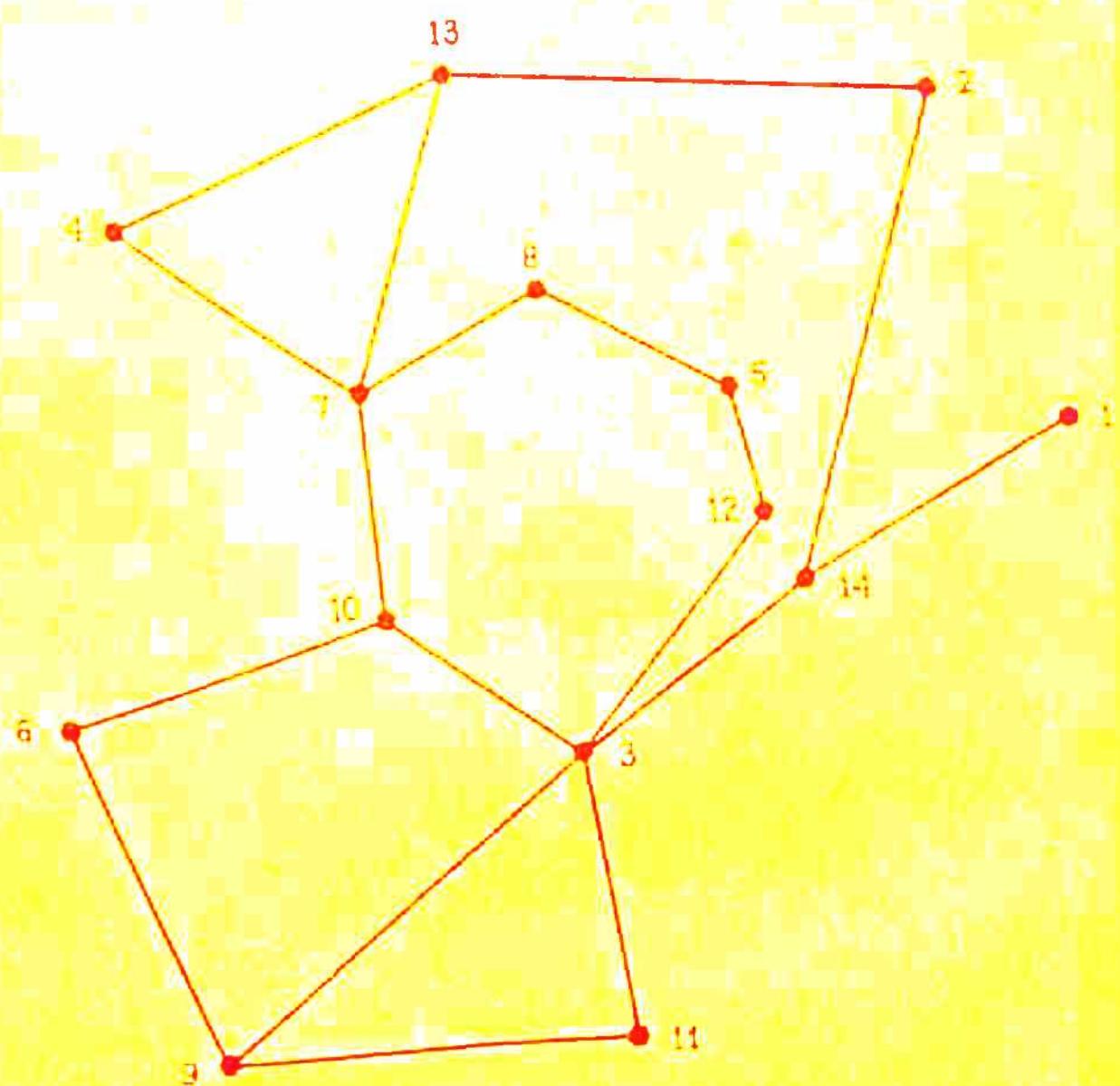
- (ii) List the MACS of node 1 as the last numbered nodes.
- (iii) Assign the next lower valued numbers to the MACS of the numbered nodes.
- (iv) Assign the next high valued numbers to the MACS of the numbered nodes and return to step (iii).

Linear graph of AEP 14 bus test system and the non-zero elements of the Y-matrix and table of factors are shown respectively in Fig. 2.3.6 and Fig. 2.3.7.

Obviously the diagonal banding schemes do not work for the system shown in Fig. 2.3.1a of which the absolute optimal order is in Fig. 2.3.2a in which the non-zero element of the table of factors remain the same as the non-zero elements of Y-matrix.

2.3.5 Tinney^{1,31,33,45} has given three schemes for near optimal renumbering. These schemes are claimed to be generally more effective than the banding schemes. These are listed in increasing order of complexity, execution time and optimality. It is assumed that the nodes are originally numbered according to some external criterion and then renumbered according to the schemes.

I. Number the rows according to the number of non-zero off diagonal elements before elimination. In this scheme the rows with only one off diagonal element are numbered first, those with two elements second, etc., and those



LINEAR GRAPH OF AEP 14 BUS TEST SYSTEM, NUMBERED THROUGH
MINOR DIAGONAL BANDING SCHEME.

| | 1 | 2 | 3 | 4 | 5 | 6 | 7 | 8 | 9 | 10 | 11 | 12 | 13 | 14 |
|----|---|---|---|---|---|---|---|---|---|----|----|----|----|----|
| 1 | x | | | | | | | | | | | | | x |
| 2 | | x | | | | | | | | | | x | x | |
| 3 | | | x | | | | | | x | x | x | x | x | x |
| 4 | | | | x | | | x | | | | | | x | |
| 5 | | | | | x | | | x | . | | | x | | |
| 6 | | | | | | x | | | x | x | | | | |
| 7 | | | x | | | | x | x | | x | | | x | |
| 8 | | | | x | | | x | x | | o | | o | | |
| 9 | | | x | | | x | | | x | o | x | o | o | |
| 10 | | | x | | | x | x | o | o | x | o | o | o | |
| 11 | | | x | | | | | x | o | x | o | | o | |
| 12 | | | x | | x | | | o | o | o | o | x | | o |
| 13 | | x | x | | | | x | | | . | | x | o | |
| 14 | x | x | x | | | | | | o | o | o | o | o | x |

x NON ZERO ELEMENTS OF Y-MATRIX

o ADDITIONAL NON ZERO ELEMENTS OF TABLE OF FACTORS

NON ZERO ELEMENTS OF AEP 14 BUS TEST SYSTEM , NUMBERED THROUGH
MINOR DIAGONAL BANDING SCHEME .

with the most elements last. This scheme does not take into account any of the subsequent effects of the elimination process. The only information needed is the number of non-zero element in each row of the original matrix.

II. Number the rows such that at each step of the process the next row, to be operated upon is the one with the fewest non-zero elements. If more than one row meet this criterion, select any one. This scheme needs a simulation of the effects on the accumulation of non-zero elements of elimination by columns. Required input information is a list by rows of the column numbers of the non-zero off diagonal elements.

III. Number the rows such that at each step of the process, the next row to be operated upon is the one that will introduce the fewest new non-zero elements.

This involves a trial simulation of every feasible alternative at each step. Input information is the same as for scheme (II).

It is hard to evaluate the three schemes since the performance of any scheme depends upon the type of system. However scheme I, though simplest, does not appear to be efficient; since subsequent built up of non-zero elements are not at all taken care of, if many solutions of the matrix are needed.

Scheme III needs large computer time since it involves trying many possible permutations.

2.3.6 At present author's subroutine works as follows:

'Number the rows such that the next row is the one having fewest non-zero elements in the table of factors of that row',

The author is now sure if this is the Tinney's II scheme.

A comparison of some of the techniques applied on AEP 14 bus test system is as follows:

Non zero elements in the Y-matrix = 50

Additional non-zero elements with major diagonal banding scheme

(Fig. 2.3.5) = 54

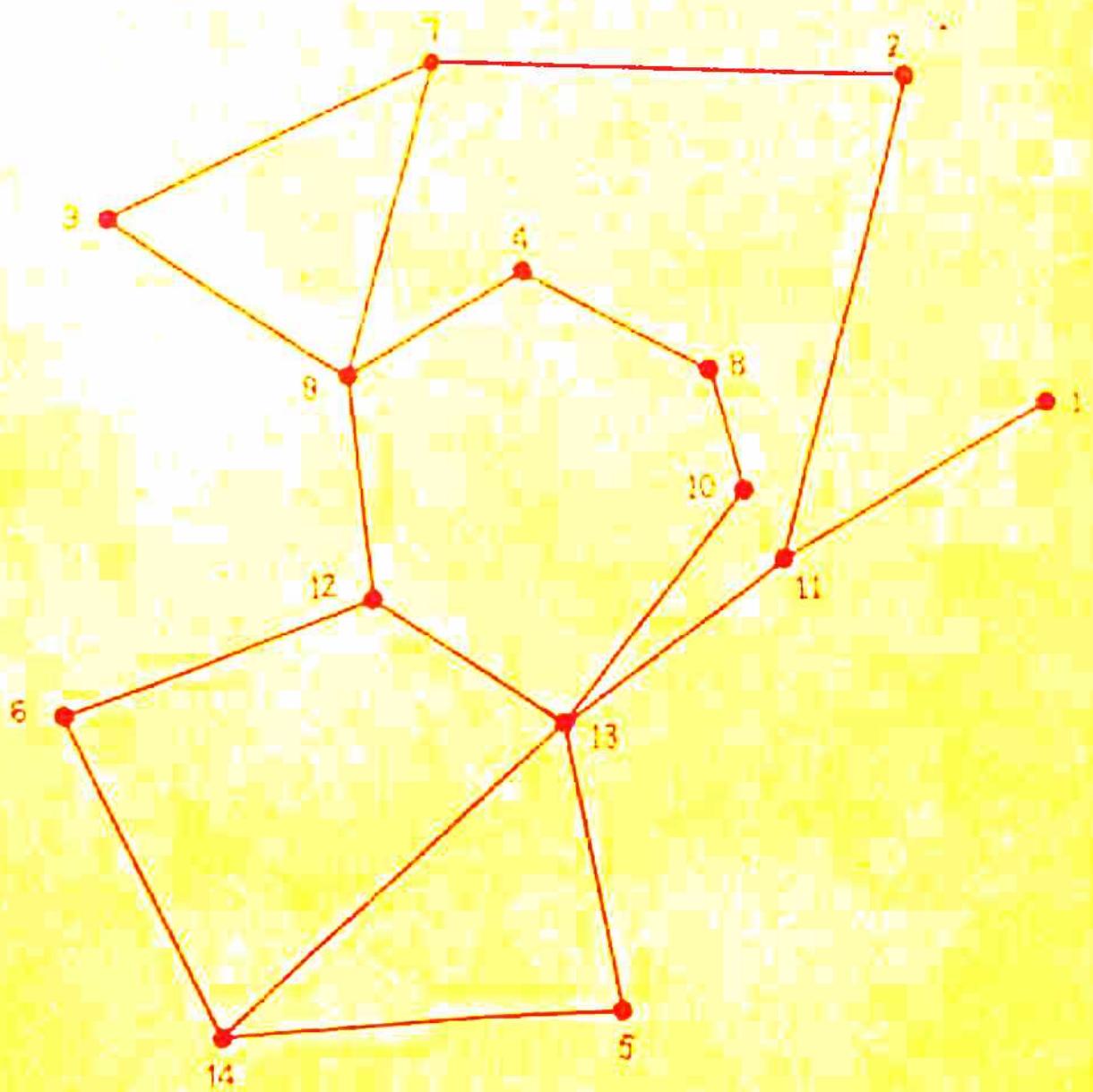
Additional non-zero elements with minor diagonal banding scheme

(Fig. 2.3.7) = 24

Additional non-zero elements with author's subroutine (Fig.2.3.9) = 16

Computer logic of author's subroutine for renumbering the buses is as follows:

For the set 'S' of n nodes; the nodes already renumbered from 1, ..., n_1 , be in the subset S_1 ; where n_1 is the total number of nodes already renumbered; remaining nodes are in subset S_2 such that $S = S_1 + S_2$.



LINEAR GRAPH OF AEP 14 BUS TEST SYSTEM, NUMBERED THROUGH
AUTHORS SUBROUTINE.

FIG.-2-3-8

| | 1 | 2 | 3 | 4 | 5 | 6 | 7 | 8 | 9 | 10 | 11 | 12 | 13 | 14 | |
|----|---|---|---|---|---|---|---|---|---|----|----|----|----|----|--|
| 1 | x | | | | | | | | | | | | | x | |
| 2 | | x | | | | | x | | | | | x | | | |
| 3 | | x | | | | | x | | x | | | | | | |
| 4 | | | x | | | | | x | x | x | | | | | |
| 5 | | | | x | | | | | | | | | x | x | |
| 6 | | | | | x | | | | | | | x | | x | |
| 7 | x | x | | | | | x | | x | | o | | | | |
| 8 | | | x | | | | | x | o | x | | x | | | |
| 9 | | x | x | | | | x | o | x | o | o | x | | | |
| 10 | | | | | x | | | x | o | x | o | o | x | | |
| 11 | x | x | | | | o | | o | o | x | o | x | | | |
| 12 | | | | x | | | | x | o | o | x | x | x | o | |
| 13 | | | | | x | | | | x | x | x | x | x | x | |
| 14 | | | | x | x | | | | | | o | x | x | | |

x NON ZERO ELEMENTS OF Y-MATRIX

o ADDITIONAL NON ZERO ELEMENTS OF TABLE OF FACTORS

NON ZERO ELEMENTS OF AEP 14 BUS TEST SYSTEM MATRIX,
NUMBERED THROUGH AUTHOR'S SUBROUTINE.

- (i) Let $n_1 = 0$; and all the nodes be in S_2 .
- (ii) For all the remaining nodes, having their old numbers as $i = 1, \dots, n$; $i \notin S_1$; perform up to (vi).
- (iii) Obtain the columns of non-zero elements of i th row from the row indices and column locations; and let them be in set C_1 . Other nodes are in set C_2 such that $S = C_1 + C_2$.
- (iv) Perform up to (v) for $j = 1, \dots, n_1$, where j is the new number of buses already renumbered. This way all the columns of non-zero elements of table of factor is obtained if the next node is i .
- (v) Obtain the old number for j and if it lies in set C_1 , all the columns of the elements of the table of factors on the right hand side of the diagonal for the j th row obtained in step (viii) are also brought in set C_1 .
- (vi) Obtain the total number of elements n_2 in the set $C_1 + S_2$. If they are less than the least such numbers obtained for other values of i ; let $m = n_2$ and $k = 1$.
- (vii) $n_1 = n_1 + 1$
- (viii) The next row to be operated upon is k and its new number is n_1 . Obtain for this row the old column numbers of the non-zero elements on the right side in the table of factors as obtained in step (iv) and (v) and store them

in the compact storage scheme for n_1 th row, as shown
in table 2.2.2 a.

(ix) If $n_1 < n$; repeat from step (ii). If $n_1 = n$, renumbering
job is complete.

CHAPTER III

FORMATION OF NETWORK ADMITTANCE MATRIX

In this chapter algorithm for forming the admittance matrix and storing it in the compact form is described. Effects of tap changing transformer and phase shifting transformer on the admittance matrix are described but the algorithm takes care of only tap changing transformer. With phase shifting transformer, the admittance matrix is no more symmetrical. Hence it is considered desirable to account for additional terms due to the phase shifting transformer separately.

3.1 Admittance matrix of a line connected between two nodes

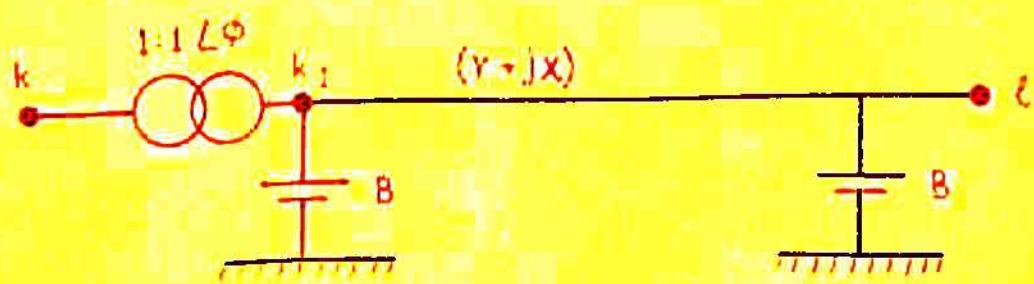
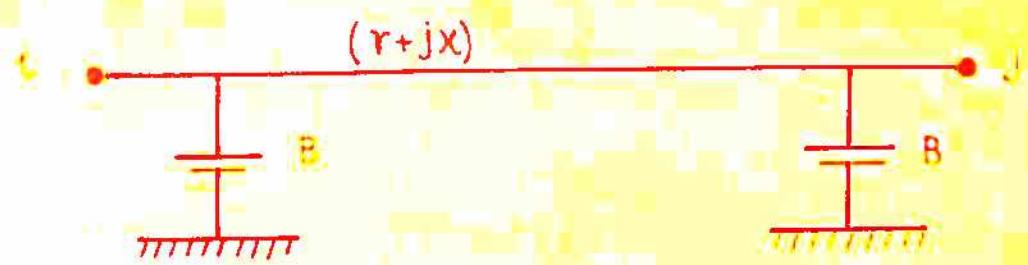
(Fig. 3.1)

$$\begin{bmatrix} I_1 \\ I_J \end{bmatrix} = \begin{bmatrix} y+jB & -y \\ -y & y+jB \end{bmatrix} \begin{bmatrix} V_1 \\ V_J \end{bmatrix} \quad (3.1)$$

where $y = 1/(r+jx)$ (3.2)

$$= \frac{r}{(r^2+x^2)} - \frac{jx}{(r^2+x^2)}$$

3.2 Admittance matrix of a line connected between two nodes through an adjustable transformer (Fig. 3.2)



$$\begin{bmatrix} I_{k1} \\ I_1 \end{bmatrix} = \begin{bmatrix} y+jB & -y \\ -y & y+jB \end{bmatrix} \begin{bmatrix} V_{k1} \\ V_1 \end{bmatrix} \quad (3.3)$$

If transformer losses are ignored;

$$V_{k1} = t V_k$$

$$I_k = t I_{k1}$$

$$\begin{bmatrix} I_k \\ I_1 \end{bmatrix} = \begin{bmatrix} t^2(y+jB) & -ty \\ -ty & (y+jB) \end{bmatrix} \begin{bmatrix} V_k \\ V_1 \end{bmatrix} \quad (3.4)$$

where y is defined by equation 3.2

k = near node of the transformer

l = far node of the transformer

t = off nominal turn ratio

For every adjustable tap setting transformer, its near node, far node, $(y+jB)$, $(-y)$ and location in the admittance array are stored separately.

If the transformer tap setting is modified; only one diagonal element and one stored off diagonal element is modified as follows.

$$y_{kk}^2 = y_{kk}^1 + (t_2 - t_1)(y + jB) \quad (3.5)$$

$$y_{kl}^2 = y_{kl}^1 - (t_2 - t_1)(y) \quad (3.6)$$

where y_{kk}^1 and y_{kl}^1 refer the tap setting t_1 , and y_{kk}^2 and y_{kl}^2 refer to the tap setting t_2 .

3.3 Admittance matrix of a line connected between two nodes through phase shifting transformer (Fig. 3.3)

$$\begin{bmatrix} I_{k1} \\ I_1 \end{bmatrix} = \begin{bmatrix} y+jB & -y \\ -y & y+jB \end{bmatrix} \begin{bmatrix} V_{k1} \\ V_1 \end{bmatrix} \quad (3.7)$$

$$V_{k1} = V_k e^{j\theta} \quad (3.8)$$

$$I_k = I_{k1} e^{-j\theta}$$

$$\begin{bmatrix} I_k \\ I_1 \end{bmatrix} = \begin{bmatrix} (y+jB) & -y e^{-j\theta} \\ -y e^{j\theta} & (y+jB) \end{bmatrix} \begin{bmatrix} V_k \\ V_1 \end{bmatrix} \quad (3.9)$$

Thus with the presence of phase shifting transformer, the admittance matrix is no more symmetrical.

3.4 Computer logic for forming the admittance matrix and storing it in compact form

The admittance matrix is formed by first setting the diagonal and off diagonal elements to zero. Data for one line which include the two nodes i and j to which the

line is connected, line resistance and reactance, shunt admittance and off nominal turn ratio and near node of the transformer (if present) are read and its admittance matrix formed and the elements added to the corresponding elements of compactly stored matrix. It is assumed that the number of first bus is smaller than the number of second bus and data are read in the ascending order of i. Data need not be arranged in the ascending order of j.

Notations

$$y_d^i = g_d^i + j b_d^i = \text{diagonal element of } i\text{th node}$$

$$I_1^i = \text{index sequence of } i\text{th row right off diagonal elements}$$

$$I_2^i = \text{index sequence of } i\text{th row, left off diagonal elements}$$

$$L_1^i = \text{column of } i\text{th right off diagonal element}$$

$$L_2^i = \text{column of } i\text{th left off diagonal element}$$

$$y^i = g^i + j b^i = i\text{th right off diagonal element}$$

$$L_3^i = \text{corresponding location of } i\text{th left off diagonal element in the array of right off diagonal elements}$$

$$n = \text{expected maximum number of lines}$$

$$n_t = \text{total number of adjustable tap setting transformers}$$

$$y_{td}^i = g_{td}^i + j b_{td}^i = (y + jB) \text{ of equation 3.3, for } i\text{th adjustable tap setting transformer}$$

k_1 = count of rows

k_2 = count of non-zero elements

t^i = tap setting of i th adjustable tap setting transformer

y_t^i = $\frac{g^i}{t} + j \frac{b^i}{t} = (-y)$ of equation 3.3 corresponding to i th adjustable tap setting transformer

I_n^i near and far nodes of i th adjustable tap setting
and n transformer respectively.

I_f^i

L_t^k = location of $i - j$ th right off diagonal element in array y if $i - j$ th element are connected through k th adjustable transformer.

The logic is as follows :

(i) Set y_t^1 and y_{td}^1 ; $l = 1, \dots, \bar{n}_t$, y_d^1 ; $l = 1, \dots, n$ and y^1 ; $l = 1, \dots, \bar{n}; n_t$ to zero.

(ii) Set $k_1 = 1$ and $k_2 = 0$

(iii) Set the real variables y_1^1 and y_2^1 ; and integer variables i_1^1 to zero for $l = 1, \dots, n$.

(iv) Read i , j , r , x , B , tap setting and near node of the transformer (if present). If i is equal to k_1 , proceed from (x), if i is greater than k_1 , proceed

from (v), and if i_1 less than k_1 , indicates data not arranged in ascending order of i and the programme is terminated.

$$(v) I_1^{k_1} = k_2 + 1$$

(vi) Perform (vii) for $l = k_1, \dots, n$; such that $I_1^l \neq 0$.

$$(vii) k_2 = k_2 + 1, L_1^{k_2} = 1, g^{k_2} = y_1^1, b^{k_2} = y_2^1, y_1^1 = 0, y_2^1 = 0, I_1^1 = 0.$$

(viii) $k_1 = k_1 + 1$ and if k_1 is less than 1 (of step iv) go to (ix); otherwise to (x).

(ix) $I_1^{k_1} = k_2 + 1$ and proceed from (viii).

(x) Check if j is greater than i . If not, data are not as required and the programme is terminated.

(xi) $I_j^j = 1, b_j = 1/(r^2 + x^2), b_1 = rb_3$, and $b_2 = -xb_3$.

(xii) If there is no transformer tap, proceed from (xiv), if the transformer tap is adjustable, proceed from (xiii), if not adjustable, proceed from (xv).

(xiii) $n_t = n_t + 1$

$$y_{td}^{nt} = b_1 + jb_2 + jb$$

$$y_t^{nt} = -b_1 - jb_2$$

$$I_n^{nt} = I_n = \text{near end}$$

$$I_f^{nt} = i_f = \text{far end}$$

t^{nt} = off nominal turn ratio.

Proceed from (xvi).

$$(xiv) \quad y_d^i = y_d^i + b_1 + jb_2 + jb$$

$$y_d^j = y_d^j + b_1 + jb_2 + jb$$

$$y_1^j = y_1^j - b_1$$

$$y_2^j = y_2^j - b_2$$

Proceed from (xvii).

(xv) Obtain the near end and far ends i_n and i_f respectively.

$$(xvi) \quad c_1 = b_1 t$$

$$c_2 = b_2 t$$

$$y_1^j = y_1^j - c_1$$

$$y_2^j = y_2^j - c_2$$

$$y_d^i = y_d^i + b_1 + jb_2 + jb$$

$$y_d^n = y_d^n + t(c_1 + jc_2) + jb t^2$$

(xvii) Check if this is the last cord of line data, if not proceed from (iv).

(xviii) Store the non-zero right off diagonal elements of i th row as in (vi) and (vii).

(xix) Search for L_t^k , $k = 1, \dots, n_t$, L_2^i , L_3^i for $i = 1, \dots, k_2$ and I_2^k , $k = 1, \dots, n$.

(xx) Read shunt admittance data and modify the diagonal elements.

CHAPTER IVPOWER FLOW EQUATIONS

In this chapter some properties of power flow equations are discussed. Using some of the identities, a lot of computer time can be saved in computing the system jacobian and powers.

4.1 Steady state relations

Under steady state power flow equations in terms of the admittance matrix may be written as follows:

$$P_i = V_i \sum_j V_j Y_{ij} \cos(\theta_i - \theta_j - \psi_{ij}) \quad (4.1)$$

$$Q_i = V_i \sum_j V_j Y_{ij} \sin(\theta_i - \theta_j - \psi_{ij}) \quad (4.2)$$

Using the notation of Vanness and Griffin⁵⁰:

$$H_{ij} = \frac{\partial P_i}{\partial \theta_j} \quad (4.3)$$

$$H_{ij} = V_j \frac{\partial P_i}{\partial V_j} \quad (4.4)$$

$$J_{ij} = \frac{\partial Q_i}{\partial \theta_j} \quad (4.5)$$

$$L_{ij} = V_j \frac{\partial Q_i}{\partial V_j} \quad (4.6)$$

Thus

$$H_{ij} = v_i v_j y_{ij} \sin(\theta_i - \theta_j - \psi_{ij}) \quad (4.7)$$

$$H_{ij} = v_i v_j y_{ij} \cos(\theta_i - \theta_j - \psi_{ij}) \quad (4.8)$$

$$J_{ij} = -v_i v_j y_{ij} \cos(\theta_i - \theta_j - \psi_{ij}) \quad (4.9)$$

$$L_{ij} = v_i v_j y_{ij} \sin(\theta_i - \theta_j - \psi_{ij}) \quad (4.10)$$

$$H_{ii} = -v_i \sum_{j \neq i} v_j y_{ij} \sin(\theta_i - \theta_j - \psi_{ij}) \quad (4.11)$$

$$H_{ii} = 2 v_i^2 b_{ii} + v_i \sum_{j \neq i} v_j y_{ij} \cos(\theta_i - \theta_j - \psi_{ij}) \quad (4.12)$$

$$J_{ii} = v_i \sum_{j \neq i} v_j y_{ij} \cos(\theta_i - \theta_j - \psi_{ij}) \quad (4.13)$$

$$L_{ii} = -2 v_i^2 b_{ii} + v_i \sum_{j \neq i} v_j y_{ij} \sin(\theta_i - \theta_j - \psi_{ij}) \quad (4.14)$$

From these equations, following identities result :

$$s_{ij} = e_{ij} s_{ij} - c_{ij} b_{ij} \quad (4.15)$$

$$s_{ij} = c_{ij} s_{ij} + e_{ij} b_{ij} \quad (4.16)$$

where,

$$s_{ij} = v_i v_j \sin(\theta_i - \theta_j) = f_i e_j - e_i f_j \quad (4.17)$$

$$c_{ij} = v_i v_j \cos(\theta_i - \theta_j) = e_i e_j + f_i f_j \quad (4.18)$$

$$J_{ij} = -H_{ij} \quad (4.19)$$

$$L_{ij} = H_{ij} \quad (4.20)$$

$$a_{11} = - \sum_{j \neq 1}^n a_{1j} \quad (4.21)$$

$$c_{11} = - \sum_{j \neq 1}^n c_{1j} \quad (4.22)$$

$$P_1 = v_1^2 e_{11} + J_{11} \quad (4.23)$$

$$C_1 = -v_1^2 b_{11} - H_{11} \quad (4.24)$$

$$E_{11} = v_1^2 e_{11} + P_1 \quad (4.25)$$

$$L_{11} = -v_1^2 b_{11} + C_1 \quad (4.26)$$

The jacobian ' H ' of the set of equations 4.3 to 4.6 is a square matrix of order $2n$:

$$H = \begin{bmatrix} M & N \\ J & L \end{bmatrix} \quad (4.27)$$

where M , N , J and L are $n \times n$ square matrix whose ij th elements are H_{ij} , N_{ij} , J_{ij} and L_{ij} , respectively.

4.2 Theorem

Whatever the values of V and θ ; matrices M , J and N are singular.

Proof:

From equations 4.21 and 4.22; sum of elements of any row of matrix M and J is zero. Hence the matrix M , J and N are singular.

4.3 Theorem

There do not exist in general vectors V and Θ such that the equations 4.1 and 4.2 are satisfied for any given value of P_i and Q_i ; $i = 1, \dots, n$.

Proof:

Let a set of variables V_0 , Θ_0 , P_0 and Q_0 satisfy equations 4.1 and 4.2. If the active and reactive powers are perturbed by ϵn_p and ϵn_q respectively where the scalar $\epsilon \rightarrow 0$ and n_p and n_q are vector quantities, such that

$$P = P_0 + \epsilon n_p$$

$$Q = Q_0 + \epsilon n_q$$

If corresponding per unit variation in voltage magnitude and variation in phase angles are respectively ΔV and $\Delta\Theta$; these are given by the following equations

$$\epsilon \begin{bmatrix} n_p \\ n_q \end{bmatrix} = K \begin{bmatrix} \Delta\Theta \\ \Delta V \end{bmatrix} \quad (4.28)$$

Since K is singular; $\Delta\Theta$ and ΔV cannot be obtained for any arbitrary n_p and n_q . Hence Θ and V cannot be obtained for any arbitrary P and Q .

4.4 Load flow problem

In order to be able to solve for bus bar voltages and angles; given the complex bus bar powers; phase angle of one bus bar must be taken as reference. This deletes one column of M and J . By taking active power of one bus as dependent variable, one row of M is deleted. This makes the matrix non-singular. However in order to be able to fix the order of bus bar voltage magnitudes; it is desirable that voltage magnitude of one bus be fixed. Thus three types of buses are represented in the load flow calculations.

Slack bus ($v \theta$): V and θ are prescribed and P and Q are the dependent variables. Selection of this bus is necessary for the reasons already mentioned.

Voltage regulated bus ($p v$): P and V are prescribed; Q and θ are the dependent variables.

Load bus ($p q$): P and Q are prescribed; V and θ are dependent variables.

4.5 Compatibility relations

Since M is a singular matrix it has zero as one of the eigen values; and let $(\lambda^t, \mu^t)^t$ be the eigen vector of M^t corresponding to this eigen value; where λ and μ are $n \times 1$ column vectors.

$$M \begin{bmatrix} \Delta \theta \\ \Delta V \\ \vdots \\ \Delta Q \end{bmatrix} = \begin{bmatrix} \Delta P \\ \Delta Q \\ \vdots \\ \Delta P \end{bmatrix} \quad (4.29)$$

After multiplying by (λ^T, μ^T) on both sides; following compatibility relations are satisfied:

$$\sum_i \lambda_i \Delta P_i + \sum_k \mu_k \Delta Q_k = 0 \quad (4.30)$$

4.6 General compatibility relation

Variational equations of the power flow problem can be written as follows:

$$\begin{bmatrix} \Delta P \\ \vdots \\ \Delta Q_k \\ \vdots \\ \Delta V_1 \end{bmatrix} = \begin{bmatrix} H & & & \\ & \ddots & & \\ J_{k1} & \dots & J_{kn} & L_{k1} & \dots & L_{kn} \\ & & & & & \\ 0 & 0 & 0 & \dots & 1 & \dots & 0 \end{bmatrix} \begin{bmatrix} \Delta \theta \\ \Delta V \end{bmatrix} \quad (4.31)$$

where $k \in$ set of bus bars with Q as independent variable

$l \in$ set of bus bars with V as independent variable.

The jacobian of equation 4.31 is also singular, hence the compatibility relation is as follows:

$$\sum_i \lambda_i \Delta P_i + \sum_k \mu_k \Delta Q_k + \sum_l u_l \Delta V_l = 0 \quad (4.32)$$

where $(\lambda^T, \mu^T, u^T)^T$ is the eigen vector of the transposed jacobian corresponding to zero eigen value.

4.7 Comments

All λ 's in equations 4.30 and 4.32 will have the same sign.

If only powers at two nodes i and j are changed by ΔP_i and ΔP_j and other incremental quantities do not undergo any change;

$$\begin{aligned} \Delta P_i \lambda_i + \Delta P_j \lambda_j &= 0 \\ \frac{P_i}{P_j} &= -\frac{\lambda_j}{\lambda_i} \end{aligned} \quad (4.33)$$

It is intuitively expected for a power system under normal loading that if power at one node increases, the power at the other node decreases, if all other powers remain unchanged. Hence $\Delta P_i / \Delta P_j < 0$. This is possible if all λ 's have the same sign.

4.8 Correlation between compatibility and loss formula relation

Through the application of tensor theory, the total system losses can be expressed in terms of all plant outputs, after a number of assumptions are made, as follows⁵³:

$$P_L = \sum_{i \in G} \sum_{j \in G} P_i B_{ij} P_j \quad (4.34)$$

Incremental changes in generation, with demand unchanged will result in the following incremental change in loss:

$$\begin{aligned} \sum_{i \in G} \Delta P_i &= \Delta P_L \\ &= 2 \sum_{i \in G} \left[\sum_{j \in G} S_{ij} P_j \right] \Delta P_i \\ \text{or } \sum_{i \in G} (1 - 2 \sum_{j \in G} S_{ij} P_j) \Delta P_i &= 0 \end{aligned} \quad (4.35)$$

Comparing equations 4.32 and 4.35;

$$\lambda_i = 1 - 2 \sum_{j \in G} S_{ij} P_j \quad (4.36)$$

CHAPTER V

LOAD FLOW SOLUTIONS

In this chapter, various load flow programmes are lumped into very general categories and compared. Sufficiency conditions for convergence of iterative procedures are discussed through the fixed point principle and applied to admittance matrix, impedance matrix and Newton's method for load flow.

5.1 Classification

Load flow programmes can be broken into the following broad categories⁹:

- (a) Low accuracy special purpose load flows³⁰: Such load flow programmes are needed where emphasis is not on accurate detailed modelling of power systems. Emphasis instead is on simple criteria such as what variables are out of limit and for what reasons. Usually this is followed by performing a series of input data changes to simulate possible system contingencies. For each such contingency the operator will like to know if any other components will be affected which may cause another system failure and eventually lead to a cascading system

breakdown. The contingencies studied are usually in the form of line outages, generator outages, short circuits and various other effects.

- (b) Conventional load flows: Such load flows need accurate set of input data and are run to appropriately small tolerance. The trade offs include programming convenience, speed and memory requirement.
- (c) Very accurate load flows: Many applications such as transient stability studies and static optimisation need in many cases a very accurate load flow programme.

5.2 Programming algorithms⁹

These algorithms for load flow solutions are almost as numerous as the number of authors who have written about load flows. These can broadly be lumped in the following categories:

- (a) Admittance matrix method: These methods use Gauss or Gauss Seidal iterative technique and system admittance matrix with ground as reference. Chief advantages of this method are ease of programming and most efficient utilization of core memory of any presently known load flow methods. It is usually used when a very large system need be studied on a computer with limited core memory. However it needs long running time and has possible non-convergence.

- (b) Impedance matrix method: This method also uses Gauss or Gauss Seidal iterative technique and system impedance matrix with one of the nodes as reference. Convergence of this method is better behaved than the admittance matrix method but Z-matrix is a full matrix while Y-matrix sparse. The author has been able to overcome the disadvantages of the full matrix by making use of L.U. factors, of the admittance matrix formed by taking slack bus as reference. This works much faster and needs much smaller storage if compact storage is used, compared to conventional Z-matrix. Many authors have suggested the application of 'piece-wise methods'^{6,22,23,24} to overcome the disadvantages of excessive storage requirement. The real advantages of this method are apparent when the user wishes to study the effect of short circuits, line and generator outages or changes in tie line flows.
- (c) Newton's method: This technique has the best convergence characteristics of all the presently available methods. Solution accuracy is restricted only by the round off errors. This method if efficiently programmed and buses optimally renumbered and compact storage used for the jacobian and admittance matrix, can be made as the fastest. This method is probably best suited for optimal power flow studies because of very high accuracy and the programme needs small modifications for obtaining incremental costs of adjustable

parameters. Main disadvantages of the method are

- (i) the programming logic is considerably more complex
- (ii) large memory requirement even if compact storage is used. Credit for making this technique practical goes to Tinney³³ and his associates of B.P.A.

- (d) Non-linear programming methods: Load flow problem is treated as the problem of optimization where the objective function is the sum of squares of the total real and reactive mismatch. Optimization is achieved through any of the gradient techniques. The author's experience with gradient techniques is that an accurate solution is hard to obtain. Fletchers Powell's gradient method⁴⁴ may yield a better convergence but needs a very large storage and time for the non-sparse Hessian matrix which is built iteratively. Chief advantages of this method appear to be only small modification needed for optimal power flow solution and less computer logic compared to Newton's method. Credit for perfecting this method goes to Casson^{4,15,16}.

Most of the research work in the field of load flow has been concerned with variation on these methods, which include better correction and acceleration methods, better ways to utilize the matrix by tearing, triangular factorization, compact storage, near optimal renumbering and various methods of representing the system as a d.c. circuit and dealing only with real power flows to obtain approximate solution.

5.3 Convergence of Gauss and Gauss-Seidel iterative techniques for linear simultaneous equations

If $n \times n$ linear simultaneous equation under consideration is as follows:

$$Ax = b \quad (5.1)$$

$$\text{Let } A = D + L + U \quad (5.2)$$

where D , L and U are $n \times n$ matrices. D contain the diagonal elements, L the elements on the left side of the diagonal and U the elements on the right hand side of diagonal.

Formulation of Gauss iterative procedure is as follows:

$$D x^{k+1} = b - (L + U) x^k \quad (5.3)$$

$$\text{if } x^k = x^* + \varepsilon \eta \quad (5.4)$$

where η is $n \times 1$ column vector and x^* is the solution of equation 5.1

$$\begin{aligned} D x^{k+1} &= A x^* - (L + U) (x^* + \varepsilon \eta) \\ &= D x^* - \varepsilon (L + U) \eta \\ x^{k+1} &= x^* - \varepsilon D^{-1} (L + U) \eta \end{aligned} \quad (5.5)$$

If absolute value of any eigen value of $D^{-1}(L + U)$ is more than 1, the Gauss iterative method will not converge, does not matter how close the initial guess is to the actual solution.

Formulation of Gauss Seidal iterative technique is as follows:

$$\mathbf{x}^{k+1} = \mathbf{x}^* - \varepsilon (\mathbf{L} + \mathbf{D})^{-1} \mathbf{U} \eta \quad (5.6)$$

Necessary and sufficient condition for the technique to converge is that the absolute value of no eigen value of $(\mathbf{L} + \mathbf{D})^{-1} \mathbf{U}$ be greater than 1.

Convergence may sometimes be improved by the application of acceleration or damping, as follows:

$$\mathbf{x}^{k+1} \text{ (modified)} = \mathbf{x}^k + \alpha(\mathbf{x}^{k+1} - \mathbf{x}^k) \quad (5.7)$$

where $\alpha > 0$

$\alpha > 1$ represents acceleration and $\alpha < 1$ represent damping.

$$\mathbf{x}^k = \mathbf{x}^* + \varepsilon \eta$$

$$\mathbf{x}^{k+1} = \mathbf{x}^* - \varepsilon \mathbf{B} \eta \quad (5.8)$$

where \mathbf{B} is the principal matrix and is as follows:

$$\mathbf{B} = \mathbf{D}^{-1}(\mathbf{L} + \mathbf{U}) \text{ for Gauss method}$$

$$\text{and } \mathbf{B} = (\mathbf{L} + \mathbf{D})^{-1} \mathbf{U} \text{ for Gauss Seidal method}$$

$$\mathbf{x}^{k+1} \text{ (modified)} = \mathbf{x}^* + \varepsilon [(1-\alpha) \mathbf{I} - \alpha \mathbf{B}] \eta \quad (5.9)$$

where \mathbf{I} is the identity matrix.

For the modified solution to converge:

$$|1 - \alpha(1 + \beta)| < 1 \quad (5.10)$$

where β is an eigen value of B .

Speed of convergence for unaccelerated solution depends upon β and for the accelerated solution upon $|1 - \alpha(1 + \beta)|$.

If $\beta > 1$, unaccelerated solution will not converge, but the damped solution satisfying equation (5.10) will converge.

If $\beta < -1$, no acceleration or damping can make it converge.

If $\beta = -1 + \delta$; for small positive δ ; convergence is poor but the accelerated solution will have better convergence.

5.4 Convergence of iteration process for non-linear functions ^{37,40}

If the equations are written in the following form,

$$x_i = F_i(x) ; i = 1, \dots, n \quad (5.11)$$

where $x = (x_1, x_2, \dots, x_n)^T$

First iteration is as follows:

$$x_1^2 = F_1(x^1) \quad (5.12)$$

$$\begin{aligned}
 x_1^* - x_1^2 &= F_1(x^*) - F_1(x^1) \\
 &= \int_0^1 \frac{\partial F_1}{\partial x} [F_1(x^1 + \epsilon(x^* - x^1))] d\epsilon \\
 &= \int_0^1 \sum_j \frac{\partial F_1}{\partial x_j} (x_j^1 + \epsilon(x_j^* - x_j^1)) (x_j^* - x_j^1) d\epsilon \\
 &= \sum_j (x_j^* - x_j^1) \int_0^1 \frac{\partial F_1}{\partial x_j} (x_j^1 + \epsilon(x_j^* - x_j^1)) d\epsilon
 \end{aligned}$$

Let $\epsilon_{ij}(x^*, x^1) = \int_0^1 \frac{\partial F_1}{\partial x_j} (x_j^1 + \epsilon(x_j^* - x_j^1)) d\epsilon$ (5.13)

$$x_1^* - x_1^2 = \sum_j (x_j^* - x_j^1) \epsilon_{ij} \quad (5.14)$$

Adding equation 5.14;

$$\sum_i (x_i^* - x_i^2) = \sum_i \sum_j (x_j^* - x_j^1) \epsilon_{ij} \quad (5.15)$$

$$= \sum_j (x_j^* - x_j^1) \sum_i \epsilon_{ij} \quad (5.16)$$

Considering only the absolute values

$$\sum_j |x_j^* - x_j^2| \leq \sum_j |x_j^* - x_j^1| + \sum_i |\epsilon_{ij}| \quad (5.17)$$

If $\sum_i \left| \frac{\partial F_1}{\partial x_j} \right| \leq \delta < 1$ for all i in the region

(x^1, x^*)

$$\sum_j |x_j^* - x_j^2| \leq \delta \sum_j |x_j^* - x_j^1| \quad (5.18)$$

This relation holds good for the first iteration. For successive iterations; similarly

$$\begin{aligned} \sum_j |x_j^* - x_j^{k+1}| &\leq \delta \sum_j |x_j^* - x_j^k| \\ \sum_j |x_j^* - x_j^{k+1}| &\leq \delta^k \sum_j |x_j^* - x_j^1| \end{aligned} \quad (5.19)$$

If δ is less than 1; right hand side of this inequality can be made as small as desirable, by repeating the iteration process sufficient number of times.

Thus the iteration process converges if

$$\sum_j \left| \frac{\partial F_j}{\partial x_1} \right| < 1 \quad (5.20)$$

for all i.

5.5 Uniqueness of solution

If x^* and y satisfy equation 5.13;

since $y_1 = F_1(y)$

$$\sum_j |x_j^* - F_j(y)| \leq \delta \sum_j |x_j^* - y_j|$$

$$\text{or } \sum_j |x_j^* - y_j| \leq \delta \sum_j |x_j^* - y_j|$$

This is a contradiction since $\delta < 1$.

5.6 Convergence of admittance matrix method using
Gauss iterative technique

$$v_i = \frac{1}{y_{ii}} \left[\frac{P_i - J v_i}{V_i} - \sum_{j \neq i} y_{ij} v_j \right] \quad (5.21)$$

for all i except slack bus.

$$\text{Let } g_{ij} + j b_{ij} = y_{ij}/y_{ii} \quad (5.22)$$

$$c_i + j d_i = (P_i - J v_i)/y_{ii} \quad (5.23)$$

$$e_i + j f_i = v_i \quad (5.24)$$

$$e_i = \frac{e_i c_i - f_i d_i}{e_i^2 + f_i^2} - \sum_{j \neq i} (g_{ij} e_j + b_{ij} f_j) \quad (5.25)$$

$$f_i = \frac{f_i c_i + e_i d_i}{e_i^2 + f_i^2} - \sum_{j \neq i} (b_{ij} e_j + g_{ij} f_j) \quad (5.26)$$

Criteria of convergence are

$$\delta_e = \sum_{\substack{i=1 \\ i \neq s}} \left| \frac{\partial e_i}{\partial e_k} \right| + \sum_{\substack{k=1 \\ k \neq s}} \left| \frac{\partial e_i}{\partial f_k} \right| < 1; \quad k \neq s \quad (5.27)$$

$$\delta_f = \sum_{\substack{i=1 \\ i \neq s}} \left| \frac{\partial f_i}{\partial e_k} \right| + \sum_{\substack{k=1 \\ k \neq s}} \left| \frac{\partial f_i}{\partial f_k} \right| < 1; \quad k \neq s \quad (5.28)$$

where e_i and f_i are the right hand side of equations 5.25 and 5.26.

$$\delta_e = \sum_{k \neq i, s} \left\{ |b_{ik}| + |e_{ik}| \right\}.$$

$$\frac{\left| f_k^2 c_k - e_k^2 c_k + 2e_k f_k d_k \right| + \left| -e_k^2 d_k + f_k^2 d_k - 2e_k f_k d_k \right|}{(e_k^2 + f_k^2)^2} \quad (5.29)$$

$$\delta_f = \sum_{k \neq i, s} \left\{ |b_{ik}| + |e_{ik}| \right\}$$

$$\frac{\left| f_k^2 d_k - e_k^2 d_k - 2f_k e_k c_k \right| + \left| e_k^2 c_k - f_k^2 c_k - 2e_k f_k d_k \right|}{(e_k^2 + f_k^2)^2} \quad (5.30)$$

For a reliable convergence δ_e and δ_f should be sufficiently less than 1. However as we know

$$- \sum_j y_{ij} + y'_i = y_{ii} \quad (5.31)$$

where y'_i is the equivalent shunt admittance at i node.

this suggest

$$\left| \sum_{j \neq i} \frac{y_{ij}}{y_{ii}} \right|$$

is nearly equal to 1.

Hence $\sum_{k \neq i} |b_{ik}| + |e_{ik}|$ and subsequently

$$\sum_{k \neq i, s} |b_{ik}| + |e_{ik}|$$

may be on either side of 1. Over and above it, there will be contribution due to additional term depending on system loading. This means that the sufficiency condition for convergence is either not satisfied or satisfied by a low margin. This explains for poor and less reliable convergence of the method.

5.7 Convergence of impedance matrix method with Gauss iterative technique

The fixed point formulation is as follows:

$$I_i = \frac{P_i - J S_i}{V_i} - y_i V_i \quad (5.32)$$

$$V_i = V_s + \sum_{\substack{j \\ j \neq s}} z_{ij} I_j \quad (5.33)$$

These equations may also be written as follows:

$$V_i = V_s + \sum_{\substack{j \\ j \neq s}} z_{ij} \left[\frac{P_j - J S_j}{(V_j)^0} - y_j V_j \right] \quad (5.34)$$

for all i except slack.

$$\text{If } r_{ij} + J z_{ij} = z_{ij} \quad (5.35)$$

$$e_j + J f_j = V_j \quad (5.36)$$

$$z_{ij} + J b_{ij} = z_{ij} y_j \quad (5.37)$$

$$e_1 = e_e + \sum_{\substack{j \\ j \neq 0}} \frac{e_j(r_{1j} p_j + x_{1j} q_j) + f_j(x_{1j} p_j - r_{1j} q_j)}{(e_j^2 + f_j^2)}$$

$$= \sum_{\substack{j \\ j \neq 0}} (r_{1j} e_j - b_{1j} f_j) \quad (5.38)$$

$$f_1 = f_e + \sum_{\substack{j \\ j \neq 0}} \frac{e_j(x_{1j} p_j - r_{1j} q_j) - f_j(r_{1j} p_j + x_{1j} q_j)}{(e_j^2 + f_j^2)}$$

$$= \sum_{\substack{j \\ j \neq 0}} (b_{1j} e_j + r_{1j} f_j) \quad (5.39)$$

Let

$$a'_{ik} = r_{ik} p_k + x_{ik} q_k \quad (5.40)$$

$$b'_{ik} = x_{ik} p_k - r_{ik} q_k \quad (5.41)$$

$$c_{ik} = e_k a'_{ik} + f_k b'_{ik} \quad (5.42)$$

$$d_{ik} = e_k b'_{ik} - f_k a'_{ik} \quad (5.43)$$

$$\varepsilon_e = \sum_{\substack{k \\ k \neq 0}} \left| \frac{(e_k^2 + f_k^2) a'_{ik} - 2e_k c_{ik}}{(e_k^2 + f_k^2)^2} - g_{ik} \right|$$

$$\sum_{\substack{i \\ i \neq k}} \left| \frac{(e_k^2 + f_k^2) b'_{ik} - 2e_k d_{ik}}{(e_k^2 + f_k^2)^2} + b_{ik} \right| \quad (5.44)$$

$$\delta_f = \sum_{\substack{i \\ i \neq k}} \left| \frac{(e_k^2 + f_k^2) b'_{ik} - 2f_k c_{ik}}{(e_k^2 + f_k^2)^2} + b_{ik} \right|$$

$$\sum_{\substack{i \\ i \neq k}} \left| \frac{-(e_k^2 + f_k^2) e'_{ik} - 2 f_k d_{ik}}{(e_k^2 + f_k^2)^2} - g_{ik} \right| \quad (5.45)$$

g_{ik} and b_{ik} are small quantities due to small shunt admittance at node k . Other elements in δ_e and δ_f are proportionate to loading. If loading is small, δ_e and δ_f will be less than 1. This assures the convergence. This explains for better convergence of Z-bus method over Y-bus method.

5.6 Convergence of Newton's method

Consider a set of n equations with n unknowns

$$f_1(x_1, x_2, \dots, x_n) = 0$$

The fixed point formulation through Newton's method is as follows:

$$x = x - J^{-1}(x) f(x) = \beta(x) \quad (5.46)$$

$$\text{where } J(x) = \begin{bmatrix} \frac{\partial f_1}{\partial x_1} & \dots & \frac{\partial f_1}{\partial x_n} \\ \vdots & \ddots & \vdots \\ \frac{\partial f_n}{\partial x_1} & \dots & \frac{\partial f_n}{\partial x_n} \end{bmatrix} \quad (5.47)$$

Because of the presence of $J(x)$; analysis through 5.4 will need a three dimensional Hessian matrix; explanation of resulting expression is hard to understand. Kantorovich's⁴⁰ analysis, attempted by Neisel and Barnard³ also has the same drawback. A simplified analysis is as follows.

$$\begin{aligned} x^{k+1} - x^* &= x^k - J^{-1}(x^k) f(x^k) - x^* \\ &= x^k - x^* - J^{-1}(x^k) [f(x^k) - f(x^*)] \end{aligned} \quad (5.48)$$

where x^* is solution of the equation

$$f(x^k) - f(x^*) = G(x^*, x^k) (x^k - x^*) \quad (5.49)$$

where G is $n \times n$ matrix whose i,j th element is given by equation 5.13

$$\begin{aligned} x^{k+1} - x^* &= x^k - x^* - J^{-1}(x^k) G(x^*, x^k) (x^k - x^*) \\ &= \{ I - J^{-1}(x^k) G(x^*, x^k) \} (x^k - x^*) \end{aligned} \quad (5.50)$$

Since

$$J^{-1}(x^*) G(x^*, x^*) = I \quad (5.51)$$

If x^1 is close enough to x^* , matrix $J^{-1}(x^1)G(x^*, x^1)$ is quite close to I and $I - J^{-1}(x^1)G(x^*, x^1)$ is quite close to zero.

If the norm of a column vector y is defined as follows

$$\|y\| = \sum_j |y_j| \quad (5.52)$$

and norm of matrix A as follows

$$\|A\| = \max_j \sum_i |a_{ij}| \quad (5.53)$$

$$\|\Delta y\| = \sum_i \left| \sum_j a_{ij} y_j \right| \leq \sum_i \sum_j |a_{ij}| |y_j|$$

$$\leq \sum_j |y_j| \max_j \sum_i |a_{ij}|$$

$$\leq \|y\| \|A\| \quad (5.54)$$

Thus

$$\|x^2 - x^*\| \leq \|I - J^{-1}(x^1) G(x^*, x^1)\| + \|x^1 - x^*\|$$

Since norm of a zero matrix is zero and since

$I - J^{-1}(x^1) G(x^*, x^1)$ is quite close to zero, there therefore exists $\epsilon > 0$ such that if $\|x^1 - x^*\| < \epsilon$

Then

$$\|I - J^{-1}(x) G(x^*, x)\| < \delta < 1$$

If $\|x^1 - x^*\| < \epsilon$ then

$$\|x^2 - x^*\| \leq \delta \|x^1 - x^*\| + \delta \epsilon < \epsilon \quad (5.55)$$

Therefore,

$$\|x^{k+1} - x^*\| \leq \delta^k \|x^1 - x^*\|$$

and $\lim_{k \rightarrow \infty} \|x^{k+1} - x^*\| = 0 \quad (5.56)$

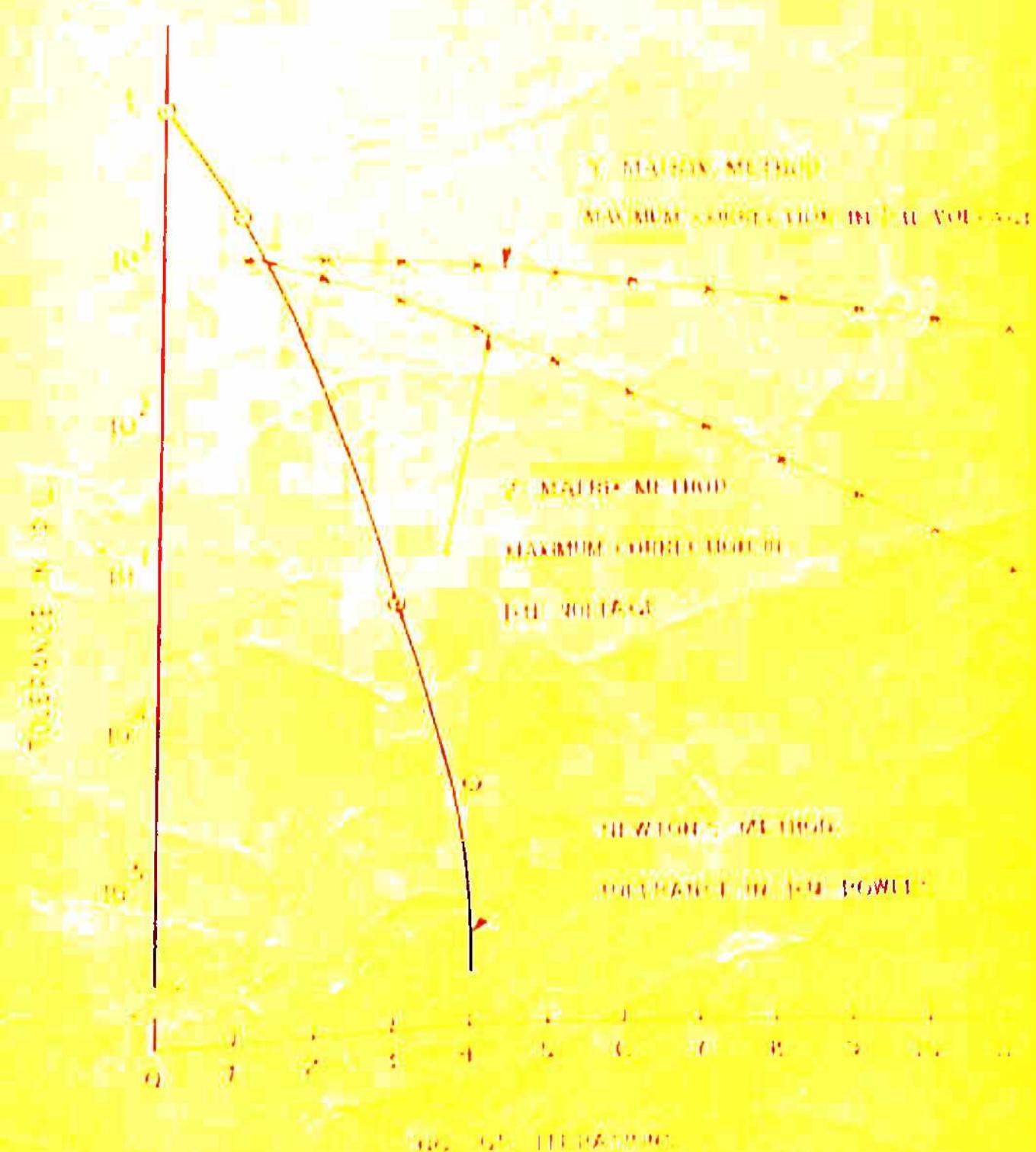
5.9 Example

The value of δ obtained on AEP 30 bus system for Y-matrix method was 1.126. This indicates doubtful convergence. However it converged to a voltage tolerance of 0.001 pu in 44 iterations with the initial voltages of $1 + j0$ for all buses except $1.06 + j0$ for the slack bus.

The same value for Z-matrix method was 0.869. This indicates sure convergence. The solution converged in 11 iterations within the tolerance indicated above.

For Newton's method the value of δ and maximum mismatch for successive iterations was as follows:

| <u>Iteration</u> | <u>δ</u> | <u>Maximum mismatch</u> |
|------------------|----------------------------|-------------------------|
| Initial start | .3863 | .94199 |
| 1 | .1254 | .20980 |
| 2 | .0327 | .26968 $\times 10^{-1}$ |
| 3 | .0043 | .73812 $\times 10^{-4}$ |
| 4 | .0002 | .53589 $\times 10^{-5}$ |



Comparison of convergence rate of matrix method with that of the Y-MATRIX METHOD

$$1 - C_1 \cdot C_2 \cdot \frac{1}{\alpha}$$

Maximum mismatch of fourth iteration is not considered reliable since the word size of IBM 1130 computer with standard precision is about 6.9 decimal digits. This should be much lower as judged from the trend and also from the corresponding S.

5.10 Conclusion

Hence following conclusions are drawn through fixed point principles to which the experimental results also agree.

- (1) Convergence of Y-matrix method is less reliable even if the initial point is close enough and loading is small.
- (2) Convergence of Z-matrix method is more reliable if the loading and shunt admittance is small.
- (3) Convergence of Newton's method is very reliable if the initial point is sufficiently close to the final solution.

It has been experimentally established in literature that the convergence of Newton's method in polar form is most reliable and fast. Convergence characteristics of load flow methods for AEP 30 bus system is as given in Fig. 5.1. A similar characteristics is also given by Dommel in correspondence for reference 33 for a 17 bus system.

CHAPTER VI

COMPUTER PROGRAMME FOR LOAD FLOW

SOLUTION BY NEWTON'S METHOD

In this chapter the load flow programme written by the author is discussed. It has not been possible for us to compare the computational time per iteration with that of BPA programme because of availability of a small computer and non-availability of BPA system test data to us. Time per iteration for AEP 30 bus system is about 10 seconds.

The programme is based on formulating and triangularizing the equation 6.1 node by node. This equation is not stored in the computer. The notation followed is that of Vanness and Griffin^{5C} given in Chapter IV. If incremental costs of maintaining the variables to the prescribed value need not be computed directly, L_{is} and N_{is} for $i \neq s$ are replaced by zeros. Similarly for decidedly pv node numbered k; L_{ik} and N_{ik} are also replaced by zero. Since the programme has been used for optimal power flow studies, this has not been done since incremental cost of maintaining the voltage to a particular value also need be computed.

6.1 Changing bus bar type

In addition to the conventional buses of p,q, p,v and slack; the computer programme provides for the following additional types of buses:

| ΔP_1 | $\Delta \theta_1$ | $\Delta V_1/V_1$ | $\Delta \theta_n$ | $\Delta V_n/V_n$ |
|---|------------------------------------|---|---|---|
| $b_{11} x_{11} + \dots + b_{1n} x_{1n}$ | $b_{11} \dots + b_{1n} x_{1n}$ | $b_{11} x_{11} + \dots + b_{1n} x_{1n}$ | $b_{nn} x_{nn} + \dots + b_{1n} x_{1n}$ | $b_{nn} x_{nn} + \dots + b_{1n} x_{1n}$ |
| $d_{11} L_{11} + \dots + d_{1n} L_{1n}$ | $L_{11} \dots + L_{1n} x_{1n}$ | $L_{11} \dots + L_{1n} x_{1n}$ | $L_{nn} \dots + L_{1n} x_{1n}$ | $L_{nn} \dots + L_{1n} x_{1n}$ |
| \vdots | \vdots | \vdots | \vdots | \vdots |
| $c_1 c_2 \dots c_n$ | $c_1 \dots c_n$ | $c_1 \dots c_n$ | $c_n \dots c_1$ | $c_n \dots c_1$ |
| \vdots | \vdots | \vdots | \vdots | \vdots |
| $b_{kk} x_{kk} + \dots + b_{kn} x_{kn}$ | $x_{kk} \dots + x_{kn} b_{kn}$ | $b_{kk} x_{kk} + \dots + b_{kn} x_{kn}$ | $b_{nn} x_{nn} + \dots + b_{kn} x_{kn}$ | $b_{nn} x_{nn} + \dots + b_{kn} x_{kn}$ |
| $d_{nn} L_{nn} + \dots + d_{kn} L_{kn}$ | $L_{nn} \dots + L_{kn} d_{kn}$ | $L_{nn} \dots + L_{kn} d_{kn}$ | $L_{nn} \dots + L_{kn} d_{kn}$ | $L_{nn} \dots + L_{kn} d_{kn}$ |
| $\text{slack bus } (n)$ | | | | |
| $P_{\text{bus}}^{\text{bus}}$ | $\theta_{\text{bus}}^{\text{bus}}$ | | | |

- (i) Bus bar with assigned reactive power provided the bus voltage is within the prescribed upper and lower limits.
- (ii) Bus bar with assigned voltage provided the reactive power is within the prescribed upper and lower limits.

This has been done as follows for kth bus in each iteration:

Obtain L_{kj} for $j \neq k$ from equations 4.15, 4.17, 4.18 and 4.20.

$$Q_k = -v_k^2 b_{kk} + \sum_{j \neq k} L_{kj} \quad (6.2)$$

$$L_{kk} = Q_k - v_k^2 b_{kk} \quad (6.3)$$

If the bus is with assigned voltage magnitude, desired change in voltage ΔV_k is obtained.

With this change in voltage, expected change in Q_k is approximately as follows:

$$Q_k \approx (\Delta V_k L_{kk}) / v_k \quad (6.4)$$

$(Q_k + \Delta Q_k)$ is now checked if this is within the prescribed bounds, the bus is treated as PV bus; if not it is treated as PQ bus with prescribed reactive power equal to the upper or lower limit, depending upon which limit is violated for that iteration.

If k th bus is with prescribed reactive power, ΔQ_k is obtained from the prescribed reactive power and the calculated reactive power.

Approximate value of ΔV_k is obtained as follows:

$$\Delta V_k = (V_k \Delta Q_k) / L_{kk} \quad (6.5)$$

$(V_k + \Delta V_k)$ is now checked for the prescribed limits of bus bar voltage. If the limit is violated, the bus is treated as p v bus with prescribed voltage equal to the upper or lower limit, depending upon which limit is violated, otherwise it is a pq bus for that iteration.

Obtaining L_{kj} does not mean additional computation time for p v bus since L_{kj} will have to be computed and $L_{kj} = H_{kj}; j \neq k$.

Convergence with the above method has been found very satisfactory.

6.2 Formation of working row

In order to reduce the dimensionality of the problem, following notation has been used.

If the new number of the bus is i and its external number is i' ; for i th node, if it a pq;

$$\Delta S_{2i-1} = \Delta P_{i'} \quad (6.6)$$

$$\Delta s_{21} = \Delta \theta_1, \quad (6.7)$$

(before triangularisation)

and

$$\Delta s_{2i-1} = \Delta \theta_i, \quad (6.8)$$

$$\Delta s_{21} = \Delta v_i / v_i, \quad (6.9)$$

(after triangularization)

$$a_{2j-1}^1 = h_{1,j} = s e_{1,j} + c b_{1,j} \quad (6.10)$$

$$a_{2j}^1 = h_{1,j} = c e_{1,j} + s b_{1,j} \quad (6.11)$$

where $s = v_i, v_j, \sin(\theta_i - \theta_j)$

$$= e_{1,j} - e_{1,i} f_{ji} \quad (6.12)$$

$$c = v_i, v_j, \cos(\theta_i - \theta_j)$$

$$= e_{1,i} e_{1,j} + f_{1,i} f_{1,j} \quad (6.13)$$

$$a_{2j-1}^2 = j_{1,j} = -a_{2j}^1 \quad (6.14)$$

$$a_{2j}^2 = L_{1,j} = a_{2j-1}^1 \quad (6.15)$$

Equations 6.10 to 6.15 are valid for all j except $j \neq 1$.

$$a_{2i-1}^1 = - \sum_{j \neq 1} a_{2j-1}^1 = h_{1,i} \quad (6.16)$$

$$a_{2i-1}^2 = \sum_{j \neq 1} a_{2j}^1 = j_{1,i} \quad (6.17)$$

$$r_{11} = V_1^2, s_{1111} + a_{21-1}^2 \quad (6.18)$$

$$-r_{11} = -V_1^2, s_{1111} - a_{21-1}^1 \quad (6.19)$$

$$a_{21}^1 = V_1^2, s_{1111} + r_{11} \quad (6.20)$$

$$a_{21}^2 = -V_1^2, s_{1111} + r_{11} \quad (6.21)$$

These elements are obtained without making use of trigonometrical functions.

H_{11} and J_{11} are now replaced by zero.

New numbers are used only in forming and triangularising equation 6.1, while admittance matrix, voltages and powers are recognised by the external numbers. However a little time per iteration can be saved by reforming the admittance matrix with the new numbers and recognising all the bus bar quantities (voltages, powers, etc.) with the new numbers.

For a PV bus; whether predecided or decided during iteration; equation 6.7 is replaced by the following equation:

$$a_{21} = V_1/V_{11} \quad (6.22)$$

the second row for that node is as follows:

$$a_{21}^2 = 1 \quad (6.23)$$

$$a_j^2; j = 1, \dots, 2n; j \neq 21 = 0 \quad (6.24)$$

This working row represented by equations 6.10 to 6.17; 6.20 and 6.21 is not represented explicitly. The total number of non-zero elements, column location of each element and the numerical value of these elements are stored instead. This can be explained as follows.

Suppose the node 4 under consideration has non-zero element in the admittance matrix at column 2, 5 and 8.

The working row formed are as follows.

Table 6.1

| S.No. | 1 | 2 | 3 | 4 | 5 | 6 | 7 | 8 |
|---------------------------------------|---------|---------|---------|---------|---------|------------|------------|------------|
| Non-zero element of first working row | a_3^1 | a_4^1 | a_7^1 | a_8^1 | a_9^1 | a_{10}^1 | a_{15}^1 | a_{16}^1 |
| Column location | 3 | 4 | 7 | 8 | 9 | 10 | 15 | 16 |

Table 6.2

| S.No. | 1 | 2 | 3 | 4 | 5 | 6 | 7 | 8 |
|--|---------|---------|---------|---------|---------|------------|------------|------------|
| Non-zero element of second working row | a_3^2 | a_4^2 | a_7^2 | a_8^2 | a_9^2 | a_{10}^2 | a_{15}^2 | a_{16}^2 |
| Column location | 3 | 4 | 7 | 8 | 9 | 10 | 15 | 16 |

Non-zero elements in the first working row $n_1 = 8$

Non-zero elements in the second working row $n_2 = 8$

Table 6.2 need not be prepared if it is a PV node.
If it is a slack node none of these rows need be prepared.

6.3 Triangularization

After the working rows are formed; the first working row is triangularised with the help of table of factors. For this fourth node under consideration; the first column of non-zero element is 3. Assuming that the non-zero elements in the third row of table of factors are in column 5, 6, 9 and 10; the working is now modified as follows.

| Col.no. | 1 | 2 | 3 | 4 | 5 | 6 | 7 | 8 | 9 | 10 |
|------------------|--------|--------|--------|--------|--------|--------|--------|---------|---------|---------|
| Non-zero element | a'_3 | a'_4 | a'_5 | a'_6 | a'_7 | a'_8 | a'_9 | a'_10 | a'_15 | a'_16 |
| Column | 3 | 4 | 5 | 6 | 7 | 8 | 9 | 10 | 15 | 16 |

Non-zero elements $m_1 = 10$

Further modifications will be made to this working row while this is triangularized with row 4, 5, 6, etc. For this example no new non-zero elements will be added during triangularisation with the 4th row.

After the triangularisation is complete; contents of this array are transferred to the table of factors. Contents of second working row (Table 6.2) are now transferred to the

first working row, and the triangularisation is repeated for the 8th row corresponding to the node 4.

This process is repeated for every node. Corrections in voltage magnitude and angles are obtained through backward substitution. The iteration is repeated till the maximum mismatch is less than the prescribed error. If maximum mismatch of one iteration is found more than that of previous iteration, divergence in the solution has been assumed. The iteration may then be repeated with modified generation and tap settings. In the programme maximum mismatch is obtained only after calculation of jacobian elements and forward triangularization is complete. Since jacobian elements are obtained while computing the calculated powers; it was considered wasteful to check for mismatch before the triangularization starts.

Non-zero left subdiagonal elements need not be stored if only load flow solution is needed, without skipping the formation of jacobian during any iteration. However as will be seen later; the effect of $(\Delta t)^{-1}$ is also to be obtained, for which the left subdiagonal elements are also needed. Compact storage scheme of table 2.2.3 has been used for the table of factors.

The load flow solution can be made even faster if it does not have a bus with indefinite mode of operation as explained in section 6.1, by adopting the storage scheme of table 2.2.4 and exploiting the symmetrical pattern of non-zero

elements of the jacobian, which are obtained as follows:

$$H_{ij} = sg_{ij} - cb_{ij}$$

$$H_{ji} = -sg_{ij} - cb_{ij}$$

$$N_{ij} = cg_{ij} + sb_{ij}$$

$$N_{ji} = cg_{ij} - sb_{ij}$$

This way computation time of obtaining jacobian elements can be reduced practically to half, but triangularization time remains the same. However this scheme needs storage for the elements on the left of diagonal even if only load flow solution is needed.

CHAPTER VII

NONLINEAR PROGRAMMING

In this chapter description of the nonlinear programming problem, criteria and techniques of identifying the optimal point are discussed. Initially the discussion is restricted to the unconstrained problem of minimising the scalar $f(x)$. Effect of constraint is included later on.

7.1 Description of the problem

$f(x)$ is a scalar function of n-dimensional vector x . Subject to the following constraints;

$$h(x) = 0 \quad (7.1)$$

$$\text{and} \quad g(x) \leq 0 \quad (7.2)$$

a minimum of $f(x)$ is to be obtained.

Where $g(x)$ and $h(x)$ are vector functions of x .

7.2 The unconstrained solution

If $f(x)$ is a twice continuously differential function on \mathbb{S}^n ; then

$$\nabla f(x^*) = 0 \quad (7.3)$$

if x^* minimises f over \mathbb{R}^n and the quadratic form

$$\sum_i \sum_j \frac{\partial^2 f}{\partial x_i \partial x_j} a_i a_j \geq 0 \quad (7.4)$$

for every point x in the neighbourhood of x^* .

The techniques for finding the minimum (if it exists) may be based upon solving the set of simultaneous equations 7.3. Solution may be obtained through iterative techniques such as Newton's, Gauss or Gauss-Seidal methods if the equations are nonlinear. In order that such techniques work, it is necessary that the guess is such that $f(x)$ is convex and remains in convex domain from iteration to iteration.

7.3 Second order gradient method

The second order gradient method is based upon obtaining the solution of equation 7.3 through Newton's method. Fixed point formulation is as follows:

$$x = x - (f_{xx}(x))^{-1} \nabla f(x) \quad (7.5)$$

In order that the algorithm converges to the minimum, Hessian matrix $f_{xx}(x)$ must be positive definite and norm of $(I - f_{xx}^{-1}(x^*) f_{xx}(x))$ should be less than 1 in the ε neighbourhood of x^* in which the initial point lies.

7.4 First order gradient methods

Assuming we are at a point x in \mathbb{R}^n and we wish to move

along some direction vector d also in \mathbb{R}^n a certain distance r . The objective is to successfully approach the minimum. In order to achieve rapid convergence the direction d and distance r must be properly decided.

$$f(x + rd) = f(x) + r \sum_1^n \frac{\partial f}{\partial x_i} d_i + \frac{r^2}{2} \sum_1^n \sum_j \frac{\frac{\partial^2 f}{\partial x_i \partial x_j}}{d_i d_j} d_i d_j \quad (7.6)$$

If x is the previous iteration point and we wish to compute the new iteration point $x+rd$ which is nearer to the minimum, we must then have

$$f(x+rd) - f(x) < 0 \quad (7.7)$$

Since the coefficient of r^2 in equation 7.6 is positive semidefinite if $f(x)$ is convex; coefficient of r should be negative or d is such that

$$\sum_1^n \frac{\partial f}{\partial x_i} d_i < 0 \quad (7.8)$$

Optimal value of r may be obtained by equating the partial derivative of $f(x+rd)$ with respect to r to zero or

$$r^* = - \frac{\nabla f^T d}{d^T f_{xx} d} \quad (7.9)$$

7.5 Steepest descent method

In this method

$$d = -\nabla f(x)$$

Thus equation 7.8 is satisfied. r is obtained such that $f(x+rd)$ is minimum along the direction d . If $f(x)$ is quadratic, it may be obtained from equation 7.9.

For a quadratic convex function,

$$f(x) = a + b^T x + \frac{1}{2} x^T C x \quad (7.10)$$

It is proved in reference (44) that

$$(s^k)^T (s^k) \leq \frac{2}{m} (1 - m/M)^k (f(x^0) - f(x^*)) \quad (7.11)$$

where x^0 is the initial point and

$$s^k = x^k - x^0$$

m and M are respectively the minimum and maximum of non-zero eigen values of C . Since the function is convex, C is positive semidefinite, which implies that each eigen value is also positive semidefinite.

Thus,

$$\lim_{k \rightarrow \infty} s^k = 0 \quad (7.12)$$

Another important conclusion drawn by author is that a function with larger m/M will have a better convergence. One method to improve this ratio is scaling. This may be illustrated with the help of following example,

$$f(x) = x_1^2 + 4x_2^2$$

Hessian matrix for this example is

$$\begin{bmatrix} 2 & 0 \\ 0 & 8 \end{bmatrix}$$

for which $m/R = .25$.

If the initial guess is $(1,1)^t$, the direction vector is $(-2, -8)^t$ for which

$$r^0 = .1308$$

$$x^1 = (.7384, -.0464)^t$$

$$r^1 = .426$$

$$x^2 = (.110, .1116)^t$$

If the variables are scaled as follows

$$y_1 = x_1$$

$$y_2 = 2x_2$$

$$f(x) = f(y) = y_1^2 + y_2^2$$

for which the corresponding initial point is $(1, 2)^t$, the direction vector $(-2, -4)^t$, Hessian matrix is

$$\begin{bmatrix} 2 & 0 \\ 0 & 2 \end{bmatrix}$$

for which $m/R = 1$

$$r^* = -5$$

$$x^* = (0, 0)^t$$

which is obviously the optimal point.

7.6 Conjugate gradient method⁴³

While convergence of steepest descent method depends upon n/k ; the conjugate gradient method guarantees convergence to the exact solution, disregarding the round off errors, in n iterations for a quadratic function of equation 7.10. In this method, a set of linearly independent directors $d^0, d^1, d^2, \dots, d^{n-1}$ are generated such that these are C conjugate ($d_i^t \circ d_j = 0$ if $i \neq j$) and the iteration is given by

$$x^{k+1} = x^k + r^k d^k \quad (7.13)$$

where r^k is for minimum of $f(x^k + rd^k)$.

The algorithm as given in reference (43) is as follows:

$$d^k = -\nabla f(x^k) + \alpha^k d^{k-1} \quad (7.14)$$

where $\alpha^k = \frac{\nabla f^t(x^k) \cdot \nabla f(x^k)}{\nabla f^t(x^{k-1}) \cdot \nabla f(x^{k-1})} \quad (7.15)$

$$d^0 = -\nabla f(x^0) \quad (7.16)$$

x^{k+1} is obtained from equation 7.13.

This may be explained for the example of (7.5) as follows:

$$f(x) = x_1^2 + 4x_2^2$$

$$x^0 = (1, 1)^t$$

$$\nabla f(x^0) = (2, 8)^t$$

$$d^0 = (-2, -8)^t$$

$$r^0 = .1308$$

$$x^1 = (.7384, -.0464)^t$$

$$\nabla f(x^1) = (1.4768, -.3712)^t$$

$$\alpha = .0343$$

$$d^1 = (-1.5454, +.0968)^t$$

$$r^1 = .478$$

$$x^2 = (0, 0)^t$$

7.7 Fletcher and Powell's method⁴⁴

This gradient method has been found more powerful than other first order gradient methods, but need additional computer time and memory for the non-sparse inverse of Hessian, which is not evaluated directly as is done in the second order gradient method. Instead it is built iteratively. For

a quadratic function of equation 7.10, it is shown to converge in n iterations as is the case with conjugate gradient method.

The algorithm is as follows:

$$d^k = -H^k \nabla f(x^k) \quad (7.17)$$

where H^k is the positive semidefinite symmetrical matrix, H^0 may be chosen as identity matrix.

$$x^{k+1} = x^k + r^k d^k \quad (7.18)$$

where the scalar r^k is chosen to minimize $f(x^k + r^k d^k)$.

$$H^{k+1} = H^k + A^k + B^k \quad (7.19)$$

where A^k and B^k are square symmetrical matrices as follows:

$$A^k = \frac{r(d^k)(d^k)^t}{(d^k)^t(y^k)} \quad (7.20)$$

$$B^k = -\frac{(H^k y^k)(H^k y^k)^t}{(y^k)^t H^k y^k} \quad (7.21)$$

$$y^k = \nabla f(x^{k+1}) - \nabla f(x^k) \quad (7.22)$$

7.6 Single dimensional minimisation^{43,44}

The first order gradient methods mentioned above need the step length r along the direction d^k which minimises

$f(x^k + rd^k)$. Techniques used are either an interpolative procedure or a symmetrical search such as Fibonacci method^[10].

7.8.1 Parabolic extrapolation

$$y(r) = f(x^k + rd^k) \quad (7.23)$$

$$y'(r) = (d^k)^t \nabla f(x^k + rd^k) \quad (7.24)$$

assuming $y(r)$ as a quadratic function of r

$$y(r) = a_1 r^2 + a_2 r + a_3 \quad (7.25)$$

$$y'(r) = 2a_1 r + a_2$$

$$r^* = -a_2/2a_1 \quad (7.26)$$

a_1 , a_2 and a_3 of equation 7.25 may be estimated by obtaining $y(0)$, $y'(0)$ and $y(a)$ for some $r = a$ as follows:

$$y(0) = a_3$$

$$y'(0) = a_2$$

$$y(a) = a_1 a^2 + a_2 a + a_3$$

$$a_1 = \frac{y(a) - y(0) - ay'(0)}{a^2}$$

$$r^* = \frac{-y'(0) a^2}{2(y(a) - y(0) - ay'(0))} \quad (7.27)$$

where r^* is the estimated value of r^* .

Tommel and Tinney²¹ have used this method.

7.8.2 Savindon's cubical interpolation^{23,44}

In this method $y'(r)$ obtained from (7.24) is examined for $r = 0, h, 2h, 4h, \dots, a, b$ where r is doubled each time and b is the first of these values at which either y' is non negative or y has not decreased. It then follows $a < r^* < b$.

Defining

$$s = 3 \frac{y(a) - y(b)}{b - a} + y'(a) + y'(b) \quad (7.28)$$

and

$$w = \sqrt{s^2 - y'(a)y'(b)} \quad (7.29)$$

$$r_e = b - \left(\frac{y'(b) + w - s}{y'(b) - y'(a) + 2w} \right) (b-a) \quad (7.30)$$

If neither $y(a)$ nor $y(b)$ is less than $y(r_e)$, r_e is an estimate of r^* , otherwise the interpolation needs be repeated over the subinterval (a, r_e) or (r_e, b) according to as $y'(r_e)$ is +ve or -ve.

7.9 Optimization under equality constraints

If some of the elements of x could be considered as independent variables u , the problem may be restated as follows:

$$\min f(x, u) = 0 \quad (7.31)$$

$$\text{subject to } h(x, u) = 0 \quad (7.32)$$

a set of dependent variables x may be obtained from equation 7.32 for a set of independent variables u .

$$\frac{\partial f(x, u)}{\partial u_k} = \frac{\partial f(x, u)}{\partial u_k} + \sum_j \frac{\partial f(x, u)}{\partial x_j} \frac{\partial x_j}{\partial u_k} \quad (7.33)$$

$$\frac{\partial h_1(x, u)}{\partial u_k} = \frac{\partial h_1(x, u)}{\partial u_k} + \sum_j \frac{\partial h_1(x, u)}{\partial x_j} \frac{\partial x_j}{\partial u_k} = 0 \quad (7.34)$$

From equation 7.34 following matrix equation results

$$\left(\frac{\partial x_j}{\partial u_k} \right) = -\left(\frac{\partial h_1(x, u)}{\partial x_j} \right)^{-1} \left(\frac{\partial h_1(x, u)}{\partial u_k} \right) \quad (7.35)$$

where

$$\left(\frac{\partial x_j}{\partial u_k} \right) \text{ is a column vector whose } \\ \text{jth element} = \frac{\partial x_j}{\partial u_k} \quad (7.36)$$

$$\left(\frac{\partial h_1(x, u)}{\partial x_j} \right) \text{ is a square matrix whose } \\ \text{ijth element} = \frac{\partial h_1(x, u)}{\partial x_j} \quad (7.37)$$

$$\left(\frac{\partial h_1(x, u)}{\partial u_k} \right) \text{ is a column vector whose } \\ \text{ith element} = \frac{\partial h_1(x, u)}{\partial u_k} \quad (7.38)$$

$\left(\frac{\partial f(x,u)}{\partial x_1} \right)$ is a column vector whose

$$\text{ith element} = \frac{\partial f(x,u)}{\partial x_i} \quad (7.39)$$

From (7.33) to (7.39) following equations result

$$\frac{\partial f(x,u)}{\partial u_k} = \frac{\partial f(x,u)}{\partial u_k} - \left(\frac{\partial f(x,u)}{\partial x_1} \right)^{-1} \left(\frac{\partial h_j(x,u)}{\partial x_j} \right)^{-1} \left(\frac{\partial h_j(x,u)}{\partial u_k} \right) \quad (7.40)$$

Defining the dual variables λ ,

$$(\lambda) = -\left(\frac{\partial h_j(x,u)}{\partial x_j} \right)^{-1} \left(\frac{\partial f(x,u)}{\partial x_1} \right) \quad (7.41)$$

Equation 7.40 is as follows

$$\begin{aligned} \frac{\partial f(x,u)}{\partial u_k} &= \frac{\partial f(x,u)}{\partial u_k} + \sum_j \lambda_j \frac{\partial h_j(x,u)}{\partial u_k} \\ &= \frac{\partial}{\partial u_k} (f(x,u) + \sum_j \lambda_j b_j(x,u)). \end{aligned} \quad (7.42)$$

Defining the Lagrangian

$$L(x,u) = f(x,u) + \sum_j \lambda_j b_j \quad (7.43)$$

where λ may be obtained from equation 7.41 or from equation 7.44 in which case the resulting expression will be similar

$$\left(\frac{\partial L(x,u)}{\partial x_j} \right) = 0 \quad (7.44)$$

Thus if the dual variables are obtained from equation

7.44, the partial derivative of Lagrangian with respect to u_k is in fact the partial derivative of the objective function with respect to u_k under the constraint domain of 7.32. After obtaining the partial derivatives in this way, any of the gradient techniques described in sections 7.5, 7.6 and 7.7 may be used.

Bryson and Denham⁴⁸, and Dommel and Tinney²¹ have used this technique.

Sufficiency condition for the minima may be stated as follows

$$\begin{bmatrix} \Delta x^t & \Delta u^t \end{bmatrix} \begin{bmatrix} \frac{\partial^2 L}{\partial x^2} & \frac{\partial^2 L}{\partial x \partial u} \\ \frac{\partial^2 L}{\partial u \partial x} & \frac{\partial^2 L}{\partial u^2} \end{bmatrix} \begin{bmatrix} \Delta x \\ \Delta u \end{bmatrix} \geq 0 \quad (7.45)$$

Under the constraint space of 7.32, or

$$(\Delta x) = -\left(\frac{\partial h_1(x, u)}{\partial x_j}\right)^{-1} \left(\frac{\partial h_1(x, u)}{\partial u_j}\right) (\Delta u) \quad (7.46)$$

Where $\left(\frac{\partial h_1(x, u)}{\partial u_j}\right)$ has its
ijth element = $\frac{\partial h_1(x, u)}{\partial u_j}$ (7.47)

7.10 Fuhn Tucker's theorem

For the nonlinear programming problem stated in 7.1, let

$$L = f(x) + \sum_i \alpha_i g_i(x) + \sum_i \lambda_i h_i(x) \quad (7.48)$$

The optimal solution satisfies following equations:

$$\frac{\partial L}{\partial x_i} = 0 \quad (7.49)$$

$$\alpha_i \geq 0 \quad (7.50)$$

$$\alpha_i g_i(x) = 0 \quad (7.51)$$

$$h(x) = 0 \quad (7.52)$$

$$g_i(x) \leq 0 \quad (7.53)$$

α_i and λ_i are the dual variables associated with the inequality constraints and equality constraints respectively. Equations 7.50 and 7.51 imply that $\alpha_i = 0$ if the constraint 7.53 is not reached and $\alpha_i > 0$ in the alternative.

7.11 Penalty and Barrier methods

These methods seek the solution of following nonlinear programming problem

$$\min f(x) \quad (7.54)$$

$$\text{subject to } g(x) \leq 0 \quad (7.55)$$

7.11.1 Penalty method

Defining a continuous function $P(x)$ termed as the penalty function, such that

$$P(x) = 0 \quad \text{if } x \in F \quad (7.56)$$

$$P(x) > 0 \quad \text{if } x \notin F \quad (7.57)$$

where F represents the feasible domain of 7.55.

One commonly used penalty function is as follows:

$$P(x) = \sum_{i=1}^m \max(0, g_i(x))^2 \quad (7.58)$$

Let

$$C(x, R) = f(x) + \frac{1}{R} P(x) \quad (7.59)$$

The penalty method uses a sequence R^k , $k = 1, \dots$ such that $R^k > 0$; $R^k > R^{k+1}$ and

$$\lim_{k \rightarrow \infty} R^k = 0$$

Defining x^k as an x that minimizes

$$C(x, R^k) = f(x) + \frac{1}{R^k} P(x) \quad (7.60)$$

The optimal point via the penalty method is obtained as follows. For a certain $R^1 > 0$ solve for x^1 using the unconstrained minimization technique. x^2 is obtained for another $R^2 < R^1$, and the process is repeated.

Following results can be easily obtained -

$$C(x^k, R^k) \leq C(x^{k+1}, R^{k+1})$$

$$P(x^k) \geq P(x^{k+1})$$

$$f(x^k) \leq f(x^{k+1})$$

$$x^* = \lim_{k \rightarrow \infty} x^k$$

7.11.2 Powell's penalty method¹⁶

With the penalty function approach mentioned earlier, there is a tendency to involve very large numbers as $R_i \rightarrow 0$, making the optimization difficult.

Powell's modification is as follows:

$$W(x, s) = 0 \quad \text{if } x \in F$$

$$W(x, s) = \sum_{i | g_i(x) > 0} \frac{(g_i(x) + s_i)^2}{R_i} \quad \text{if } x \notin F \quad (7.61)$$

$$C(x, s) = f(x) + W(x, s) \quad (7.62)$$

s_i is varied from iteration to iteration as follows

$$s_i^{k+1} = s_i^k + g_i^k(x^k) \quad (7.63)$$

$$\text{Thus } s_i^{k+1} > s_i^k$$

Following results can be easily obtained

$$C(x^k, s^k) \leq C(x^{k+1}, s^{k+1}) \quad (7.64)$$

$$g_i(x^k) \geq g_i(x^{k+1})$$

$$f(x^k) \leq f(x^{k+1})$$

$$\bar{x} = \lim_{k \rightarrow \infty} x^k$$

7.11.3 Barrier method

A function $B(x)$ is introduced such that it becomes infinity if $g(x) = 0$. This method needs that the initial guess be strictly in the constraint region which may be very difficult to obtain. The method forces the solution in the constraint region from iteration to iteration.

Piacco McCormic¹⁶ has suggested the following function

$$B(x) = - \sum_i \frac{1}{g_i(x)} \quad (7.65)$$

Lootma's¹⁶ function is as follows:

$$B(x) = - \sum_i \log(-g_i(x)) \quad (7.66)$$

The transformation is as follows:

$$D(x, k) = f(x) + R B(x) \quad (7.67)$$

The method uses a sequence R^k , $k = 1, \dots$ such that

$$R^k > C \text{ for all } k \text{ and}$$

$$R^k > R^{k+1} \text{ and}$$

$$\lim_{k \rightarrow \infty} R^k = 0$$

following results can be easily obtained:

$$D(x_1^k, \bar{x}^k) \geq D(x_1^{k+1}, \bar{x}^{k+1})$$

$$D(x^k) \leq D(x^{k+1})$$

$$f(x^k) \geq f(x^{k+1})$$

$$x^* = \lim_{k \rightarrow \infty} x^k$$

-x-

CHAPTER VIII

A REVIEW OF OPTIMAL POWER FLOW TECHNIQUES

This chapter critically reviews the published work on optimal power flow studies. The convergence of some of the techniques is compared through fixed point analysis.

6.1 Objective function

The cost f in rupees per unit time of operating the power system depends upon the active powers P_i^G , the reactive powers Q_i^G being free of cost once the equipment for their production e.g. capacitors etc. has been installed. Thus

$$f = \sum_{i \in G} f_i(P_i^G) \quad (6.1)$$

Optimization of equation 6.1 is subject to the constraints described later.

6.2 Classification

The techniques for finding the optimal point can be broadly classified as follows:

(1) Loss formula formulation

(2) Formulation with full a.c. simulation of network.

8.3 Loss formula method

In this method optimization of equation 8.1 is achieved through the following constraint:

$$P_L - \sum_I P_I = 0 \quad (8.2)$$

System loss P_L is expressed as a quadratic function of power generations. This method though most widely in use because of simplicity in application is based on the following assumptions, none of which really hold good in practice.

- (a) Bus voltages remain constant in magnitudes and angles.
- (b) Individual loads remain a constant complex fraction of the total load.
- (c) Ratios of reactive to real power generations remain constant.

This method optimizes for real powers only.

Because of the limitations mentioned above a need for a more accurate method was felt.

8.4 Popov, Elitin, Stagg and Watson's³² method (1967)

This method, essentially a modification of loss formula method, expresses losses as a quadratic function of active and reactive bus bar powers as follows:

$$P_L + \sum_j Q_j = I^T Z_{bus} I \quad (8.3)$$

$$\begin{aligned} P_L &= \sum_j \sum_k (P_j V_k \alpha_{jk} + Q_j V_k \beta_{jk}) \\ &\quad + P_j Q_k \beta_{jk} - Q_j P_k \beta_{jk} \end{aligned} \quad (8.4)$$

where,

$$\alpha_{jk} = \frac{r_{jk}}{V_j V_k} \cos \theta_{jk} \quad (8.5)$$

$$\beta_{jk} = \frac{-r_{jk}}{V_j V_k} \sin \theta_{jk} \quad (8.6)$$

$$\frac{\partial P_L}{\partial P_j} = 2 \sum_{k=1}^n (\alpha_{jk} P_k + \beta_{jk} Q_k) \quad (8.7)$$

$$\frac{\partial P_L}{\partial Q_j} = 2 \sum_{k=1}^n (\alpha_{jk} Q_k - \beta_{jk} P_k) \quad (8.8)$$

Augmenting the objective function of equation 8.1 with the constraint equation of 8.2 through dual variable the Lagrangian is as follows:

$$L = \sum_{i \in G} f_i(P_i) + \lambda (P_L - \sum_j P_j) \quad (8.9)$$

Resulting optimization equations are -

$$\frac{\partial f_i}{\partial P_i} + \lambda \frac{\partial P_L}{\partial P_i} = \lambda \quad \text{for } i \in G \quad (8.10)$$

$$\frac{\partial P_L}{\partial Q_i} = 0 \quad \text{for } i \in R \quad (8.11)$$

The suggested algorithm is as follows.

- (a) Estimate a generation schedule for real and reactive powers and obtain α_{jk} and β_{jk} through load flow solution.
- (b) With the active generations unchanged, obtain new reactive power schedule, making use of steepest descent method (section 7.5) with the gradients obtained from equation 8.6. This step needs repeated load flow solution and computation of α_{jk} and β_{jk} .
- (c) With the reactive power schedule computed above; obtain α_{jk} , β_{jk} and system losses.
- (d) Estimate λ and obtain the new power generations from equation 8.10 and the incremental Loss equation.
- (e) Check if equation 8.2 is satisfied, if not repeat step (d) with another λ .
- (f) Repeat from step (b) till solution converges.

8.4.1 Author's comments

The algorithm is based on equations 8.7 and 8.8. Evidently though not admitted in the paper these equations are obtained on the assumption that α_{jk} and β_{jk} remain unchanged for the incremental change in generation, which amounts to assuming the bus bar voltage remains constant in magnitude and angle.

The reactive power optimization is achieved through the steepest descent method, with the gradients obtained from equation 8.6. This does not appear reasonable since the slack bus reactive power is a dependent variable. If the incremental change in the reactive power losses is ignored;

$$\begin{aligned} \frac{\partial P_1}{\partial x_1} &= \frac{\partial P_1}{\partial V_1} - \frac{\partial P_1}{\partial \theta_1} \\ &= 2 \sum_{k=1}^n (\alpha_{1k} - \alpha_{sk}) c_k - (\beta_{1k} - \beta_{sk}) r_k \quad (8.12) \end{aligned}$$

Gradient method based on equation 8.12 should be more reasonable. A more accurate representation taking incremental change in the reactive power loss into account will however need additional coefficients ($x_{ij} \cos \theta_{ij}/V_i V_j$) and ($x_{ij} \sin \theta_{ij}/V_i V_j$).

Fixed point analysis

Fixed point analysis to test the convergence for real power optimization has been made as follows:

Assuming the incremental cost of generation a straight line;

$$\frac{\partial f_i}{\partial P_i^G} = a_1^1 + a_1^2 p_i^G \quad ; \quad i \in G \quad (8.13)$$

In the suggested algorithm, this is made equal to

$$\lambda \left(1 - \frac{\partial P_1}{\partial V_1} \right) \quad (8.14)$$

Fixed point representation is as follows:

$$\lambda + \frac{\alpha_{11}}{a_{11}^2} = \frac{\alpha_{11}}{a_{11}^2} - \frac{r_1}{a_{11}^2} \quad (6.15)$$

$$r_1 = \frac{\alpha_{11}}{a_{11}^2} - \frac{r_1}{a_{11}^2}$$

$$= r_1(P^G) \quad (6.16)$$

λ obtained from equations 6.7 and 6.15 is as follows:

$$\lambda = \frac{\sum_{k \in S} \alpha_{kk} + \sum_{k \in S} \frac{a_k}{a_k^2}}{\sum_{k \in S} \frac{(1 - \partial r_1 / \partial r_k)}{a_k^2}} \quad (6.17)$$

$$\frac{\partial \lambda}{\partial r_1} = \frac{2 \lambda \sum_{k \in S} \frac{\alpha_{kk}}{a_k^2} + \frac{\partial r_1}{\partial r_1}}{\sum_{k \in S} \frac{(1 - \partial r_1 / \partial r_k)}{a_k^2}} \quad (6.18)$$

$$\frac{-\partial r_1(P^G)}{\partial r_1} = \frac{1}{a_1^2} \left[-2 \alpha_{11} \lambda + \frac{\partial \lambda}{\partial r_1} (1 - \frac{\partial r_1}{\partial r_1}) \right] \quad (6.19)$$

The sufficiency condition for the Gaussian type iteration to converge (section 5.4) is

$$\sum_i \left| \frac{\partial f_i(P_i)}{\partial P_1} \right| < 1 \quad (8.20)$$

This condition will not be satisfied if any of the generation units has a near flat incremental cost characteristic. This establishes the unreliable convergence of the method.

A better method will be to obtain the real powers also through the gradient technique of section (7.9) with the following Lagrangian.

$$L = \sum_i f_i(P_i) + \lambda_1(i_1 - \sum_i p_i) + \lambda_2(i_1 - \sum_i \bar{p}_i) \quad (8.21)$$

Ignoring the incremental change of i_1 ; λ_1 and λ_2 are as follows:

$$\lambda_1 = \frac{\frac{\partial f_n}{\partial P_n} \lambda_n \bar{P}_n^G}{(1 - \frac{\partial P_1}{\partial P_n^G})} \quad (8.22)$$

$$\lambda_2 = \lambda_1 \frac{\partial P_1}{\partial \bar{P}_n} \quad (8.23)$$

$$\frac{\partial f}{\partial P_1^G} = \frac{\partial f_1}{\partial P_1^G} + \lambda_1 \left(\frac{\partial P_1}{\partial P_1^G} \right) = 1 \quad (8.24)$$

$$\frac{\partial f}{\partial P_1^G} = \lambda_1 \frac{\partial P_1}{\partial \bar{P}_n} + \lambda_2 \quad (8.25)$$

While the convergence with Gauss iterative method was found doubtful, a first order gradient method is stable since

This ensures the objective function reduce from step to step.

Additional drawback of the method is large storage requirement for non-sparse γ -matrix and α_{jk} , B_{jk} coefficients.

8.5 Formulation with full a.c. simulation of network^{19,36,39,46,47}

These methods optimize equation 8.1 under the equality constraints of equations 4.1 and 4.2 which are restated as follows:

$$P_i = P_i^+ - P_i^- = p_i(V, \theta) = V_i \sum_j V_j Y_{ij} \cos(\theta_i - \theta_j - \Psi_{ij}) \quad (8.26-i)$$

$$Q_i = Q_i^G - Q_i^L = q_i(V, \theta) = V_i \sum_j V_j Y_{ij} \sin(\theta_i - \theta_j - \Psi_{ij}) \quad (8.27-i)$$

Inequality constraints on generation nodes are as follows:

$$(P_i^G)^2 + (Q_i^G)^2 - (\bar{P}_i)^2 \leq 0 \quad (8.28)$$

$$\underline{P}_i - \bar{P}_i \leq 0 \quad (8.29)$$

$$P_i^G - \bar{P}_i \leq 0 \quad (8.30)$$

$$Q_i^G - \bar{P}_i \leq 0 \quad (8.31)$$

$$\underline{Q}_i - \bar{Q}_i \leq 0 \quad (8.32)$$

Constraints 8.28 to 8.32 are determined by current magnitude, boiler instability, turbine output, generator excitation limits and generator instability respectively.

Additional constraints determined primarily by insulation and transformer tap ranges for consumption as well as generation nodes are as follows:

$$v_i - \bar{v}_i \leq 0 \quad (8.33)$$

$$\underline{v}_i - v_i \leq 0 \quad (8.34)$$

In addition to these, upper limit of the line currents also need be fixed.

$$I_{ij} = \left| \frac{v_i / e_i - v_j / e_j}{z_{ij}} \right|$$

where I_{ij} is the current and z_{ij} the impedance of line connected between nodes i and j.

$$= \left| \frac{\frac{e_i - e_j}{2} / \frac{\pi - e_i - e_j}{2} + (v_i - v_j) / e_j}{z_{ij}} \right|$$

$$\approx \left| \frac{2 v_i \sin(\frac{e_i - e_j}{2})}{z_{ij}} \right|$$

for normal differences in line voltages and phase angles.

Thus prescribing the upper limit of line current almost amounts to prescribing the upper limit of $|e_i - e_j|$. This constraint can be described as follows:

$$|e_i - e_\alpha| - T_{i\alpha} = 0 \quad (8.35)$$

where α is the adjacent node.

8.6 Subclassification

Methods making use of a.c. representation of network can be further classified as follows:

- (1) Variational equations and linear programming formulation²⁸
- (2) Kuhn Tuckers formulation^{11,13,19,45}
- (3) Penalty function formulation^{4,15,16}
- (4) Mixed formulation, making use of both Kuhn Tucker's theorem or equivalent formulation and penalty functions^{17,21}

8.7 Lauppin, Feingold and Pohn's method²⁶ (1967)

This method makes use of variational equations and linear programming technique, for which the objective function is as follows:

$$f = \sum_{i \in G} (a_i + b_i P_i^0) \quad (8.36)$$

$$\Delta f = \sum_{i \in G} b_i \Delta P_i^0 \quad (8.37)$$

The general compatibility relation of equation 4.32

is as follows:

$$\sum_{i=1}^n \lambda_i \Delta P_i + \sum_{k \in R} \mu_k \Delta V_k + \sum_{j=1}^m u_j \Delta V_j = 0 \quad (8.33)$$

λ, μ, \dots, k and l are as defined in section 4.6.

The method seeks the minimization of equation 8.37 under constraint of equations 8.36 and 8.29 to 8.34.

Since the effect of second order terms is ignored, this in general causes ΔP_i^G , ΔQ_i^G and ΔV_i to fail to satisfy the power flow equations. Therefore the distribution is recalculated after obtaining the new coefficients of compatibility equations. Possibility of new values of the objective function greater than the previous value has also been admitted, in which case recalculations on the basis of half the correction has been recommended.

8.7.1 Author's comments

From equations 8.37 and 8.38,

$$\begin{aligned} \Delta f = & \sum_{\substack{i \in G \\ i \neq s}} \left(b_i - \frac{b_s \lambda_i}{\lambda_s} \right) \Delta P_i - \sum_{k \in R} \frac{\mu_k b_k}{\lambda_s} \Delta V_k \\ & - \sum_{j \in R} \frac{u_j b_j}{\lambda_s} \Delta V_j \end{aligned} \quad (8.39)$$

Assuming the inequality constraints of slack bus not violated for minimum Δf ; following equations hold good:

$$\Delta P_1 = \underline{P}_1^G - \bar{P}_1^G \quad \text{if} \quad b_1 - \frac{b_e \lambda_1}{\lambda_s} < 0$$

$$\Delta P_1 = \underline{P}_1^G - \bar{P}_1^G \quad \text{if} \quad b_1 - \frac{b_e \lambda_1}{\lambda_s} > 0$$

$$\Delta \gamma_k = \bar{\gamma}_1 - \underline{\gamma}_1 \quad \text{if} \quad \frac{b_e u_k}{\lambda_s} > 0$$

$$\Delta \gamma_k = \underline{\gamma}_1 - \bar{\gamma}_1 \quad \text{if} \quad \frac{b_e \bar{\gamma}_k}{\lambda_s} < 0$$

$$\Delta V_1 = \bar{V}_1 - V_1 \quad \text{if} \quad \frac{u_1 b_e}{\lambda_s} > 0$$

$$\Delta V_1 = V_1 - \underline{V}_1 \quad \text{if} \quad \frac{u_1 b_e}{\lambda_s} < 0$$

This amounts to substituting for either maximum value or minimum value of independent variables for most buses in the next iteration. This indicates that jamming of the algorithm is quite likely to occur.

8.8 Carpentier's method^{36,37,46,47} (1962-1963)

This method is based upon the following Lagrangian:

$$\begin{aligned} L = & \sum_{i \in G} r_i(\underline{P}_i^G) + \sum_i \lambda_i(P_i(V, e) - \bar{P}_i) \\ & + \sum_i \mu_i(Q_i(V, e) - \bar{Q}_i) + \sum_{i \in B \cup G} r_i((\underline{P}_i^G)^2 \\ & + (\bar{Q}_i^G)^2 - \varepsilon_i^2) + \sum_{i \in G} m_i(\underline{V}_i - \bar{V}_i) \end{aligned}$$

$$\begin{aligned}
 & \cdot \sum_{i \in A} e_i (\bar{e}_i - \bar{e}_1) + \sum_{i \in B} e'_i (\bar{e}_1 - \bar{e}'_i) \\
 & + \sum_i \sum_{\alpha} r_{i\alpha} (e_i - e'_{\alpha} - T_{i\alpha}) \\
 & + \sum_i \pi_i (v_i - \bar{v}_i) + \sum_i \pi'_i (v_i - v'_i) \quad (8.40)
 \end{aligned}$$

Necessary conditions for f to be minimum are given in section (7.10).

$$\frac{\partial L}{\partial p_i^G} = 0 \quad \text{gives}$$

$$\lambda_i = \frac{\partial f_i}{\partial p_i^G} + 2 K_i p_i^G - \pi_i \quad \text{for } i \in G. \quad (8.41-i)$$

$$\frac{\partial L}{\partial e_i^G} = 0 \quad \text{gives}$$

$$\mu_i = 2 h_i e_i^G + e_i - e'_i \quad \text{for } i \in R \quad (8.42-i)$$

$$\frac{\partial L}{\partial \theta_i} = 0 \quad \text{gives}$$

$$\sum_j \lambda_j \frac{\partial p_j}{\partial \theta_i} + \sum_j \mu_j \frac{\partial q_j}{\partial \theta_i} + \sum_{\alpha} (r_{i\alpha} - r'_{i\alpha}) = 0 \quad (8.43-i)$$

$$\frac{\partial L}{\partial v_i} = 0 \quad \text{gives}$$

$$\sum_j \lambda_j \frac{\partial p_j}{\partial v_i} + \sum_j \mu_j \frac{\partial q_j}{\partial v_i} + \pi_i - \pi'_i = 0 \quad (8.44-i)$$

Since slack bus voltage and angle is fixed; equations 8.43 and 8.44 do not apply for this bus.

Apart from these equations; power flow equations of 8.26 and 8.27 must be satisfied.

8.8.1 Verification of number of equations and unknowns

(i) Consumption nodes: The unknowns are $\theta_i, v_i, \lambda_i, \mu_i$ and the non-zero dual variables corresponding to the inequality constraints of equations 8.26 to 8.35.

The equations are 8.26, 8.27, 8.43 and 8.44 plus one equation each arising from the active inequality constraints.

(ii) Production nodes (excluding slack node): The unknowns are $P_i^G, Q_i^G, \epsilon_i, v_i, \lambda_i, \mu_i$ and the non-zero dual variables of inequality constraints. The equations are 8.41, 8.42, 8.43, 8.44, 8.26 and 8.27 plus one equation each arising from the corresponding active inequality constraints.

(iii) slack node: The unknowns are P_i^G, Q_i^G, λ_i and μ_i and the non-zero dual variables of the inequality constraints. The equations are 8.41, 8.42, 8.26 and 8.27 plus one equation each from the active inequality constraint.

8.8.2 Algorithm

(i) Assume a set of unknown P, Q, V, θ, μ and λ . These may not satisfy the power flow equations.

- (2) Assuming $\Delta \theta_{\alpha} = 0$ and $\Delta V_{\alpha} = 0$; obtain Δe_i and ΔV_i from the variation equations of 8.26-1 and 8.27-1 respectively as follows:

$$\Delta e_i = \frac{P_i^G - P_i^L - p_i(V, \theta)}{\partial p_i(V, \theta) / \partial e_i} \quad (8.45-1)$$

$$\Delta V_i = \frac{V_i^G - V_i^L - q_i(V, \theta)}{\partial q_i(V, \theta) / \partial V_i} \quad (8.46-1)$$

Check if $e_i + \Delta e_i$ and $V_i + \Delta V_i$ does not violate the inequality constraints. If these constraints are violated, fix e_i and V_i to the constraint value.

- (3) If constraints in (2) are not active, $r_{i\alpha}$, $r_{\alpha i}$, π_i and π'_i are equal to zero, obtain λ_i and μ_i respectively from equations 8.43-1 and 8.44-1, if i is not a slack node.
- (4) If constraints in (2) are not active and it is not a slack node, obtain P_i^G and Q_i^G from equations 8.41-1 and 8.42-1 respectively, if it is a production node. If these powers are within limits, b_i , m_i , e_i and e'_i are zero. If power constraints are violated, these are equated to the constraint value and corresponding dual variables obtained from equations 8.41-1 and 8.42-1 respectively.

(5) If constraint in (2) are active and this is a production node, generations are obtained from equations 8.26-i and 8.27-i. λ_i and μ_i are obtained from equations 8.41-i and 8.42-i respectively. If it is not a production node, λ_i or $r_{i\alpha}$ and λ_j or $r_{j\beta}$ are obtained from equations 8.43-i and 8.44-i respectively.

The iteration is repeated from (2) till the solution converges.

8.3.3 Author's comments

The method turns out to be the Gauss Seidel iterative method applied to each node in succession. The convergence of this iterative technique is not reliable. A fixed point analysis of the method also appears difficult. The method assumes that a solution in the feasible domain satisfying all equality and inequality constraints exist which may not always be so.

The solution does not remain in the feasible domain from iteration to iteration. Hence final values are of little use if solution fails to converge. Investigations on this method by Pernchon¹⁹ et al show divergence in solutions where the initial guess is not close enough to the optimal solution. For systems with unusually high or low reactive powers, the method failed to give the solution even if the initial guess is very close to the optimal solution.

This paper can be regarded as a historical break through. Later on number of other papers making use of Carpentier's formulation have appeared. Numerical technique adopted to solve system of equations are based on either Gauss Seidal or gradient method. In many cases not all the inequality constraints are considered.

8.9 El-Abiad and Jaimes' method¹³ (1969)

In this paper inequality constraints are excluded. Partial derivative of the Lagrangian with respect to the slack bus voltage is also equated to zero, which evidently means that this voltage is also adjustable. Equations 8.41 to 8.44 thus reduce to the following:

$$\lambda_i = \frac{\partial f_i}{\partial P_i} , \quad i \in G \quad (8.47-1)$$

$$\mu_i = 0 ; \quad i \in R \quad (8.48-1)$$

$$\sum_j \lambda_j \frac{\partial P_j}{\partial \theta_i} + \sum_j \mu_j \frac{\partial Q_j}{\partial \theta_i} = 0 ; \quad i \neq s \quad (8.49-1)$$

$$\sum_j \lambda_j \frac{\partial P_j}{\partial V_i} + \sum_j \mu_j \frac{\partial Q_j}{\partial V_i} = 0 \quad (8.50-1)$$

From equations 8.48, 8.49 and equations of 8.50 for nodes with no reactive power generation resulting equations is 8.51.

$$\begin{array}{|c|c|c|c|c|} \hline
 \text{a}_1^1 = \frac{\partial P_j}{\partial e_1} & \text{a}_{1j}^2 = \frac{\partial Q_j}{\partial e_1} & \lambda_j & & \\ \hline
 \text{i.e. } j \in S & i \neq 1; j \in R & j \neq 0 & -\lambda_j & \\ \hline
 \text{a}_{1j}^3 = \frac{\partial P_j}{\partial V_1} & \text{a}_{1j}^4 = \frac{\partial Q_j}{\partial V_1} & \mu_j & & \\ \hline
 \text{i.e. } j \in R & i \in R, j \notin R & j \neq 0 & j \neq 0 & \\ \hline
 \end{array}$$

(8.51)

Jacobian on the right hand side of equation 8.51 is the transfer of jacobian needed for load flow solution by Newton's method with voltage as independent variable for reactive power production nodes.

The algorithm is as follows:

- (1) Assuming a set of feasible generations and voltages and a value of λ_1 ; λ and μ are obtained from equations 8.51 and 8.48. Real generations are obtained from equation 8.47 in terms of a multiplier λ' included to assure a proper level of generation. For the incremental costs given by equation 8.13 the value of λ' such that generations are equal to the pre-estimated losses and demand is given by 8.52

$$\lambda' = \frac{P_L + P_D + \sum_{i \in G} a_i^1/a_i^2}{\sum_{i \in G} \left\{ \frac{\lambda_1}{a_i^2} \right\}} \quad (8.52)$$

$$\text{and } \frac{\partial f_1}{\partial P_1} = \lambda' \lambda_1 \quad (8.53)$$

- (2) Obtain voltage of the nodes with reactive power generations from the equations 8.50-i, $i \neq 1$, which after rearranging is as follows:

$$AB = b \quad (8.54)$$

where A is a square matrix whose ij th element is as follows:

$$a_{ij} = \lambda_j v_{ij} \cos(\theta_i - \theta_j - \psi_{ij}) \\ + \mu_i v_{ij} \sin(\theta_i - \theta_j - \psi_{ij}) \text{ for } i \neq j; i \in R, \\ j \in R$$

$$a_{ii} = \lambda_i E_{ii}$$

$$b_i = - \sum_{j \notin R} \lambda_j v_j v_{ij} \cos(\theta_i - \theta_j - \psi_{ij}) \\ - \sum_{j \notin R} \mu_j v_j v_{ij} \sin(\theta_i - \theta_j - \psi_{ij}) \text{ for } i \in R$$

v is a column vector of voltages of nodes with reactive power generation.

8.9.1 Author's comments

Though the exact fixed point analysis appears difficult, an analysis for real power dispatch has been made as follows.

Premultiplying both sides of equation 8.51 by a row vector $(\Delta \theta_i, i \neq s; \Delta V_i, i \neq R)$; following relation is

obtained

$$\sum_i \lambda_i \alpha_i + \sum_{i \notin R} \mu_i \alpha_i = 0 \quad (8.55)$$

Thus the dual variables obtained in this method are in fact coefficients of the compatibility relation of equation 4.30. These coefficients are related to the loss formula expression of equation 8.4 as follows:

$$\lambda_1 = \lambda' \left(1 - \frac{\partial P_1}{\partial \alpha_1} \right) \quad (8.56)$$

$$\mu_1 = -\lambda' \frac{\partial P_1}{\partial \alpha_1} \quad (8.57)$$

Step 1 of the algorithm thus amounts to

$$\frac{\partial f_1}{\partial \alpha_1} = \lambda_1 = \lambda' \left(1 - \frac{\partial P_1}{\partial \alpha_1} \right) \quad (8.58)$$

This step is similar to that of equation 8.13 and 8.14. Hence the fixed point analysis of 8.4.1 is valid for this method too. Thus convergence of the method is doubtful if any of the generation unit has a near flat incremental cost characteristics. Poor convergence and need of acceleration technique has been admitted for the five node problem given in the paper.

8.10 Shen and Laughton's method¹¹ (1969)

This method also utilizes the Carpentier's formulation and uses Gauss Seidel type iterative technique. Optimal

adjustment for phase shifting transformer and a tap setting is also included. Apart from the constant load demand, a variable component of load demand proportionate to the voltage magnitude has been included, however objective function is cost of real power generation. Excluding some of these additional features, the system Lagrangian is as follows:

$$L_1 = L + \sum_h C_h(t_h - \bar{t}_h) + \sum_h C'_h(\underline{t}_h - t_h) \\ - \sum_h B_h(\phi_h - \bar{\phi}_h) + \sum_h B'_h(\underline{\phi}_h - \phi_h) \quad (8.59)$$

where L is given by equation 8.4C.

Equations obtained are 8.41, 8.42, 8.43 and 8.44. Equations 8.43 and 8.44 are initially written for all the nodes including the slack node. Additional equations for the transformers are as follows:

$$\frac{\partial L_1}{\partial t_h} = \lambda_i \frac{\partial P_1}{\partial t_h} + \lambda_j \frac{\partial P_1}{\partial t_h} + \mu_i \frac{\partial Q_1}{\partial t_h} \\ + \mu_j \frac{\partial Q_j}{\partial t_h} + C_h - C'_h = 0 \quad (8.60)$$

$$\frac{\partial L_1}{\partial \phi_h} = \lambda_i \frac{\partial P_1}{\partial \phi_h} + \lambda_j \frac{\partial P_1}{\partial \phi_h} + \mu_i \frac{\partial Q_1}{\partial \phi_h} \\ + \mu_j \frac{\partial Q_j}{\partial \phi_h} + B_h - B'_h = 0 \quad (8.61)$$

where i and j are two nodes to which the k th transformer is connected.

In such iteration the computation is divided in four steps involving in turn computation of

(1) active generation and voltage phase angles

(2) phase shifting transformer tap settings

(3) reactive generation and voltage magnitudes

(4) non-phase shifting transformer tap settings

8.10.1 Determination of active generation and voltage phase angles

In this step, an estimate of power generations, voltage phase angles and dual variables λ is made, assuming other variables as fixed.

Let

$$e_{pi} = p_i(V, \theta) - P_i^G + P_i^L \quad (8.62)$$

$$e_{\theta i} = \frac{\partial \lambda}{\partial \theta_i} = \text{l.h.s. of equation 8.43-i} \quad (8.63)$$

$$\Delta e_{pi} + \Delta P_i^G = \sum_j \frac{\partial p_i(V, \theta)}{\partial \theta_j} \Delta \theta_j \quad (8.64)$$

$$\Delta^e \theta_1 = \sum_j \frac{\partial P_j}{\partial \theta_1} \Delta \lambda_j + \sum_j \frac{1}{\partial \theta_1} \left(\frac{\partial L_j}{\partial \theta_1} \right) \Delta e_j \quad (8.65)$$

Without any explanation, the paper indicates that $\frac{\partial}{\partial \theta_1} \left(\frac{\partial L_j}{\partial \theta_1} \right)$ elements are small; therefore equation 8.65 modified is as follows:

$$\Delta^e \theta_1 = \sum_j \frac{\partial P_j}{\partial \theta_1} \Delta \lambda_j \quad (8.66)$$

The method is as follows:

- (1) Since the matrix associated with 8.66 is singular; $\Delta \lambda_j$ are obtained from this equation in terms of $\Delta \lambda_1$, which is to be adjusted for proper level of generations obtained from equation 8.41.
- (2) Phase angles are obtained from 8.64.

8.10.2 Author's comments

This step is similar to the real power optimization of 8.9 in the following respect. Both obtain the dual variables as functions of λ_s , which is chosen for proper level of generations. Thus fixed point analysis of 8.9.1 is valid for this method too for real power optimization.

This indicates that convergence is doubtful if any of the generation unit has a near flat incremental cost characteri-

tice. λ and ϵ in this method have been obtained by assuming μ and ϵ respectively as constants. Convergence of a partitioned iterative method has been found poorer than convergence without partitioning (well established for Z-matrix load flow).

8.10.3 Obtaining phase shifting transformer settings

Equation 8.61 written explicitly is as follows:

$$\begin{aligned} \frac{\partial I_1}{\partial \beta_h} = & -\lambda_i v_i v_j y_{ij} \sin(\theta_i - \theta_j - \psi_{ij} + \beta_h) \\ & + \lambda_j v_i v_j y_{ij} \sin(\theta_j - \theta_i - \psi_{ij} - \beta_h) \\ & + \mu_i v_i v_j y_{ij} \cos(\theta_i - \theta_j - \psi_{ij} + \beta_h) \\ & - \mu_j v_i v_j y_{ij} \cos(\theta_j - \theta_i - \psi_{ij} - \beta_h) \\ & + \beta_h - \beta'_h \end{aligned} \quad (8.67)$$

Thus β_h is at a limit $\bar{\beta}_h$ or $\underline{\beta}_h$ with β_h or β'_h

positive or

$$(\theta_i - \theta_j + \beta_h) = \tan^{-1} \frac{(\lambda_i - \lambda_j) \sin \psi_{ij} + (\mu_i - \mu_j) \cos \psi_{ij}}{(\lambda_i + \lambda_j) \cos \psi_{ij} - (\mu_i + \mu_j) \sin \psi_{ij}} \quad (8.68)$$

New values of voltage phase angles and phase shifts are obtained from equation 8.68 and following variational equation obtained from equation 8.68 and following variational equation

$$\Delta P_1 = 0 = \sum_{k \neq i} \frac{\partial P_1}{\partial \beta_k} \wedge e_k + \sum_h \frac{\partial P_1}{\partial \beta_h} \Delta \beta_h \quad (8.69)$$

8.10.4 Obtaining reactive power generation and voltage magnitudes

Let

$$e_{vi} = q_i(v_i, t) = \frac{G}{V_i} + \frac{\mu}{V_i} \quad (8.70)$$

$$e_{vi} = \frac{\partial}{\partial V_i} = \text{left hand side of equation 8.44} \quad (8.71)$$

$$\Delta e_{qi} + \Delta \frac{\mu}{V_i} = \sum_j \frac{\partial q_i}{\partial V_j} \Delta V_j \quad (8.72)$$

$$\Delta e_{vi} = \sum_j \frac{\partial G_j}{\partial V_i} \Delta \mu_j + \sum_j \frac{\partial}{\partial V_j} (e_{vi}, \Delta V_j + \Delta \pi_i - \Delta \pi'_i) \quad (8.73)$$

Neglecting $\frac{\partial}{\partial V_j} (e_{vi})$ for $i \neq j$, $\Delta \frac{\mu}{V_i}$, $\Delta \mu$ and

ΔV are obtained from equations 8.72, 8.73 and 8.42, remembering

$$\Delta \frac{G}{V_j} = 0 \quad \text{for load buses}$$

$\mu_j = \Delta \mu_j = 0$ for buses with reactive power generation not violating V or Q constraints.

8.10.5 Obtaining optimal transformer settings

The transformer setting t_h is either at \bar{t}_h or \underline{t}_h with σ_h or σ'_h positive or following relation obtained from equation 8.60 hold good

$$G_{h1} t_h - \sigma_{h2} = 0 \quad (8.74)$$

where $G_{h1} = 2\lambda_1 V_1^2 y_{h1} \cos \psi_{h1}$

$$- 2 \mu_1 y_{h1} V_1^2 \sin \psi_{h1} \quad (8.75)$$

$$\begin{aligned}
 G_{h2} &= \lambda_i V_i V_j Y_{h2} \cos(\theta_i - \theta_j - \psi_{h2}) \\
 &+ \lambda_j V_i V_j Y_{h2} \cos(\theta_j - \theta_i - \psi_{h2}), \\
 &+ \mu_i V_i V_j Y_{h2} \sin(\theta_i - \theta_j - \psi_{h2}) \\
 &+ \mu_j V_i V_j Y_{h2} \sin(\theta_i - \theta_j - \psi_{h2}) \quad (8.76)
 \end{aligned}$$

$$y_{11} / \psi_{11} = y_{11}^o / \psi_{11}^o + t_h^2 y_{h1} / \psi_{h1} \quad (8.77)$$

$$y_{1j} / \psi_{1j} = y_{1j}^o / \psi_{1j}^o - t_h y_{h2} / \psi_{h2} \quad (8.78)$$

If σ_t is the value of left hand side of equation 8.74;

$$\begin{aligned}
 \Delta \sigma_t &= t_h \frac{\partial G_{h1}}{\partial V_1} \Delta V_1 + G_{h1} \Delta t_h \\
 &- \frac{\partial G_{h2}}{\partial V_1} \Delta V_1 - \frac{\partial G_{h2}}{\partial V_j} \Delta V_j \quad (8.79)
 \end{aligned}$$

$$\Delta q_1 = \frac{\partial q_1}{\partial V_1} \Delta V_1 + \frac{\partial q_1}{\partial V_j} \Delta V_j + \frac{\partial q_1}{\partial t_h} \Delta t_h \quad (8.80)$$

$$\Delta q_j = \frac{\partial q_1}{\partial V_1} \Delta V_1 + \frac{\partial q_1}{\partial V_j} \Delta V_j + \frac{\partial q_1}{\partial t_h} \Delta t_h \quad (8.81)$$

Δt_h , ΔV_1 and ΔV_j are obtained from equations 8.79, 8.80 and 8.81 with $\Delta q_1 = \Delta q_j = 0$ if voltage constraints are not violated.

8.10.6 Author's comments

The author has very extensively tested the method of tap adjustment described in 8.10.5 on IEEE 14 bus and 30 bus system and finds that this does not work. In most cases Δt_h is found of the same sign as the gradient obtained from equation 8.60. This indicates that the objective function will in fact increase if this adjustment is made, and optimal solution will not be obtained even with damping or acceleration.

Similar to Carpentier's method, the method also turns out to be the Gauss-Seidel iterative method, with the iteration not remaining in feasible domain. The method also assumes that a feasible solution exists. Certain cases where the solution fails to converge for real power scheduling and transformer settings have been pointed out. For active and reactive power scheduling, the method turns out a partitioned version of 8.9 in which case the authors themselves have pointed out poor convergence.

As pointed out earlier, the load includes a variable component proportionate to the voltage magnitude while the objective function is cost of real power generation. To the author this does not sound very reasonable. With change in voltage, revenue also changes because of this component of demand. Hence a more reasonable objective function is cost of generation minus revenue for the variable component of load demand.

• 11 Panichon, Pierrey, Tinney, Tveit and Cuenod's method 19,36

1962.

This method is in fact a historical development of the research carried out for the BPA system. The system has 1005 hydro-generation. Thus objective function is system loss. Method provides for reactive power optimization. Use has been made of the Lagrangian of equation 8.40 with objective function as slack bus power and constraints on active power, apparent power and angles ignored. Resulting equations are as follows:

$$\lambda_p = 1 \quad (8.82)$$

$$\mu_i = e_i - e'_i \quad (8.83)$$

$$\sum_j \lambda_j \frac{\partial p_i}{\partial e_j} + \sum_j \mu_j \frac{\partial q_i}{\partial e_j} = 0; i \neq n \quad (8.84)$$

$$\sum_j \lambda_j \frac{\partial p_i}{\partial v_j} + \sum_j \mu_j \frac{\partial q_i}{\partial v_j} \cdot \bar{v}'_i - \bar{v}'_i = 0 \quad (8.85)$$

In view of the obvious fact that efficient transmission of power implies high voltage levels throughout the grid, it is initially guessed that at the optimal point nodes with reactive power generation, except where the upper limit of generation is violated, will be at the maximum voltage. The algorithm is as follows:

- (1) With $V = \bar{V}$ at the reactive production nodes, all remaining primary variables P , Q , E and V are obtained from power flow equations.

- (2) For nodes where $\beta > \bar{\beta}$; $\beta = \bar{\beta}$ is imposed and primary variables recomputed.
- (3) Obtain dual variables from equations 8.82 to 8.85 and check for the values. If $e_i = \bar{e}_i > 0$. If $V_i = \bar{V}_i$; $\pi_i > 0$. For nodes where e_i or π_i are negative; corresponding C_i or V_i is decreased to make these dual variables as zero. A linear relationship is assumed between V_i and π_i or C_i and e_i as follows:

$$V_i(k) = V_i(k-1) - \frac{V_i(k-1) - V_i(k-2)}{\pi_i(k-1) - \pi_i(k-2)} \pi_i(k-1) \quad (8.86)$$

$$e_i(k) = e_i(k-1) - \frac{C_i(k-1) - C_i(k-2)}{e_i(k-1) - e_i(k-2)} e_i(k-1) \quad (8.87)$$

8.11.1 Author's comments

This method turns out to be a first order gradient method with a set of dependent variables adjusted through linear extra-polation. Convergence depends essentially upon the choice of $V_1(2) - V_1(1)$ and $C_1(2) - C_1(1)$. The iteration is in the feasible domain of power flow equations. It is admitted that convergence has been good where there are few reactive production nodes with the optimal voltage not at the maximum, which essentially means that the initial guess is sufficiently close to the optimal solution.

Two separate load flow solutions have been suggested in

step (1) and (2). With the authors load flow programme described in Chapter 6, this can be achieved through a single solution.

d.12 Fasson's method^{4,15,16}

This method¹⁶ makes use of Powell's penalty function formulation of section (7.11.2) to represent power flow equations and other inequality constraints, all expressed in rectangular form. Optimization is achieved by Fletcher Powell's gradient method⁴⁴ (sec. 7.7) in terms of real and quadrature components of line voltages, and active and reactive power generations. The approach needs less computer logic compared to Carpentier-type methods. Because of very large number of variables, optimization may be very difficult to achieve in presence of discontinuous penalty functions. Fletcher Powell's gradient method need a non-sparse though symmetrical Hessian matrix. For the 30 node problem 230 Fletcher Powell's steps were needed¹⁶. With 4 nodes for real power generation and 6 nodes for reactive power generation, a 68×68 matrix need be recalculated for every Fletcher Powell's step. Iteration is not in the feasible domain.

In a later paper⁴, it is admitted that the method becomes very inefficient in terms of storage requirement and time for large systems. For such systems use of decomposition technique has been suggested. The paper demonstrates that this way the computer time is cut down to 10% for a system of 57 buses and

5. For 11C buses, efficiency of decomposition depends upon how the system is partitioned, which still is an open question.

8.13 Bonnel and Tinney's method²¹

This method makes use of mixed formulation. The power flow equations in polar form have been handled through the dual variables and other inequality constraints by augmenting the objective function through penalty function approach, of 7.11.1. The independent variables (called control variables) are the active power generations and line voltage magnitude or reactive power of the generation nodes except the powers for the slack node. Slack bus active power generation is expressed as a function of voltage magnitudes and angles. Thus the objective function of equation 8.1 is rewritten as follows:

$$f = \sum_{\substack{i \in G \\ i \neq s}} f_i(p_i^G) + f_s(p_s(V, \theta) + p_s^L) \quad (8.88)$$

Sensitivity of the objective function with respect to the decision variables is obtained by the method of section (7.9) making use of the following Lagrangian:

$$\begin{aligned} L = & \sum_{\substack{i \in G \\ i \neq s}} f_i(p_i^G) + f_s(p_s^*(V, \theta) + p_s^L) + \text{penalty terms} \\ & + \sum_{\substack{i \\ i \neq s}} \lambda_i(p_i(V, \theta) - p_i^G + p_i^L) \\ & + \sum_{\substack{i \\ i \neq s}} \mu_i(q_i(V, \theta) - z_i^G + z_i^L) \end{aligned} \quad (8.89)$$

It may be noted that this Lagrangian is different from that of Carpentier's. Use has been made of the steepest descent technique of section (7.5) with the optimal step length obtained through parabolic extrapolation (section 7.8-1). Whenever the gradient changes the sign from iteration to iteration, the control variable is re-obtained through linear interpolation. This modification of the gradient method has been called a mixed method.

C.13.1 Author's comments

This technique though not the latest the author finds as the most promising of all the discussed techniques for the following reasons:

- (1) The iteration is in the feasible domain of power flow equations.
- (2) Gradient technique has been used instead of Gauss Seidel iterative technique. While convergence of the later technique is doubtful, convergence of the former is more reliable since ensures that the objective function reduce from step to step, even if the initial guess is not in the convex domain.
- (3) Dependent variables (called state variables) have been obtained through load flow solution and not through gradient technique. Convergence of the gradient technique becomes more difficult as the number of

variables to be adjusted by this technique increase. Application of Newton's load flow assures that the power flow equations are satisfied to a very high degree of accuracy.

- (4) Algorithm makes the application of sparsity techniques possible.

The author's experience over the method while trying

PP 14 bus and 30 bus systems is as follows:

- (1) Application of parabolic interpolation in the presence of penalty functions did not work satisfactorily.
- (2) Though the paper suggests the use of either voltage magnitude or reactive power for nodes with such generations as independent variable, the author finds that the step length obtained with voltage magnitude as independent variable is much smaller (about 50%) compared to one obtained with reactive power as such variable. This makes the change in tap settings and real power generation smaller, resulting in the need for more iterations with the former as independent variable.
- (3) Application of mixed gradient method did not work satisfactorily in the presence of adjustable reactive power and tap settings possibly because the off diagonal elements of the Hessian $\frac{\partial^2 f}{\partial \Delta t^2}$ could not be ignored.

S.14 Zamangorthry and Gopala Rao's method¹⁷

In this paper voltage magnitudes and angles of generation buses are adjusted through gradient method. Inequality constraints on other variables are represented by square type penalty functions. In the analysis, application of dual variables has been avoided. It is as follows:

$$\begin{bmatrix} \Delta P^G \\ \Delta Q^G \\ \Delta P^L \\ \Delta Q^L \end{bmatrix} = \begin{bmatrix} c_1 & & & c_2 \\ & \text{---} & & \\ & & c_3 & c_4 \end{bmatrix} \begin{bmatrix} \Delta \theta^G \\ \Delta V^G \\ \Delta \theta^L \\ \Delta V^L \end{bmatrix} \quad (8.90)$$

With $\Delta P^L = 0$ and $\Delta Q^L = 0$,

$$\begin{bmatrix} \Delta \theta^L \\ \Delta V^L \end{bmatrix} = -c_4^{-1} c_3 \begin{bmatrix} \Delta \theta^G \\ \Delta V^G \end{bmatrix} \quad (8.91)$$

$$\begin{bmatrix} \Delta P^G \\ \Delta Q^G \end{bmatrix} = \left[c_1 - c_2 c_4^{-1} c_3 \right] \begin{bmatrix} \Delta \theta^G \\ \Delta V^G \end{bmatrix} \quad (8.92)$$

Elements of jacobian in equation 8.92 are the partial derivatives of the generation with respect to the voltage phase angles and magnitudes of generation buses in the domain of power flow equations.

$$\frac{\partial f}{\partial \theta_i^G} = \sum_j \frac{\partial f_j}{\partial \theta_j^G} \frac{\partial P^G}{\partial \theta_i^G} + \sum_j \frac{\partial f_j}{\partial V_j^G} \frac{\partial Q^G}{\partial \theta_i^G} \quad (8.93)$$

$$\frac{\Delta f_i}{f_i} = \frac{1}{2} \frac{\frac{\partial^2 f}{\partial x_i^2}}{\frac{\partial f}{\partial x_i}} \frac{x_i^{(k)}}{v_i^{(k)}} + \frac{1}{2} \frac{\frac{\partial^2 f}{\partial x_i^2}}{\frac{\partial f}{\partial x_i}} \frac{x_i^{(k+1)}}{v_i^{(k)}} \quad (6.94)$$

• 14.1 Author's comments

- (1) In the choice of independent variables, depend the minimum to maximum ratio of eigen values of the hessian matrix which determines convergence of the gradient technique (section 7.5). Though Hessian is too difficult to obtain, the convergence as judged from the results supplied is not good enough compared to that of 3.13 as found by the author.
- (2) Though the method has the advantages of 3.13 from 1 to 3 the algorithm needs C_4^{-1} explicitly in each iteration. This matrix for a 30 bus system is 52×52 with four generation nodes.

Method

Method of reference 21

As discussed in section 8.1.1, the gradient method of Powell and Tinney²¹ based on the technique of section 7.9 is the most promising. The author's computer programme is based on this method modified to take care of the experience mentioned in this section.^(8.13.1) Here the method of reference 21 need the cost of slack bus expressed in terms of V and θ , in this programme it is taken as function of power.

As discussed in section 8.1.1, it is more advantageous to treat the reactive power generation as independent variable, rather than voltage magnitudes. Treating the inequality constraints on voltage magnitudes for nodes with reactive power generation through penalty method has been found to result in considerable convergence difficulties since the solution oscillates between inside and outside of the feasible domain from iteration to iteration. Such constraints are therefore treated differently.

9.1 Effect of inequality constraint

Inequality constraints on the independent variable u_i can be considered by realising for the optimal solution;

$$\frac{\partial \omega}{\partial u_1} = 0 \quad \text{if } u_1 \leq \underline{u}_1 \leq \bar{u}_1 \quad (9.1)$$

$$\frac{\partial \omega}{\partial u_1} \geq 0 \quad \text{if } u_1 = \underline{u}_1 \quad (9.2)$$

$$\frac{\partial \omega}{\partial u_1} \leq 0 \quad \text{if } u_1 = \bar{u}_1 \quad (9.3)$$

Inequality constraints on the function $\omega_1(x, u)$ can be considered by restating the problem of section 7.1 as follows:

$$\min f(x, u) \quad (9.4)$$

$$\text{subject to} \quad h(x, u) = 0 \quad (9.5)$$

$$g(x, u) - v = 0 \quad (9.6)$$

$$v \leq 0 \quad (9.7)$$

where v is the vector of additional variables introduced. If the jacobian $\begin{bmatrix} \frac{\partial h_i}{\partial x} & \frac{\partial g_i}{\partial x} \end{bmatrix}^T$ is non-singular, v can be considered as independent variables.

Defining the Lagrangian as follows

$$L = f(x, u) + \sum \lambda_i h_i(x, u) + \sum \alpha_i (g_i(x, u) - v_i) \quad (9.8)$$

see the analysis of section 7.9; if the dual variables are obtained from equation 9.9 as follows,

$$\begin{bmatrix}
 \lambda_1 \\
 \vdots \\
 \lambda_k \\
 - \\
 x_1 \\
 \vdots \\
 x_n \\
 - \\
 x_1
 \end{bmatrix} =
 \begin{bmatrix}
 \frac{\partial f_1}{\partial x_1} & \dots & \frac{\partial f_1}{\partial x_n} & \frac{\partial f_1}{\partial \lambda_1} & \dots & \frac{\partial f_1}{\partial \lambda_k} \\
 \vdots & & \vdots & & & \vdots \\
 \vdots & & \vdots & & & \vdots \\
 \vdots & & \vdots & & & \vdots \\
 \vdots & & \vdots & & & \vdots \\
 \vdots & & \vdots & & & \vdots \\
 \vdots & & \vdots & & & \vdots \\
 \vdots & & \vdots & & & \vdots \\
 \vdots & & \vdots & & & \vdots \\
 \frac{\partial f_1}{\partial x_{k+1}} & \dots & \frac{\partial f_1}{\partial x_{n+1}} & \frac{\partial f_1}{\partial \lambda_1} & \dots & \frac{\partial f_1}{\partial \lambda_k}
 \end{bmatrix}^{-1} \begin{bmatrix}
 \frac{\partial f_1}{\partial x_1} \\
 \vdots \\
 \vdots \\
 \vdots \\
 \frac{\partial f_1}{\partial x_{k+1}}
 \end{bmatrix}$$

(9.9)

then

$$\frac{\partial}{\partial v_i} = -x_1 - \frac{\partial f(x, u)}{\partial x_1} \quad \text{in the constraint domain} \quad (9.10)$$

Hence for the optimal solution the following relations are satisfied:

(9.11)

$$\text{if } v_i < 0 ; \quad \alpha_i = 0$$

(9.12)

$$\text{if } v_i = 0 ; \quad \alpha_i > 0$$

9.2 Variables

The 14 known variables have been classified as follows:

- (a) Independent variables: These variables include active and reactive powers of all generation nodes treated as p q nodes, slack bus voltage magnitude and adjustable transformer tap setting.
- (b) Dependent variables: These variables are slack bus active and reactive powers, voltage magnitudes and angles of p q nodes, reactive powers and phase angles of p v nodes..

9.3 Lagrangian

The computer programme is based upon the following Lagrangian :

$$\begin{aligned}
 L = & \sum_{i \in G} f_i(p_i^G) + \text{penalty terms} \\
 & + \sum_i \lambda_i (p_i(v, \theta) - p_i^G + p_i^I) \\
 & + \sum_i \mu_i (\eta_i(v, \theta) - \eta_i^G + \eta_i^I) \\
 & + \sum_{i \in P} \pi_i (v_i - \bar{v}_i) + \sum_{i \in S} \pi_i' (v_i - \underline{v}_i) \\
 & + \alpha_S (v_r - v_s') + \beta_r (\theta_r - \theta_s') \quad (9.13)
 \end{aligned}$$

penalty terms include voltage inequality constraints on

notes other than with reactive power generation, inequality constraints on the slack bus powers and inequality constraints on voltage angle differences. After some experimentation, Powell's penalty functions of equation 7.61 have been adopted. These are restated as follows:

$$(x, \alpha) = C \quad \text{if } x \leq \bar{x}$$

$$= \frac{(x - \bar{x} + \epsilon_x^L)^2}{\epsilon_x^L} \quad \text{if } x > \bar{x}$$

$$= \frac{(x - \bar{x} + \epsilon_x^U)^2}{\epsilon_x^U} \quad \text{if } x < \underline{x} \quad (9.16)$$

Dual variables α_L and β_U have been introduced with the slack bus voltage magnitude and angle to make the Lagrangian compatible with the load flow programme of Chapter VI.

If the objective function is different from the cost of real power generation, the same is substituted in place of

$$\sum_{i \in G} f_i(r_i^0).$$

9.4 Dual Variables

The variables have been obtained by making the partial derivatives of the Lagrangian with respect to the dependent variables equal to zero as described in section 7.9.

$$\frac{\partial L}{\partial \theta_i} = C - \frac{\partial f_i(\theta_i)}{\partial \theta_i} - \lambda_i = 0 \quad (9.15)$$

$$\frac{\partial L}{\partial \theta_i} = C - \lambda_i = -\mu_i \quad (9.16)$$

The reactive power generation nodes have been considered as p-q nodes except if the voltage constraint is violated for that node, in which case these are considered as p-v nodes with the assigned voltages equal to V or \underline{V} depending upon which constraints are violated. For such nodes λ_i^* is a dependent variable.

$$\frac{\partial L}{\partial \theta_i} = 0 = u_i \text{ for p-v node} \quad (9.17)$$

The dual variables λ_i^* and $\lambda_i^{*\prime}$; $i \in R$ are zero if the voltage constraint is not reached.

Thus for every node except the slack node, there are two constraint equations and two non-zero dual variables as follows:

(a) p-q nodes (Set A)

The constraint equations are as follows:

$$h_{21-i} = p_i(V, \theta) - P_i^G + P_i^L = 0 \quad (9.18)$$

$$h_{21} = q_i(V, \theta) - Q_i^G + Q_i^L = 0 \quad (9.19)$$

Associated non-zero dual variables are λ_i and μ_i respectively.

(b) p.v nodes (Set 1)

The constraint equations are as follows:

$$n_{2i-1} = p_i(v, \epsilon) - p_i^+ + p_i^- = 0 \quad (9.20)$$

$$n_{2i} = v_i - \bar{v}_i \text{ or } v_i - \underline{v}_i = 0 \quad (9.21)$$

The non-zero dual variables are λ_1 and π_1 or π_1' .

(c) slack node (Set C)

The constraint equations included for the reasons mentioned earlier are as follows:

$$n_{2g-1} = c_g - \bar{c}_g = 0 \quad (9.22)$$

$$n_{2g} = v_g - \bar{v}_g = 0 \quad (9.23)$$

Associated dual variables are β_g and α_g respectively.

In addition the slack bus has a non-zero dual variable λ_n which is obtained from equation 9.15.

By equating the partial derivative of L with respect to v_j to zero and pre-multiplying the resulting equation by v_j ,

$$\begin{aligned}
 -\frac{\partial F}{\partial v_j} &= \lambda_1 v_j \frac{\partial h_{21}}{\partial v_j} \\
 &= \sum_{1 \leq i \leq n} (\alpha_i v_j \frac{\partial h_{21-i}}{\partial v_j} + \beta_i v_j \frac{\partial h_{21}}{\partial v_j}) \\
 &= \sum_{1 \leq i \leq n} (\alpha_i v_j \frac{\partial h_{21-i}}{\partial v_j} + \gamma_i^j v_j \frac{\partial h_{21}}{\partial v_j}) \\
 &= \sum_{1 \leq i \leq n} (\beta_i v_j \frac{\partial h_{21-i}}{\partial v_j} + \alpha_1 v_j \frac{\partial h_{21}}{\partial v_j}) \quad (9.24-j)
 \end{aligned}$$

where γ is the objective function augmented with the penalty terms.

$\gamma_i^j = \pi_i$ or π_i' depending upon which constraint is active.

Similarly by equating the partial derivative of Lagrangian with respect to e_j to zero, resulting equation is as follows:

$$\begin{aligned}
 -\frac{\partial F}{\partial e_j} - \lambda_1 \frac{\partial p_a}{\partial e_j} &= \sum_{1 \leq i \leq n} (\alpha_i \frac{\partial h_{21-i}}{\partial e_j} + \beta_i \frac{\partial h_{21}}{\partial e_j}) \\
 &+ \sum_{1 \leq i \leq n} (\alpha_i \frac{\partial h_{21-i}}{\partial e_j} + \gamma_i^j \frac{\partial h_{21}}{\partial e_j}) \\
 &+ \sum_{1 \leq i \leq n} (\beta_i \frac{\partial h_{21-i}}{\partial e_j} + \alpha_1 \frac{\partial h_{21}}{\partial e_j}) \quad (9.25-j)
 \end{aligned}$$

Thus the $2n$ non-zero dual variables can be obtained from equations 9.24-j and 9.25-j, $j = 1, \dots, n$. The associated

try to find transpose of the Jacobian used in load flow solution. The non-zero dual variables have been obtained by main use of the table of factor obtained during the load flow solution and the computer logic described in section 2.17.

9.5 Sensitivity of the augmented objective function

The gradient vector of the augmented objective function with respect to the independent variables, in the constraint domain of power flow equations is as follows:

$$\frac{\partial F}{\partial t_i} = \frac{\partial G}{\partial t_i} - \lambda_1 \quad \text{for } i \in S; i \neq s \quad (9.26)$$

$$\frac{\partial F}{\partial t_i} = -\gamma_1 \quad \text{for } i \in N \cup R \quad (9.27)$$

$$\frac{\partial F}{\partial t_i} = -\gamma_1^T \quad \text{for } i \in E \cup R \quad (9.28)$$

$$\frac{\partial F}{\partial t_h} = \lambda_1 \frac{\partial P_1}{\partial t_h} + \lambda_j \frac{\partial P_{ij}}{\partial t_h} + \mu_1 \frac{\partial Q_{ij}}{\partial t_h} + \mu_2 \frac{\partial Q_{h2}}{\partial t_h}$$

$$= G_{h1} t_h - G_{h2}$$

where G_{h1} and G_{h2} are respectively given by equations 8.75 and 8.76 respectively.

$$\begin{aligned} \frac{\partial F}{\partial V_h} &= -\lambda_1 v_i v_j y_{ij} \sin(\theta_i - \theta_j - \psi_{ij} + \gamma_h) \\ &\quad + \lambda_j v_i v_j y_{ij} \sin(\theta_j - \theta_i - \psi_{ij} - \gamma_h) \\ &\quad + \mu_1 v_i v_j y_{ij} \cos(\theta_i - \theta_j - \psi_{ij} + \gamma_h) \end{aligned}$$

$$= p_j V_i V_j y_{ij} \cos(\theta_j - \theta_i - \Psi_{ij} - \delta_n) \quad (9.3)$$

or the optimal solution equation 9.1 to 9.3 hold, so
or the bus are with reactive power generation, rechein the
voltage constraint; equation 9.21 holds

$$\pi_i^1 \geq 0 \quad \text{if } V_i = \bar{V}_i \quad (9.1)$$

$$\pi_i^1 \leq 0 \quad \text{if } V_i = \underline{V}_i$$

$$\frac{\partial L}{\partial \pi_i} = \lambda_i \quad (9.22)$$

With the increase in load demand, operating cost increases.
Thus $\lambda_i > 0$ for all i if the objective function is the
operating cost.

9.6 Direction vector

The direction vector can be found out by any of the methods discussed in sections 7.5, 7.6 and 7.7. The conjugate gradient and Fletcher-Powell's gradient methods guarantee the minimum can be located exactly within certain finite number of iterations for the quadratic function. Even in regions remote from minimum, these methods by taking account of the curvature of the function are expected to deal with complex situations such as presence of a long curving valley^{43,44}. As discussed

earlier conjugate gradient method needs only very little additional memory space compared to the steepest descent method, this method has therefore been adopted. The new direction vector has been obtained from equations 7.14 and 7.15.

Due to non-linear tractive on the decision variable; following modification has been made-

If the decision variable u_i^k at kth iteration is at the constraint limit and sign of gradient satisfies equations 9.2 and 9.3; the variable is kept fixed. For such variable and $i = l$; the variable is kept fixed. For such variable and $i = l$; the corresponding $\frac{\partial f}{\partial u_i^k}$ term is not added while obtaining $d_i^k = l$.

$\nabla f(u^k) \neq \nabla f(u^{k-1})$. If sign of the gradient for the variable at constraint limit do not satisfy equations 9.2 and 9.3; the following value of d_i^{k-1} is substituted

$$d_i^{k-1} = \frac{u_i^k - u_i^{k-1}}{r_e} \quad (7.13)$$

For bus bars with controllable reactive powers with constraints on V and θ , V has normally been taken as an independent variable, since this has been found to result in better convergence. For such bus bars a provision has been made in the load flow solution to check if voltage constraint is violated in which case the bus bar is considered as p.v node with the bus voltage magnitude equal to the constraint value, provided it is within the constraint limits (section 6.1).

or when node two gradient has been approximately estimated as follows:

$$\frac{\partial}{\partial r} \left(\frac{\partial V_1}{\partial r} \right) = \frac{1}{r_1^2} \frac{V_1 - V_{11}}{r_1 - r_{11}}$$

9.7 optimal step length (cf section 7.0.1)

In order to find the optimal step length through parabolic extrapolation as done in reference 21 (section 7.0.1), $y(0)$, $y'(r_1)$ and $y(r)$ need be obtained and r_e is obtained from equation (7.0.1).

In order to obtain $y(r)$ a load flow solution is needed. $y'(r)$ can be obtained by just one additional solution of the simultaneous equations for which table of factor has already been obtained in compact form. With these four values, cubical extrapolation is now possible to obtain a more accurate value of r_e . Even this was not found working satisfactorily, hence Livindon's cubical interpolation described in section 7.0.2 had to be used. This method needs $y(r)$ and $y'(r)$ for $r = 0, h, 2h, 4h, \dots, a, b$ where r is doubled each time and b is the first of these values at which y' is non-negative or y has not decreased. For every such value of r , a load flow solution and one additional regularization is needed. Hence it is desirable that $y'(r)$ is non-negative in as few trials as possible. After obtaining the direction vector, b is obtained on the basis of maximum change in the independent variable. After

obtaining the value of r , maximum change in the decision variable is reobtained. For the next trial the maximum change in the independent variable is assumed 'a' ($a < 1$) times the maximum change of this trial.

$$\begin{aligned} h_p &= \frac{\Delta_r}{\max(d_1)} \\ h_q &= \frac{\Delta_r}{\max(d_1)} \\ h_t &= \frac{\Delta_t}{\max(d_1)} \end{aligned} \quad (9.35)$$

where Δ_r , Δ_r and Δ_t are the estimated maximum changes in real power, reactive power and tap settings. The value of Δ_r is the minimum of h_p , h_q and h_t . For the next iteration ΔP_1 , Δ_r and Δ_t are estimated as follows:

$$\begin{aligned} \Delta P_1^{\text{new}} &= (a \Delta P_1^{\text{old}} r^{**}/h_p \\ \Delta_r^{\text{new}} &= (a \Delta_r^{\text{old}} r^{**}/h_q \\ \Delta_t^{\text{new}} &= (a \Delta_t^{\text{old}} r^{**}/h_t \end{aligned} \quad (9.36)$$

Typical initial values taken are $\Delta_r = .5$ pu, $\Delta_t = .1$ pu, $\Delta_t = .1$ and $a = .8$.

y/r is obtained from equation 7.24 except in case when some of independent variables violate the constraint, the following modification is made:

$$y'(r) = \nabla r^T \nabla (u + r d) \quad (9.37)$$

$$= \sum_i \gamma_i d_i \quad (9.38)$$

where γ_i is the gradient obtained as in section 9.5 if

$$v_i = u_i + r d_i$$

$$\text{and } u \leq v_i \leq \bar{u}_i.$$

If $v_i < u_i$ or $v_i > \bar{u}_i$, the new value of the independent variable is taken equal to the bounded value and $r_i = 0$ if $y'(r) < 0$.

Thus in order to obtain $y'(r)$; $\sum_{i \in F} \gamma_i d_i$ is obtained

for the independent variables within the constraint limit.

$\sum_{i \notin F} \gamma_i d_i$ for the variables violating the constraints is obtained

separately with γ_i as the gradient at the constraint limit.

The two are added if $\sum_{i \in F} \gamma_i d_i$ is non-negative.

In case load flow programme fails to converge; a solution with half the changes in the independent variables is attempted. Convergence of load flow solution is judged on the basis of maximum mismatch in power as described in section 6.3.

9.6 System Description

The system extensively tested is a "C-bus system", description of which is as follows.

$$\text{Base MVA} = 100$$

Table 9.6.1: Sequence and line charging data

| Line connection | Resistance per unit | Reactance per unit | Line charging per unit |
|--------------------|------------------------|-----------------------|---------------------------|
| 1 2 | 0.0192 | 0.6575 | 0.0574 |
| 1 3 | 0.0452 | 0.1472 | 0.0202 |
| 2 4 | 0.0570 | 0.1717 | 0.0184 |
| 2 5 | 0.0472 | 0.1963 | 0.0209 |
| 2 6 | 0.0581 | 0.1763 | 0.0107 |
| 2 7 | 0.0172 | 0.0379 | 0.0042 |
| 4 6 | 0.0111 | 0.0114 | 0.0045 |
| 4 12 | 0.0000 | 0.0256 | 0.0000 |
| 5 7 | 0.0460 | 0.1160 | 0.0102 |
| 6 7 | 0.0267 | 0.0020 | 0.0087 |
| 6 8 | 0.0120 | 0.0420 | 0.0045 |
| 6 9 | 0.0000 | 0.2000 | 0.0000 |
| 6 10 | 0.0000 | 0.5560 | 0.0000 |
| 6 28 | 0.0169 | 0.0799 | 0.0065 |
| 8 28 | 0.0636 | 0.2000 | 0.0214 |
| 9 10 | 0.0030 | 0.1100 | 0.0000 |
| 9 11 | 0.0000 | 0.2000 | 0.0000 |
| 10 17 | 0.0324 | 0.0845 | 0.0000 |
| 10 20 | 0.0936 | 0.2090 | 0.0000 |
| 10 21 | 0.0348 | 0.0749 | 0.0000 |
| 10 22 | 0.0727 | 0.1499 | 0.0000 |
| 12 13 | 0.0000 | 0.1400 | 0.0000 |
| 12 14 | 0.1231 | 0.2559 | 0.0000 |
| 12 15 | 0.0662 | 0.1304 | 0.0000 |
| 12 16 | 0.0945 | 0.1987 | 0.0000 |
| 14 15 | 0.2210 | 0.1997 | 0.0000 |
| 15 18 | 0.1073 | 0.2185 | 0.0000 |
| 16 23 | 0.1000 | 0.2020 | 0.0000 |
| 16 17 | 0.0824 | 0.1923 | 0.0000 |
| 18 19 | 0.0639 | 0.1292 | 0.0000 |
| 19 20 | 0.0240 | 0.0680 | 0.0000 |
| 21 22 | 0.0116 | 0.0236 | 0.0000 |
| 22 24 | 0.1150 | 0.1790 | 0.0000 |
| 23 24 | 0.1320 | 0.2700 | 0.0000 |
| 24 25 | 0.1885 | 0.3292 | 0.0040 |
| 25 26 | 0.2544 | 0.3800 | 0.0000 |

Table 9.0.1 (contd..)

| <u>line connection</u> | <u>current per unit</u> | <u>reactance per unit</u> | <u>line charging per unit</u> |
|----------------------------|-----------------------------|-------------------------------|-----------------------------------|
| 25 27 | 0.1693 | 0.2037 | 0.0000 |
| 27 28 | 0.0000 | 0.2960 | 0.0000 |
| 27 29 | 0.21 | 0.4173 | 0.0000 |
| 27 30 | 0.02 | -0.627 | 0.0000 |
| 28 30 | 0.2099 | 0.4173 | 0.0000 |

Table 9.0.2: Static capacitor data

| <u>bus bar r</u> | <u>uncertainty per unit</u> |
|------------------|---------------------------------|
| 10 | 0.1900 |
| 24 | 0.4300 |

Table 9.0.3: Transformer data (Initial values)

| <u>transformer designation</u> | <u>no rating</u> |
|------------------------------------|----------------------|
| 4 12 | 0.932 |
| 6 9 | 0.978 |
| 6 10 | 0.969 |
| 28 27 | 0.965 |

Table 9. 4

App. 9. 4 shows the obtaining turned in the subroutine
 to find the solution of the system of linear equations. In
 the next form of obtaining turned the sum of the rows in
 the system.

Table 9. 5

| ode | row number | Incons. elements of admittance matrix | | 1 | 2 |
|-----|---------------|--|-----------|----|----|
| | | real | imaginary | | |
| | | rt | - rt | | |
| 1 | 1 | 6.463 | -20.795 | 1 | 1 |
| 2 | 16 | 9.742 | -36.647 | 6 | 2 |
| 3 | 14 | 9.439 | -24.602 | 7 | 3 |
| 4 | 28 | 16.314 | -54.405 | 9 | 5 |
| 5 | 27 | 6.379 | -12.196 | 10 | 6 |
| 6 | 29 | 27.341 | -62.174 | 15 | 8 |
| 7 | 13 | 6.544 | -13.476 | 13 | 10 |
| 8 | 12 | 7.773 | -26.527 | 16 | 11 |
| 9 | 11 | 0.000 | -18.706 | 18 | 12 |
| 10 | 28 | 12.462 | -41.330 | 22 | 14 |
| 11 | 2 | 0.000 | -6.007 | 22 | 15 |
| 12 | 30 | 6.573 | -24.424 | 26 | 16 |
| 13 | 2 | 0.000 | -7.142 | 26 | 17 |
| 14 | 11 | 4.017 | -8.424 | 27 | 18 |
| 15 | 27 | 0.765 | -16.011 | 29 | 20 |
| 16 | 25 | 3.034 | -8.497 | 30 | 21 |
| 17 | 10 | 5.838 | -14.710 | 30 | 23 |
| 18 | 9 | 4.886 | -9.906 | 31 | 24 |
| 19 | 24 | 0.958 | -12.903 | 31 | 25 |
| 20 | 1 | 0.258 | -15.750 | 32 | 27 |
| 21 | 7 | 7.657 | -45.103 | 32 | 28 |
| 22 | 31 | 21.77 | -45.432 | 33 | 30 |
| 23 | 20 | 21.316 | -6.956 | 34 | 31 |
| 24 | 6 | 7.421 | -8.801 | 35 | 31 |
| 25 | 19 | 5.511 | -8.064 | 36 | 33 |
| 26 | 18 | 4.495 | -1.817 | 38 | 34 |
| 27 | 1 | 1.116 | -9.460 | 38 | 35 |
| 28 | 17 | 1.152 | -22.342 | 41 | 36 |
| 29 | 1 | 5.06 | -0.604 | 41 | 39 |
| 30 | 1 | 1.907 | -0.017 | 42 | 40 |
| | 1 | 1.599 | | | |

Table 10

DF diagonal element

| | | | | real part | imaginary part |
|----|----|----|----|-----------|----------------|
| 1 | 2 | 1 | 1 | -0.724 | 15.646 |
| 2 | 3 | 1 | 2 | -1.243 | 5.096 |
| 3 | 6 | 2 | 6 | -1.705 | 5.197 |
| 4 | 5 | 2 | 4 | -1.135 | 4.772 |
| 5 | 6 | 2 | 5 | -1.616 | 5.116 |
| 6 | 4 | 4 | 7 | -0.190 | 23.530 |
| 7 | 6 | 4 | 9 | -6.410 | -2.111 |
| 8 | 7 | 5 | 10 | 0.000 | 2.460 |
| 9 | 7 | 7 | 11 | -2.956 | 7.440 |
| 10 | 8 | 6 | 12 | -0.239 | 22.012 |
| 11 | 9 | 8 | 13 | 0.000 | 4.701 |
| 12 | 9 | 9 | 15 | 0.000 | 1.742 |
| 13 | 10 | 9 | 17 | -6.762 | 15.463 |
| 14 | 28 | 7 | 1 | -1.443 | 4.540 |
| 15 | 23 | 8 | 22 | 0.000 | 9.590 |
| 16 | 10 | 1 | 22 | 0.000 | 4.007 |
| 17 | 11 | 12 | 23 | -1.956 | 10.317 |
| 18 | 17 | 12 | 24 | -1.784 | 3.985 |
| 19 | 20 | 14 | 25 | -5.101 | 10.980 |
| 20 | 21 | 17 | 25 | -2.610 | 5.400 |
| 21 | 22 | 10 | 18 | 0.000 | 7.147 |
| 22 | 11 | 16 | 29 | -1.526 | 2.173 |
| 23 | 14 | 15 | 27 | -3.095 | 6.097 |
| 24 | 17 | 18 | 30 | -1.51 | 4.104 |
| 25 | 16 | 10 | 19 | -2.490 | 2.250 |
| 26 | 15 | 19 | 31 | -1.010 | 3.687 |
| 27 | 18 | 10 | 20 | -1.968 | 3.976 |
| 28 | 21 | 10 | 21 | -1.882 | 4.393 |
| 29 | 17 | 21 | 32 | -3.079 | 6.111 |
| 30 | 15 | 15 | 28 | -5.002 | 11.764 |
| 31 | 20 | 22 | 33 | -16.774 | 34.127 |
| 32 | 22 | 23 | 24 | -2.540 | 7.954 |
| 33 | 21 | 24 | 35 | -1.461 | 2.989 |
| 34 | 21 | 25 | 26 | -1.309 | 2.287 |
| 35 | 25 | 25 | 37 | -1.216 | 1.317 |
| 36 | 26 | 6 | 14 | -1.965 | 3.760 |
| 37 | 27 | 8 | 15 | 0.000 | 2.444 |
| 38 | 28 | 27 | 38 | -0.995 | 1.881 |
| 39 | 29 | 27 | 39 | -6.687 | 1.293 |
| 40 | 30 | 27 | 40 | -6.912 | 1.723 |
| 41 | 36 | 29 | 41 | 0.000 | 0.000 |
| 42 | 30 | 30 | 42 | 0.000 | 0.000 |

where,

i_1 = index sequence of corresponding row, right off diagonal elements

i_2 = index sequence of corresponding row, left off diagonal element.

i_1 = column of corresponding right off diagonal element

i_2 = column of corresponding left off diagonal element

i = corresponding location of left off diagonal element in the array of right off diagonal elements.

Table 9.6.5: Transformer tap admittance

| Transformer designation | i_t | y_t | y_{td} |
|-------------------------|-------|--------------|--------------|
| 4 12 | 8 | 0.0+j3.90625 | 0.0-j3.90625 |
| 6 9 | 12 | 0.0+j4.80769 | 0.0-j4.80769 |
| 6 10 | 13 | 0.0+j1.79856 | 0.0-j1.79856 |
| 28 27 | 38 | 0.0+j2.52525 | 0.0-j2.52525 |

where, i_t = location of right off diagonal element in Table 9.6.4-b

$y_{td} = (y + j3)$ of equation 3.3

$y_t = (-y)$ of equation 3.3.

Table 9. 5: Initial or rating conditions

| Bus number | gen. unit | bus volt. re ds rear | description | | Load | |
|---------------|--------------|----------------------------|-------------|-------|--------|--------|
| | | | Mvar | Mvar | MVA | Mvar |
| 1 | 1.60 | 0.000 | 0.000 | 0.000 | 0.000 | 0.000 |
| 2 | 1.000 | 0.000 | 40.000 | 0.000 | 1.700 | 12.700 |
| 3 | 1.000 | 0.000 | 0.000 | 0.000 | 2.400 | 1.700 |
| 4 | 1.00 | 0.000 | 0.000 | 0.000 | 7.600 | 1.700 |
| 5 | 1.700 | 0.000 | 0.000 | 0.000 | 94.200 | 1.000 |
| 6 | 1.000 | 0.000 | 0.000 | 0.000 | 0.000 | 0.000 |
| 7 | 1.000 | 0.000 | 0.000 | 0.000 | 22.000 | 10.900 |
| 8 | 1.000 | 0.000 | 0.000 | 0.000 | 30.000 | 30.000 |
| 9 | 1.000 | 0.000 | 0.000 | 0.000 | 0.000 | 0.000 |
| 10 | 1.000 | 0.000 | 0.000 | 0.000 | 5.300 | 2.000 |
| 11 | 1.000 | 0.000 | 0.000 | 0.000 | 0.000 | 0.000 |
| 12 | 1.000 | 0.000 | 0.000 | 0.000 | 0.000 | 0.000 |
| 13 | 1.000 | 0.000 | 0.000 | 0.000 | 6.200 | 1.600 |
| 14 | 1.000 | 0.000 | 0.000 | 0.000 | 0.200 | 2.500 |
| 15 | 1.000 | 0.000 | 0.000 | 0.000 | 3.500 | 1.800 |
| 16 | 1.000 | 0.000 | 0.000 | 0.000 | 9.000 | 5.700 |
| 17 | 1.000 | 0.000 | 0.000 | 0.000 | 3.200 | 0.900 |
| 18 | 1.000 | 0.000 | 0.000 | 0.000 | 9.500 | 3.400 |
| 19 | 1.000 | 0.000 | 0.000 | 0.000 | 2.200 | 0.700 |
| 20 | 1.000 | 0.000 | 0.000 | 0.000 | 17.500 | 11.200 |
| 21 | 1.000 | 0.000 | 0.000 | 0.000 | 0.000 | 0.000 |
| 22 | 1.000 | 0.000 | 0.000 | 0.000 | 3.200 | 1.600 |
| 23 | 1.000 | 0.000 | 0.000 | 0.000 | 6.700 | 6.700 |
| 24 | 1.000 | 0.000 | 0.000 | 0.000 | 0.000 | 0.000 |
| 25 | 1.000 | 0.000 | 0.000 | 0.000 | 3.500 | 2.300 |
| 26 | 1.000 | 0.000 | 0.000 | 0.000 | 0.000 | 0.000 |
| 27 | 1.000 | 0.000 | 0.000 | 0.000 | 0.000 | 0.000 |
| 28 | 1.000 | 0.000 | 0.000 | 0.000 | 2.400 | 0.900 |
| 29 | 1.000 | 0.000 | 0.000 | 0.000 | 10.600 | 1.900 |
| 30 | 1.000 | 0.000 | 0.000 | 0.000 | | |

* "lack bus.

3.3 Example

Assume active power generation at buses 1, 2 and 11 and reactive power generations at buses 1, 2, 5, 7, 11 and 13 and assume following generation costs;

$$f_1(P_1^0) = .7 P_1^0 + .01(P_1^0)^2$$

$$f_2(P_2^0) = .7 P_2^0 + .02(P_2^0)^2$$

$$f_{11}(P_{11}^0) = .8 P_{11}^0 + .02(P_{11}^0)^2$$

With the reactive power generations $-0.5 \leq Q_i \leq 0.5$ and the line voltage magnitudes $1.0 \leq V \leq 1.15$; results are given below. Slack bus voltage is also considered decision variable.

Table 9.9.1: with 1 operating conditions

| Tap Number | V | S | pu | pu |
|------------|---------|-----------|----------|----------|
| 1 | 1.00000 | -0.10000 | 2.62209 | 0.14554 |
| 2 | 1.02773 | -5.21097 | 0.18300 | -0.12700 |
| 3 | 1.05617 | -0.10003 | -0.02397 | -0.01150 |
| 4 | 1.08461 | -0.77967 | -0.07600 | -0.01600 |
| 5 | 1.11275 | -16.5142 | -0.14200 | 0.20727 |
| 6 | 1.14089 | -11.5142 | -0.05000 | -0.00500 |
| 7 | 1.16845 | -1.75150 | -0.02800 | -0.16500 |
| 8 | 1.19622 | -12.2153 | -0.10000 | -0.34920 |
| 9 | 1.22399 | -15.22470 | 0.00000 | 0.00000 |
| 10 | 1.25176 | -17.1534 | -0.05700 | -0.02000 |
| 11 | 1.27953 | -19.0276 | 0.00000 | 0.4508 |
| 12 | 1.30730 | -16.17507 | -0.11200 | -0.37800 |
| 13 | 1.33507 | -16.17507 | 0.00000 | 0.16343 |
| 14 | 1.36284 | -17.2772 | -0.05199 | -0.01600 |
| 15 | 1.39061 | -17.63490 | -0.08200 | -0.02400 |
| 16 | 1.41838 | -16.92746 | -0.03500 | -0.01800 |
| 17 | 1.44615 | -17.37614 | -0.09000 | -0.05799 |
| 18 | 1.47392 | -18.21161 | -0.03200 | -0.06900 |
| 19 | 1.50169 | -0.09546 | -0.09500 | -0.04400 |
| 20 | 1.52946 | -1.4660 | -0.02199 | -0.00700 |
| 21 | 1.55723 | -13.15274 | -0.17000 | -0.11200 |
| 22 | 1.58499 | -17.80529 | 0.00000 | 0.00000 |
| 23 | 1.61276 | -17.75118 | -0.03200 | -0.01600 |
| 24 | 1.64053 | -16.67487 | -0.08699 | -0.06700 |
| 25 | 1.66830 | -19.6214 | 0.00000 | 0.00000 |
| 26 | 1.69607 | -16.12974 | -0.03500 | -0.02299 |
| 27 | 1.72384 | -16.57549 | 0.00000 | 0.00000 |
| 28 | 1.75161 | -16.02239 | 0.00000 | 0.00000 |
| 29 | 1.77938 | -12.24857 | -0.02399 | -0.00900 |
| 30 | 1.80715 | -10.26488 | -0.10599 | -0.01900 |
| | 1.83492 | -19.23135 | | |

With the above operating conditions obtained at the tap settings of Table 9.9.1; the objective function without penalty function = 2.16741 pu and with penalty function = 2.20042.

Optimal loading conditions and the corresponding tap settings are given in Tables 9.9.2 and 9.9.3 respectively.

Table 2.6.1: Estimated tin corrections

| | α | β | γ | δ |
|----|------------|-----------|-----------|-----------|
| 1 | 1.0 551 | -1.00000 | 1.0 7723 | -1.0 6712 |
| 2 | 1.0 14 | -1.09257 | 0.02304 | -0.0157 |
| 3 | 1.0 7334 | -1.04427 | -0.01399 | -0.0111 |
| 4 | 1.0 6493 | -1.02703 | -0.01790 | -0.01700 |
| 5 | 1.0 57116 | -1.01900 | -0.016200 | 0.0 1600 |
| 6 | 1.0 57117 | -1.01763 | 0.01600 | 0.01600 |
| 7 | 1.0 5746 | -1.01649 | -0.012800 | -0.01090 |
| 8 | 1.0 57043 | -1.01544 | -0.013000 | 0.0 5776 |
| 9 | 1.0 57044 | -1.01566 | 0.01600 | 0.01600 |
| 10 | 1.0 57 01 | -1.01571 | -0.015799 | -0.012000 |
| 11 | 1.0 57391 | -1.01552 | 0.01564 | -0.0111 |
| 12 | 1.0 57 7 | -1.01590 | -0.011200 | -0.01700 |
| 13 | 1.0 57 52 | -1.015909 | 0.00000 | 0.04740 |
| 14 | 1.0 57 11 | -1.015170 | -0.00199 | -0.0115 |
| 15 | 1.0 5734 | -1.015356 | -0.008200 | -0.02700 |
| 16 | 1.0 57031 | -1.015356 | -0.03500 | -0.0100 |
| 17 | 1.0 57 11 | -1.01570 | -0.03000 | -0.01900 |
| 18 | 1.0 57318 | -1.01447 | -0.09500 | -0.01400 |
| 19 | 1.0 57 69 | -1.01447 | -0.02199 | -0.00700 |
| 20 | 1.0 5792 | -1.014545 | -0.17000 | -0.11200 |
| 21 | 1.0 57777 | -1.014545 | 0.00000 | 0.00000 |
| 22 | 1.0 57 119 | -1.014545 | -0.03200 | -0.01700 |
| 23 | 1.0 57710 | -1.014545 | -0.05600 | -0.06700 |
| 24 | 1.0 11166 | -1.014545 | 0.00000 | 0.00000 |
| 25 | 1.0 57 01 | -1.014545 | -0.00000 | -0.02299 |
| 26 | 1.0 57373 | -1.014545 | 0.00000 | 0.00000 |
| 27 | 1.0 58473 | -1.014545 | 0.00000 | 0.00000 |
| 28 | 1.0 6980 | -1.014545 | -0.02399 | -0.00900 |
| 29 | 1.0 6612 | -1.014545 | -0.10599 | -0.01900 |
| 30 | 1.0 5536 | -1.014545 | -0.10599 | -0.01900 |

Table 5.2.3: Optimal settings for optimal operation:

| <u>Bus number</u> | <u>Optimal setting</u> |
|-------------------|------------------------|
| 4 | 0.961 |
| 5 | 0.371 |
| 6 | 0.97741 |
| 7 | 1.0002 |
| 8 | 27 |

The value of objective function with and without penalty function value at optimal operating conditions is found to be 2.15905.

Table 5.2.4:

Value of initial & optimal operating conditions.

Table 5.2.4-1: Initial operating conditions

| <u>Bus number</u> | <u>Initial value</u> | <u>Optimal value</u> |
|-------------------|----------------------|----------------------|
| 1 | 2.62209 | 1.97725 |
| 2 | 0.40000 | 0.6594 |
| 3 | 0.00000 | 0.33684 |
| 4 | | |

Table 9.9.4-a: Reactive power generations:

| <u>b</u> | <u>Initial value</u> | <u>Optimal value</u> |
|----------|----------------------|----------------------|
| 1 | 0.14564 | -0.6772 |
| 2 | 0.0236 | -0.11142 |
| 3 | 0.01923 | 0.50000 |
| | 0.26937 | 0.36974 |
| 11 | -0.24708 | -0.02051 |
| 12 | 0.19843 | 0.04746 |

Table 9.9.4-b: Transformer tap settings:

| <u>Transformer connection</u> | <u>Initial value</u> | <u>Optimal value</u> |
|-------------------------------|----------------------|----------------------|
| 3 12 | 0.97200 | 0.97961 |
| 4 9 | 0.97800 | 0.98071 |
| 6 10 | 0.96900 | 0.97721 |
| 2 27 | 0.96800 | 1.00802 |

Table 9.9.4-c: Slack bus voltage:

Initial value = 1.00000

Optimal value = 1.08521

Table 7.10.1: Objective Functions

| | <u>Initial value</u> | <u>Final value</u> |
|--|----------------------|--------------------|
| objective function without penalty function | 2.14741 | 2.15205 |
| objective function with penalty function | 2.15042 | 2.15905 |

7.10 Example 2:

Another application for which the computer programs was tested is planning for additional reactive power equipment needed to maintain the minimum prescribed voltage on the buses. The objective function is as follows:

$$f = \sum (\text{additional Mvar needed}) \quad (7.39)$$

The total additional capacity is to be made minimum by adjusting the controllable reactive power generation.

Table 7.10.1:

Results of initial and optimal operating conditions

Table 1.1.1-1 Initial and final values of v for various nodes

| <u>node</u> | <u>Initial v-value</u> | <u>Final v-value</u> |
|-------------|---|---------------------------------------|
| 1 | -0.0002 | -0.0577 |
| 2 | -0.0001 | -0.14234 |
| 3 | -0.33562 | -0.236.1 |
| 4 | -0.12.71 | -0.26552 |
| 11 | 0.0.06 | -0.5272 |
| 1 | -0.0001 | -0.46516 |

Table 1.1.1-2 Kvar needed at bus 1

| <u>node</u> | <u>Initial value</u> | <u>Final value</u> |
|-------------|--------------------------|------------------------|
| 1 | 0.06434 | - |
| 12 | 0.16163 | - |
| 14 | 0.04093 | - |
| 15 | 0.01040 | - |
| 16 | 0.07235 | - |
| 18 | 0.02406 | - |
| 19 | 0.04006 | - |
| 26 | 0.00260 | - |
| 32 | 0.06111 | 0.01052 |

6.11 Convergence

In the paper the code written is user oriented. Only one algorithm, in which objective function, number of the constraints, iteration and cost of generation and loadity functions are defined, is given. It defines for a different objective function. The program is able to add value to the utility companies. The program is able to add value to the utility companies. After each iteration, all objective functions and for different after each iteration, all objective functions and for different operating conditions are given. Author's comment for the following question are given.

- (1) As I wrote out in the objective function is very fast in first two or three iterations. In subsequent iterations, the movement becomes slow. In most cases there is no noticeable improvement after four or five iterations, even though there is a noticeable change in the operating conditions. For the example of section 4; the number of controllable variables is 10. The conjugate gradient method adopted in the program, should need 12 iterations for a quadratic function, but in this case there is no noticeable change in the objective function after 4 iterations. Near the optimal point, a wide range of operating conditions give nearly the same objective function.
- (2) If the limit for voltage is controllable, the convergence is affected due to the gradients since if in any iterations, upper limit of the voltage magnitude is not violated, it will lead to the either in the next iteration for which the violation to be either in the next iteration for which the upper limit will be violated for many buses and the slack upper limit will be violated for many buses and the slack

Such a situation can be expressed in the following
situation. In all cases more efficient criteria
than in this case solution is the movement in the
objective function rather than gradient.

-x-

CHAPTER XSCHEDULING OF HYDROTHERMAL SYSTEMS

This chapter describes algorithms developed by the author for optimum scheduling of combined hydrothermal system, making use of system variables in discrete form. Because of the limited amount of water available for hydrogeneration in a given period of time, the problem is that of dynamic optimization. It appears convenient to break the problem to that of long term planning and short term planning.

10.1 Long-term planning:

This involves planning for the whole year which is divided into 12 or 52 subintervals. It is assumed that the water inflow and load demand is known with complete certainty. The problem has been transformed into -stage decision process. Superscript m denotes the stage or subinterval number.

The fuel cost of the thermal plants is as follows

$$T \sum_{i \in S} \sum_m f_i (P_i^m) \quad (10.1)$$

where P_i^m is the average power output during the subinterval m . T is the time of each subinterval.

Disregarding the water overflow the water continuity relation is as follows

$$Y_i^m = Y_i^{m-1} + J_i^m - I_i^m \quad \text{for } i \in H \quad (10.2)$$

where

Y_i^m = water storage at the end of m th subinterval

J_i^m = water inflow into the reservoir during m th subinterval

I_i^m = water discharge through turbine during m th subinterval.

The average hydro power may be expressed as follows

$$P_i^m = P_i(Y_i^m, Y_i^{m-1}, I_i^m) \quad (10.3)$$

Nerita²⁹ et al. have given the following expression for a conventional hydro plant

$$P_i^m = H_i \left[1 + \frac{C_i}{2} (Y_i^m + Y_i^{m-1}) \right] (I_i^m - q_i)/T \quad (10.4)$$

Power flow relation is as follows

$$\sum_i P_i^m - P_D^m - P_L^m = 0 \quad (10.5)$$

where P_D^m is the total average load power and P_L^m is the average loss expressed as functions of P_i^m as follows

$$P_L^m = \sum_j \sum_i P_i^m \alpha_{ij} P_j^m \quad (10.6)$$

Additional constraints include upper and lower limits on I_i^m , Y_i^m and P_i^m . For hydro plants initial and final values of water storage Y_i^0 and Y_i^T is also prescribed.

10.1.1 Variables

Unknown variables have been classified as follows.

Independent variables. These variables include thermal generations of all nodes except that of slack node and water discharges through turbines of hydro plants in all subintervals, except of first which may be obtained as follows.

$$I_1^1 = Y_1^0 - Y_1^1 + \sum_m J_1^m - \sum_{m=2}^M I_1^m \quad (10.7)$$

Dependent variables. These variables include thermal generation of slack bus, water discharges of the hydro plants during the first subinterval, hydrogenerations and water storages.

10.1.2 Nonlinear programming formulation

The problem has been solved by the method of section 7.9. An initial set of independent variables have been assumed from which the dependent variables calculated. The dual variables are obtained by equating the partial derivatives of the Lagrangian with respect to the dependent variables to zero. Inequality constraints on the dependent variables are handled through penalty function approach making use of Powell's functions.

$$\begin{aligned}
 L = & \sum_m \sum_{i \in S} [f_i(P_i^m) + \lambda(P_i^m)] + \sum_m \sum_{i \in H} [-\gamma(Y_i^m) + \lambda(P_i^m)] \\
 & + \sum_{i \in H} [w(I_i^1)] + \sum_m z_1^m (\sum P_i^m - P_D^m - P_U^m) \\
 & + \sum_m \sum_{i \in H} z_{2i}^m (Y_i^m - Y_i^{m-1} - J_i^m + I_i^m) \\
 & + \sum_m \sum_{i \in H} z_{3i}^m [P_i^m - \gamma_i(Y_i^m, Y_i^{m-1}, I_i^m)] \quad (10.e)
 \end{aligned}$$

where subscript s refers to the slack bus with thermal generation. (\cdot) is the penalty function of x defined by equation (x.14). Dual variables z_1^m , z_{21}^m and z_{31}^m are obtained by equating the partial derivatives of the Lagrangian with respect to the dependent variables to zero which yield the following equations.

$$\frac{\partial \mathcal{L}}{\partial z_1^m} = 0 = \psi(\bar{x}_1) + z_1^m + \frac{z_1^m}{d\bar{x}_1^m} + \frac{\partial P_1(\bar{x}_1)}{\partial \bar{x}_1^m} \quad (10.10-s-1)$$

$$\frac{\partial \mathcal{L}}{\partial z_{21}^m}; i.e. = 0 = \psi(\bar{x}_1) + z_{21}^m \left(1 - \frac{\partial P_2(\bar{x}_1)}{\partial \bar{x}_1^m}\right) + z_{31}^m \quad (10.10-s-1)$$

$$\frac{\partial \mathcal{L}}{\partial z_{31}^m}, i.e. = 0 = \psi(\bar{x}_1) + z_{21}^m - z_{31}^m \frac{\partial P_3(\bar{x}_1, \bar{x}_{1-1})}{\partial \bar{x}_1^m} \quad (10.11-1)$$

$$\frac{\partial \mathcal{L}}{\partial \bar{x}_1^m}, i.e. \text{ and } m \neq 1 = 0$$

$$= \psi(\bar{x}_1) + z_{21}^m - z_{31}^m + \frac{\partial P_3(\bar{x}_1, \bar{x}_{1-1}, \bar{x}_3^m)}{\partial \bar{x}_1^m}$$

$$= z_{31}^m - \frac{\partial P_3(\bar{x}_1, \bar{x}_{1-1}, \bar{x}_3^m)}{\partial \bar{x}_1^m} \quad (10.12-m-1)$$

The dual variables for any subinterval may be obtained as follows.

(i) Obtain z_1^m from equation (10.7-m)

(ii) Obtain z_{31}^m ; i.e. from equation (10.10-s-1).

(iii) For $m = 1$; z_{21}^m is obtained from equation (10.11-1), and for other values from equation 10.12-(m-1)-1.

The gradient vector is as follows.

$$\frac{\partial L}{\partial z_1^m} \text{ (for } i \neq s) = \frac{\partial f_i(z_1^m)}{\partial z_1^m} + z_1^m \left(1 - \frac{\partial r_i^m}{\partial z_1^m} \right) \quad (10.13)$$

$$\frac{\partial L}{\partial I_1^m} \text{ (for } m \neq 1) = z_{21}^m - z_{31}^m - \frac{\partial P_1}{\partial I_1^m} (Y_1^m, Y_1^{m-1}, I_1^m) \quad (10.14)$$

Next set of independent variable may be obtained by any of the first order gradient techniques described in section 7.5, 7.6 and 7.7. Conjugate gradient method is best suited. The method is expected to be stable and fast enough as is true for the first order gradient techniques. Strongest source of divergence is active penalty functions. This difficulty can be minimized by initially choosing a large value of R and small s .

This technique however has the disadvantage of large memory requirement since the independent variables and gradients need be simultaneously stored for all subintervals. Decomposition technique used by in reference 6 can be applied to this problem as follows.

10.1.3 Decomposition technique

It is assumed that the water discharge of all the subintervals and the thermal generations of all buses except that of slack are decision variables. The Lagrangian is as follows.

$$\begin{aligned}
 L = & \sum_{i=1}^m \left[\left\{ \sum_{j \in H} f_j(P_i^m) + u(P_i^m) \right\} \right. \\
 & + \sum_{j \in H} \left\{ w(Y_j^m) + v(P_j^m) \right\} + z_1^m \left(\sum_1^m P_i^m - P_L^m - P_D^m \right) \\
 & + \sum_{j \in H} z_{2j}^m (Y_j^m - Y_j^{m-1} - J_j^m + I_j^m) \\
 & \left. + \sum_{j \in H} z_{3j}^m \left\{ p_i^m - p_i(Y_j^m, Y_j^{m-1}, I_j^m) \right\} \right] \\
 & + z_{4j}^m (Y_j^m - Y_j^0 - \sum_m J_j^m + \sum_m I_j^m) \quad (10.15)
 \end{aligned}$$

where Y_i^f is the prescribed value of storage at the end of i th subinterval.

It is proved in reference 42 that for the optimal solution in the feasible domain the Lagrangian is minimum for the unknown variables and maximum for the dual variables. The decomposition technique is based on this. L is minimised for all the other variables and maximised for z_{4i} , $i \in H$. Other dual variables are obtained by the technique of section 7.3.

Partial derivatives with respect to the dependent variables yield equation 10.9, 10.10 and equation 10.12.

Partial derivatives with respect to the independent variables yield following equations.

$$\frac{\partial L}{\partial I_i^m}, i \in H = z_{21}^m - z_{31}^m \frac{\frac{\partial p_i(Y_i^m, Y_i^{m-1}, I_i^m)}{\partial I_i^m}}{+ z_{4i}} \quad (10.16-m-i)$$

$$\frac{\partial L}{\partial P_i^m}, i \in S; i \neq s = \frac{\partial f_i(P_i^m)}{\partial P_i^m} + z_i^m \left(1 - \frac{\partial P_L}{\partial P_i^m} \right) \quad (10.17-m-i)$$

$$\frac{\partial L}{\partial P_i^m}, i \in S; i \neq s = \frac{\partial f_i(P_i^m)}{\partial P_i^m} + z_i^m \left(1 - \frac{\partial P_D}{\partial P_i^m} \right) \quad (10.18-m-i)$$

Suggested algorithm is as follows.

- (1) Set $z_{2i}^1 = 0$ for each storage type hydro plant since equation associated with this dual variable is redundant because of equation associated with z_{4i} .
 - (2) Assume a value of z_{4i} for each storage type hydro plant. This value will be later on modified by gradient method to maximise L of equation (10.5).
 - (3) Set $m = 1$.
 - (4) Obtain z_{1i}^m from equation (10.9-m).
 - (5) Obtain z_{3i}^m from equation (10.10-m-1).
 - (6) For $m > 1$ obtain z_{2i}^m from equation 10.1-(n-1)-1.
 - (7) Adjust I_i^m and P_i^m , by the gradient technique with the gradients obtained from equations 10.16-m-1 and 10.17-m-1; and repeat from 4 till the gradients of the above equations satisfy the optimal conditions of equations 9.1, 9.2 and 9.3.
 - (8) If $m < n$; set $m = m + 1$ and repeat from 4.
 - (9) With the gradients obtained from the following equation; adjust z_{4i} by the gradient technique and proceed from 3.
- $$\frac{dL}{dz_{4i}} = Y_i^f - Y_i^o - \sum_m J_{1i}^m + \sum_m I_{1i}^m \quad (10.18)$$

10.1.4 Comparison

- (1) The decomposition technique needs comparatively less storage.

(ii) A quadratic function need the number of iterations equal to the number of decision variables by conjugate gradient method. Assuming the quadratic behaviour of the objective function; the number of iterations for a single value of z_{4i} for the decomposition technique will be almost equal to that without decomposition. Since L is to be optimised with respect to z_{4i} ; this set of iteration will have to be repeated two to three times the number of storage type hydro-plants.

10.2 Short Term Planning

This involves planning for a comparatively short period say one day. During this period, total amount of water K_1 to be used for each reservoir has been preestimated. Head variation could be ignored. The period is again subdivided into 4 (say 24) subintervals, during which the load demand may be assumed constant. Assuming that the hydrogeneration could be expressed as follows

$$P_i^m = p_i(I_i^m) \quad i \in H \quad (10.19)$$

For a simplified model this could be expressed as follows.

$$P_i^m = H_i(I_i^m - q_i)/T \quad (10.20)$$

If losses are expressed in terms of loss formula the analysis is similar to that of section 10.1. If busbar powers are expressed in terms of bus bar voltage magnitudes and angles, the Lagrangian is as follows.

$$\begin{aligned}
 L = & \sum_i \left[\sum_{j \in S} r_j (P_j^m) + \sum_j \lambda_j^m \{ p_j (v^m, \theta^m) - P_j^m + z_j^L \} \right] \\
 & + \sum_j \mu_j^m \{ \gamma_j (v^m, \theta^m) - \lambda_j^m + z_j^L \} \\
 & + \sum_{j \in H} \pi_j^m (v_j^m - \bar{v}_j) + \sum_{j \in R} \pi_j^m (v_j^m - \underline{v}_j) \\
 & + \sum_j (v_j^m) + \alpha(P_S^m) + \sum_{j \in H} \{ U(P_j^m) + z_j^L \} \\
 & + \sum_{j \in H} z_{3j}^L [r_j^m - p_j(I_j^m)] \\
 & + \sum_{j \in H} z_{4j}^L (K_j - \sum_m I_j^m) \quad (10.21)
 \end{aligned}$$

The variables have been classified similar to the ones described in section 9.2 and 10.1.1.

The dual variables λ^m , μ^m , π^m during any subintervals are determined by the method of section 9.4. Other equations to obtain dual variables are as follows.

$$\frac{\partial L}{\partial P_1^m}, \quad i \in H = 0 = -\lambda_1^m + W'(P_1^m) + z_{31}^m \quad (10.22)$$

$$\frac{\partial L}{\partial I_1^1}, \quad i \in H = 0 = -z_{41} - z_{31}^1 \frac{\frac{\partial p_1(I_1^1)}{\partial I_1^1}}{\frac{\partial p_1(I_1^1)}{\partial I_1^1}} + v'(I_1^1) \quad (10.23)$$

z_{31}^m may be obtained from equation 10.22 and z_{41} from 10.23.

Independent variables can be adjusted by the gradient techniques with gradients obtained as follows.

$$\frac{\partial L}{\partial z_{41}^m}; \text{ i.e., } 1 \neq s = \frac{\partial f_1(P_1^m)}{\partial P_1^m} - \lambda_1^m \quad (10.24)$$

$$\frac{\partial L}{\partial z_{31}^m}; \text{ i.e., } 1 \neq s = -z_{41}^m - z_{31}^m - \frac{\partial p_1(I_1^m)}{\partial I_1^m} \quad (10.25)$$

$$\frac{\partial L}{\partial z_{11}^m}; \text{ i.e., } 1 \neq s = -\mu_1^m \quad (10.26)$$

where set A is defined in section 9.4.

Gradient for transformer tap settings is given by equation 9.29.

10.2.1 Decomposition technique

With the advantages and disadvantages mentioned in section 10.1.4, the decomposition technique for the above problem may be as follows.

The Lagrangian is minimised with respect to the unknown variables and maximised with respect to the dual variables z_{41} i.e.

$$\frac{\partial L}{\partial z_{41}} = \kappa_1 - \sum_m I_1^m \quad (10.27)$$

$\lambda(I_1^m)$ is excluded from the Lagrangian of equation 10.21. The value of z_{31}^m obtained from equation 10.22 is used to obtain the gradient vector during the subinterval m in equations 10.24, 10.25 (including for m = 1), 10.26 and 9.29. The independent variables are adjusted during the subinterval to make the gradients satisfy the set of equations 9.1, 9.2 and 9.3.

Technique of this section is parallel to that of reference 6. However the decision variables and the Lagrangian is different.

10.3 Example

The problem considered has two thermal generators with the following incremental costs,

$$\frac{d f_1}{d P_1} = .4 + .2 P_1$$

$$\frac{d f_2}{d P_2} = .6 + .3 P_2$$

and two hydro generators represented by equation 10.4 with following data

$$H_1 = 1.0, C_1 = .1, q_1 = .1$$

$$H_2 = 1.2, C_2 = .12, q_2 = .05$$

Total water available in the 12 subintervals of one unit time each is 12 and 15 respectively. Initial storages are 10 and 12 respectively. The loss formula matrix is as follows.

| | | | |
|-------|-------|--------|--------|
| 0.05 | -0.02 | -0.01 | 0.0 |
| -0.02 | 0.06 | -0.02 | -0.01 |
| -0.01 | -0.02 | +0.04 | -0.005 |
| 0.0 | -0.01 | -0.005 | +0.02 |

Water input during the 12 subintervals is as follows.

| m | 1 | 2 | 3 | 4 | 5 | 6 | 7 | 8 | 9 | 10 | 11 | 12 |
|-------------------|---|----|-----|-----|-----|-----|-----|------|-----|-----|----|----|
| J_1^{in} | 0 | .6 | 1.2 | 1.2 | 1.2 | 1.6 | 2.4 | 1.44 | 1.2 | .36 | 0 | 0 |
| J_2^{in} | 0 | 0 | 0 | 1.5 | 3.0 | 4.5 | 4.5 | 1.5 | 0 | 0 | 0 | 0 |

Initial and final values are given in Table 10.3.1 and 10.3.2 respectively.

10.4 Comments

Application of discrete maximum principle essentially will give the same set of equations since the relations of this principle have been obtained in reference 25 making use of Lagrangian formulation. Application of discrete maximum principle as attempted in reference 29 turns out to be the approach of 10.1.3 except the dual variable associated with the storage equality expression is not adjusted by gradient technique, for a system with one hydro and one thermal generation. For a system of multihydro-plants this reference (29) suggests an iterative technique of Gauss Seidal type, convergence of which is doubtful.

Analysis given by Kirchmayer⁵² also makes use of a technique similar to decomposition technique except no definite method is given to adjust the dual variable of the storage expression.

The approach of section 10.2 making use of a.c. simulation of network does not appear to be very promising

because of the need for repeated load flow solution during each subinterval of each iteration of scheduling. The author feels that the proposed methods of section 10.1 i which the losses are expressed in terms of loss formulae are more practical for short term planning too. It might be better to make use of different loss formula matrix for light and heavy loads. However, at the time of actual dispatch the thermal generation, reactive power and tap settings may be obtained making use of the computer programme of Chapter 9, with the hydrogenerations as planned.

Table 10.2.1
Initial value

| Subinterval number | $\frac{I_1}{I_2}$ | Water discharge | Hydro Power | | Internal power |
|-----------------------|-------------------|--------------------|-------------|--------|----------------|
| | | | 1 | 2 | |
| 1 | 1.00 | 1.25 | 1.7549 | 3.4655 | 1.6169 |
| 2 | 1.00 | 1.25 | 1.6813 | 3.1895 | 1.6169 |
| 3 | 1.00 | 1.25 | 1.6623 | 3.0735 | 1.6169 |
| 4 | 1.00 | 1.25 | 1.7039 | 3.0771 | 1.6169 |
| 5 | 1.00 | 1.25 | 1.7189 | 3.0591 | 1.6169 |
| 6 | 1.00 | 1.25 | 1.7039 | 3.0713 | 1.6169 |
| 7 | 1.00 | 1.25 | 1.6729 | 3.0535 | 1.6169 |
| 8 | 1.00 | 1.25 | 1.9457 | 4.3559 | 1.6169 |
| 9 | 1.00 | 1.25 | 1.9765 | 4.3295 | 1.6169 |
| 10 | 1.00 | 1.25 | 1.9817 | 4.0535 | 1.6169 |
| 11 | 1.00 | 1.25 | 1.9947 | 3.3375 | 1.1105 |
| 12 | 1.00 | 1.25 | 1.6649 | 3.6215 | 1.7403 |

Total cost of thermal generation = 14.8 pu.

Table 10.3..Final value

| Subinterval number | Water discharge l_1^m | Water discharge l_2^m | Years later | | Interest factor | |
|-----------------------|----------------------------|----------------------------|-------------|--------|-----------------|--------|
| | | | 1 | 2 | 1 | 2 |
| 1 | 0.7801 | 0.7262 | 1.3336 | 1.3647 | 1.0000 | 1.0226 |
| 2 | 0.8754 | 1.0870 | 1.0051 | 1.0415 | 1.3195 | 1.3754 |
| 3 | 0.9613 | 1.4496 | 1.6267 | 2.7936 | 1.5612 | 1.5597 |
| 4 | 0.9589 | 1.1372 | 1.0008 | 2.7003 | 1.1433 | 1.1363 |
| 5 | 1.0397 | 1.2287 | 1.6140 | 3.2676 | 1.3373 | 1.3867 |
| 6 | 1.0420 | 1.3559 | 1.6618 | 3.2940 | 1.7323 | 1.2421 |
| 7 | 0.9936 | 1.2794 | 1.6628 | 1.2291 | 6.0746 | 1.1171 |
| 8 | 0.9897 | 1.2706 | 1.9373 | 1.5023 | 5.5e18 | 1.35e2 |
| 9 | 1.0040 | 1.2904 | 1.9176 | 1.6805 | 5.0455 | 1.0450 |
| 10 | 1.0475 | 1.3240 | 2.0556 | 1.3619 | 5.0369 | 1.1165 |
| 11 | 1.0510 | 1.4643 | 2.0545 | 1.5756 | 5.0260 | 1.1465 |
| 12 | 1.1762 | 1.5163 | 2.0551 | 1.6400 | 5.0173 | 1.1865 |

Total cost of thermal generation = 1.7 ru.

Chapter XI

CONCLUSIONS AND SUGGESTIONS FOR FURTHER WORK

The author is very strongly convinced about the need of sparsity programming for systems involving large sparse matrices, absence of which, apart from the huge additional computer time, may mean failure of an algorithm due to built up round off errors. Piecewise methods have been used by many authors for large network problems such as Z-matrix load flows^{1,23,24} and optimal power system operations^{4,52}. The power system is subdivided into a number of interconnected subsystems. Z-matrix or other now sparse matrix of each subsystem is formed separately and every subsystem is solved independently, making use of the variables obtained for other subsystems. Any such techniques need Gauss-Seidal type block iteration, convergence of which may be hard in the absence of efficient subdivision. If Z-matrix is replaced by the table of factors of nodal admittance matrix, with compact storage and optimal renumbering, decomposition in many cases may not be needed.

Application of compact storage, makes the computer logic complicated. If a compiler could be written such that the programme written in conventional way could with little modification use compact storage, this will be of great value.²
BPA probably has such a programme.²

Efficiency of an iterative technique is generally judged by actually trying the technique over the system. This usually needs months of programming effort. Through the application of fixed point analysis, it is sometimes possible to judge the algorithm. The author has demonstrated the application of this analysis for load flow methods and some of the optimal power flow methods. For the optimal power flow methods making Gauss-Seidel type iterations with Carpenter's Lagrangian, such analysis is restricted for real powers only. Additional work needs be done in this direction.

The author's computer programme for load flow by Newton's method makes use of all sparsity techniques and obtains the non zero elements of both rows and active and reactive powers in one set of calculations. The method suggested in section 6.1 for changing bus bar types works very well.

The optimal power flow programme described in Chapter IX works efficiently with hardly any convergence difficulties for any objective function and any assumed set of initial conditions. The solution needs about 20 to 30 solutions of Jacobian equations for objective functions without penalty functions and 30 to 50 such solutions for objective functions with penalty functions. As pointed earlier, apart from the routine fuel cost minimization, the method is found to work well for planning addition ~~over~~ requirements as well. Further improvements are likely as more experience is obtained while working over a practical system. On-line application

of the computer programme will need the system data telemetered and processed through 'state estimation' for which additional work is needed.

Because of computer memory limitations, our study was restricted to 30 buses only. The solution needs about 4,500 words of storage for the variables and 10,200 for the programme with two links and 13 local subroutines. While memory requirement for the former will be approximately proportionate to the bus bar numbers, for the later it will remain unchanged. The author expects to be able to run a problem of about 150 buses on an IBM 1130 computer with 32 K words of storage. Additional work needs be done for simulating larger systems on small computers. Decomposition does not appear to be very useful along with optimal renumbering and compact storage.

In Chapter X algorithms for optimal hydrothermal scheduling with and without, a.c. simulation of the network, and decomposition for multithermal and multihydro plants (with and without head variation) are presented for deterministic load demands and water inflows. Feasibility studies need be made for the algorithms with the trade offs of computer time, memory and accuracy. Additional work needs be done for stochastic load demands and water inflow.

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