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# MODIFIED ELECTRIC POWER SYSTEM STATE ESTIMATION -

# **MULTI- PROCESSING TECHNIQUE**

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### **ABSTRACT**

Estimating the state of the power system in Real time is a challenging task, mainly because, solving the non-linear large power system equations by iterative solution comsumes huge computational time and memory. Online Energy Management System (EMS) these state estimators need to Estimate the state of the system at very short intervals. Conventional methods like Newton-Raphson is not suitable for such real time application.

To improve the performance of SE, certain techniques like AI, Real time, Hirarchial and Dynamic methods are listed in reference [1][4][5][11][12][13]. The most common "Two-Level" SE technique is applied by spliting the large power system in to sub-networks and each sub-networks are managed by the local control station in coordination with the Centre station – [2][3][7][8][9][14]. Even though it reduces the computational time, still it is not suitable for real time applications. In the "Two-Level SE" mainly same NR method is used, however, before applying the NR technique, the electric network is physically divided, which results to mathematical approximation.

This paper presents an approach to solve the problem in hand without physically dividing the network and/or any approximation on the existing NR method. Single large mathematical problem can be divided into smaller independent parallel task by changing the multiplication method of large sparce jacobian matrix. This leads to a whole new approach resulting in reducing in computing time as well as dynamic memory requirement.

**KEYWORDS: SE**-State Estimation, **WLS**-Weight Least Square, **RTU**-Remote Terminal Unit, **NR**-Newton Rapson, **AI**-Artificial Intelligence, **Node Area**-A Node along with its Connected Node is Referred as Node Area

#### INTRODUCTION

The power system is highly distributed in nature, Instead of dividing the network in to sub networks, it is better to divide the large single problem into small independent sub tasks so that multiple processor can be used for computation. Major challenge here to obtaing the independent sub-task. The NR method mathematical solution is procedure oriented iterative technique which involvs large number of non-linear simultaneous equations. The ideal solution would have been the vertical division of the probem into many sub-independent parallel task. By modifying the existing Newton-Raphson state estimation computational procedure requires the clear understanding of the procedural steps involed in the existing NR method. The brief insight of NR method is given below.

### NEWTON RAPSON (WLS) METHOD FOR STATE ESTIMATION –BRIEF REVIEW

The Newton Rapson final equation is  $A \Delta x = b$  (taking weighting matrix W which is a diagonal matrix  $W_{ii}$  is the standard deviation of meter i.) By applying the tylor series to the nonlinear equations of power system following equations

are derived. Here, the error vector includes nonlinear vector function f(x). In order to estimate x, an initial value  $x_0$  is assumed and after every iteration  $x_0$  will be up-dated till the convergence.

$$(J_0^T W J) \Delta x = J_0^T W \Delta z$$
Let  $A = (J_0^T W J) \& b = J_0^T W \Delta z$ 

$$A \Delta x = b$$
(1)

This is a set of linear equations, if higher order terms of the taylor expansion of f(x) were really negligible, the solution yield the correct 'x'. The jacobian J is itself a function of x. The state variable vector x can be obtained by solving the equation  $A*\Delta x = b$  iteratively. The vector x should therefore be changed accordingly after every iteration till the convergence is obtained.

 $x^{n+1} = x^n + \Delta x^1$  (1<sup>st</sup> iteration count) until convergence is reached. Weighted least squares (WLS) minimization technique is used.

No of state variables = (2\*n-1)

Total no of measurements = m

These measurements may include one or all quantities such as  $[P_i, Q_i, p_{ij}, q_{ij}, V_i, \delta_i]$ 

- P<sub>i</sub>,Q<sub>i</sub> = Real & Imaginary part of injected power respectively
- $p_{ij}$ ,  $q_{ij}$  = Real & imaginary part of line follows respectively
- $V_i \& \delta_i = Voltage magnitude \& phase angles.$
- Dimension of Jacobian matrix = m \* (2n 1)

$$\Delta z = z^{measured} - z^{calculated}$$

'W' is the diagonal weigh matrix of the order of (m\*m)

$$x = [\delta_1, \delta_2, \dots, \delta_{n-1}, v_1, v_2, \dots, v_n]$$
; state variables.

n= number of network nodes, 1,2,...n.

Elements of jacobian are derived from injected power and line flow equations.

$$\begin{bmatrix} \Delta P_{i} \\ \Delta Q_{i} \\ \Delta p_{ij} \\ \Delta q_{ij} \\ \Delta v_{i} \\ \Delta \partial_{i} \end{bmatrix} = \begin{bmatrix} |H_{1}| & |H_{2}| \\ |H_{3}| & |H_{4}| \\ |H_{5}| & |H_{6}| \\ |H_{7}| & |H_{8}| \\ |\Delta v_{i} \\ |1 & |0 & | \end{bmatrix} \begin{bmatrix} \Delta \delta_{i} \\ \Delta v_{i} \end{bmatrix}$$

$$(2)$$

### PROPOSED ALGORITHUM

### **Proposed New Algorithum**

The following modification of forming the jacobian results in independent parallel task and thereby multiple processors can be used to carry out the task parallelly. Jacobian being the large sparse matrix, the suggested new method can fully avoid storing and retrieving of large saprse matrix. the let 'm' be the total number of measurements taken for the given network. Each row of jacobian is formed based on the type of measurement underconsideration, if there are 'm' measurements, the jacobian of the network will contain 'm' rows and (2n-1) coloms.

Instead of forming the complete jacobian of the netork and then multiplying to find 'A' matrix, (see eq1), It can be computed by each row of jacobean  $[J_1 * W_1 * J_1 = A_1]$  and  $[J_1 * W_1 * Z_1 = b_1]$ , this results in first terms of each elements of the resultant matrix "A" and "b" respectively, which can be visualized as shown below.

$$\{A_1 + A_2 ... + A_i ... + A_m\} \Delta x = \{b_1 ... + b_r ... + b_m\}$$
(3)

$$\sum_{j=1}^{m} (A_j) \Delta x = \sum_{j=1}^{m} (b_j)$$
(4)

$$A = \{A_1 + A_2 + A_r + A_m\}$$
 (5)

$$b = \{b_1 + b_r + b_m\}$$
 (6)

Hence, it is not necessary to form complete two dimentonal Jacobian and then carry-out the multiplication task. New jacobian is a one dimention arry of the order of [1\*(2n-1)] is replaced by [m\*(2n-1)], thereby no special sparse matrics storage is required.

### Node Wise Grouping of Measurement: Jacobian Relation

The measurements can be grouped node wise, and it can be represented as shown below.

$$Node-1 \ \textit{Measurements} \begin{bmatrix} \left\{ \begin{matrix} \Delta P_{1} \\ \Delta Q_{1} \\ \Delta \rho_{ij}^{1} \\ \Delta \alpha_{1}^{1} \\ \Delta \delta_{1} \end{matrix} \right\} \end{bmatrix} = \begin{bmatrix} \left[ \begin{matrix} |H_{1}^{1}| & |H_{2}^{1}| \\ |H_{3}^{1}| & |H_{4}^{1}| \\ |H_{5}^{1}| & |H_{6}^{1}| \\ |H_{7}^{1}| & |H_{8}^{1}| \\ |H_{9}^{1}| & |H_{10}^{1}| \\ |H_{11}^{1}| & |H_{12}^{1}| \end{bmatrix} \end{bmatrix} \begin{bmatrix} \Delta \delta_{i} \\ \Delta \nu_{i} \end{bmatrix}$$

$$\begin{bmatrix} \Delta \delta_{i} \\ \Delta \nu_{i} \end{bmatrix}$$

$$Node-r \ measurements \begin{bmatrix} \Delta P_r \\ \Delta Q_r \\ \Delta Q_t' \\ \Delta q_{ij}^r \\ \Delta v_r \\ \Delta \frac{\partial}{\partial r_r} \\ \end{bmatrix} = \begin{bmatrix} |H_1^r| & |H_2^r| \\ |H_3^r| & |H_4^r| \\ |H_5^r| & |H_6^r| \\ |H_7^r| & |H_8^r| \\ |H_9^r| & |H_{10}^r| \\ |H_{11}^r| & |H_{12}^r| \end{bmatrix} \begin{bmatrix} \Delta \delta_i \\ \Delta v_i \end{bmatrix}$$

$$\begin{bmatrix} \Delta \delta_i \\ \Delta v_i \end{bmatrix}$$

$$Node - n^{th} \ measurements \begin{bmatrix} \Delta P_{n} \\ \Delta Q_{n} \\ \Delta p_{ij}^{n} \\ \Delta q_{ij}^{n} \\ \Delta v_{n} \\ \Delta \partial_{n} \end{bmatrix} = \begin{bmatrix} |H_{1}^{n}| & |H_{2}^{n}| \\ |H_{3}^{n}| & |H_{4}^{n}| \\ |H_{5}^{n}| & |H_{6}^{n}| \\ |H_{7}^{n}| & |H_{8}^{n}| \\ |H_{7}^{n}| & |H_{10}^{n}| \\ |H_{11}^{n}| & |H_{12}^{n}| \end{bmatrix} \begin{bmatrix} \Delta \delta_{i} \\ \Delta v_{i} \end{bmatrix}$$

$$\begin{bmatrix} \Delta \delta_{i} \\ \Delta v_{i} \end{bmatrix}$$

$$A_{NAr} = \left\{ \left( A_r^P + A_r^Q + \sum A_r^{pij} + \sum A_r^{qij} + A_r^V + A_r^{\delta} \right) \right\}$$

$$= J_{NAr}^T * W_r * J_{NAr}$$

$$(10)$$

$$b_{NAr} = \left\{ \left( b_r^P + b_r^Q + \sum b_r^{pij} + \sum b_r^{qij} + b_r^v + b_r^{\delta} \right) \right\}$$

$$= J_{NAr}^T * W_r * z_{NAr}$$
(11)

$$A = (A_{NA1} + A_{NA2} + ... + A_{NAr} + ... + A_{NAn}) = \sum_{j=1}^{n} (A_{NAj})$$

$$b = (b_{NA1} + b_{NA2} + ... + b_{NAr} + ... + b_{NAn}) = \sum_{j=1}^{n} (b_{NAj})$$

$$\sum_{j=1}^{n} (A_{NAj}) \Delta x = \sum_{j=1}^{n} (b_{NAj})$$
(12)

$$\sum_{j=1}^{m} (A_j) = \sum_{j=1}^{n} (A_{NAj}) = A$$
(13)

$$\sum_{i=1}^{m} (b_{j}) = \sum_{i=1}^{n} (b_{NAj}) = b$$
(14)

#### Hence $A\Delta x = b$

 $\left( \right)_{NAi}$  – subscript 'NAi' refers to ith node area

 $A_r^P$  - It is the sub set of matrix 'A' for the injected Real power measurement taken at 'r'th node/bus. Similarly,

 $A_r^Q$  (here Q represents reactive power) and like wise for other measurements.

Hence, from the above derivation, it is evedient that up to the formation of  ${}^{\circ}A_{NAi}{}^{\circ}$  (node level), computations can be carried-out independently without any approximation, which gives the scope for parallel processing.

### **System Design**

### **Assumptions**

- All measurements are taken at the same instant
- At each node total of measurement can be: injected power measurements of node + (node voltage measurement + node angle measurement + the line flow measurements) of that node to other connected nodes.

#### **Flow Chart**

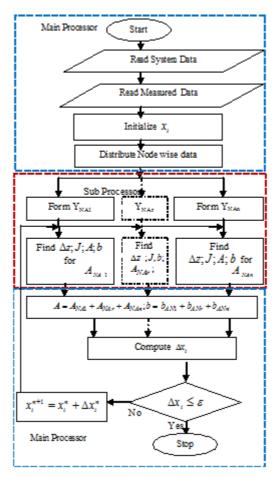


Figure 1

#### COMPUTING TIME ANALYSIS OF SE

# **General Technique**

The computational time of state estimation using NR method is directly linked with the number of nodes in the network. Generally, to solve  $A*\Delta x = b$ , if the computational time required is (say) " $t_1$ " for the network having  $n_1$  number of nodes, then the computational time required for the network having  $n_2$  number of nodes is approximately equals to  $(n_2/n_1)^2*t_1$ . Or in other words the SE computational time increases "square time" the ratio of increase in the number of nodes of the network. It's because as the number of node increases the size of the Jacobian and "A" matrix. For example

The size of matrix "A" for the network having 10 nodes is nearly equal to (20 x 20) = 400.

The size of matrix "A" for the network having 30 nodes is nearly equal to  $(60 \times 60) = 3600$ .

$$(3600/400) = 9 = (30/10)^2$$
.

#### **Optimum Computing Technique**

The fact is, these matrices are highly sparse in nature. These none zero elements of jacobian/'A" has a direct relation with the network incident matrix. Using this information one can design the computation focusing only on none zero elements to avoid operations which results in zero. This can be done by using the network incident matrix as index to compute arithmetic operation of none-zero elements. This makes the computational time required is directly proportional to the number of measurements and does not depend on the size of the matrices. SE computational time reduces drastically by using optimum computing technique.

For example, a 10 bus system having number of lines connected (say) = 14, then the size of Ybus = 10x10 = 100, whereas none zero elements in the Ybus = 10+14x2 = 38;

Hence, none zero elements in the matrix "A" = 38 x4=152; whereas size of "A" approximately equal to (20x20) =400. Say "t<sub>1</sub>" is SE computational time for the network having 10 nodes using general method, then by using optimum technique, SE computational time for the same network is approximately equal to (152/400)x t<sub>1</sub>, let this be equal to t<sub>2</sub>. It's obvious that t<sub>2</sub> << t<sub>1</sub>. Now if the network size is increased to n<sub>2</sub> (say for example n<sub>2</sub>=30, number of connected lines=41, none zero Ybus elements=30+41x2=112 hence number of none zero elements of matrices=112x4=448), then the computing time is approximately proportional to (448/152) = which is nearly equal '3', hence the computing time is nearly equal to=  $(n_2/n_1)$  x t<sub>2</sub>. Because, whatever the size of matrices, computational time depends upon the number of none zero elements which increases almost linearly with node.

Total SE computational time can be related with the following equation

$$(SE)_{t} = \left(\Delta x i = A^{-1} * b\right)_{t} + \left(J^{T} * J\right)_{t} + \left(\varepsilon\right)_{t}$$

$$(SE)_{t} = \text{total SE computational time}$$

$$\left(\Delta x i = A^{-1} * b\right)_{t} = \text{Time taken to solve } \Delta x i$$

$$\left(J^{T} * J\right)_{t} = \text{Time taken to } obtain A \& b$$

$$\left(\varepsilon\right)_{t} = \text{updating time}$$

## **EXAMPLE & RESULTS**

Note: All quantities are in pu.

Table 1: Input Line Data for 13 Bus System (ISE)

Bus No	R(Pu)	X(Pu)
1-2	.00147967	.00286760
2-3	.00043843	.00124174
3-4	.00027711	.00078428
4-5	.00059760	.00166769
4-8	.00159967	.00310017
5-6	.00034314	.00097190
5-9	.00034314	.00097190
6-7	.00032364	.00091669
7-10	.00032364	.00091669
8-9	.00029438	.00083380
9-10	.00053157	.00150562
9-12	.00037793	.00107050
10-11	.00058777	.00166488
11-13	.00032364	.00091669
12-13	.00036843	.00104355

**Table 2: Injected Power & Voltage Measurements** 

Ems No.	Real	1mg.	Bus No.	Volt
1	28.5335	13.0837	1.	1.05300
2	-4.73	-1.55	3.	0.95792
3	-1.27	-0.41	10.	0.92304
4	-0.35	-0.11	11.	0.91939
5	-4.38	-1.44	12.	0.92178
6	-2.11	-0.69	13.	0.92036
7	-0,42	-0.13	0	0
8	-4.73	-1.55	0	0
9	-1.27	-0.41	0	0
10	-0.35	-0.11	0	0
11	-4.38	-1.44	0	0
12	-2.11	-0.69	0	0
13	-0.42	-0.13	0	0

**Table 3: Line Flow Measurements** 

Bus No.	Real	1mg.
1-2	28.5327	13.0837
2-1	-27.2179	-10.5355
2-3	22.4879	8.9855
3-2	-22.2197	-8.2258
3-4	20.9000	7.7700
4-5	12.9700	5.1000
4-8	7.6200	2.1282
5-9	4.3979	1.9040
8-4	-7.4182	-1.9161
8-9	2.7900	0.3800
9-10	1.,8043	0.5194
11-13	-1.4600	-0.4582

Table 4: SE Result of L3 Bus Test System \*Slack Bus

No	Volt	Ang Rad
1	1.053269	0.101912
2	0.979368	0.041320
3	0.958231	0.015757
4*	0.945912	0.000000
5	0.928889	0.021096
6	0,925904	0.025075
7	0.924513	-0.026839
8	0.926468	0.022792
9	0.925278	0.025320
10	0.923398	0.028199
11	0.919737	0.033261
12	0.922133	0.029788
13	0.920707	0.031856

<sup>\*</sup> Convergence = .001 for voltage,

Table 5: Computational Time for 13 Bus Test System

	Total SE Time	Time-up to Jacobian Product
Conventional method	1.08 ms	0.88 ms
New method (single processore)	0.52 ms	0.32 ms
New method (paralle processores)*	0.225 ms	0.025 ms

<sup>\*</sup>Number of parallel processors are assumed to be =13

<sup>= .00001</sup> for angle.

Table	٠6٠	$\mathbf{D}^{A}$	ativa	Time
1 and	. v.	KCI	auve	1 11111

	Total SE Time Ratio	Time Ratio- Jacobian Product
Conventional /new single ≈	2	2.75
Conventional /new parallel ≈	5	35
New single/New parallel ≈	2.36	13

The above table shows actual results obtained which is same as that of the Integrated SE results obtained by the conventional method. It shows that both approach yields the same results. Thereby this new concept has been proved with this example.

### **CONCLUSIONS**

Node wise grouping of data leads to new approach for applying the NR solution. The number of state variables to be computed at each bus-level is very small as compared to the whole network, and till the last step computations can be carriedout independently & separately thereby the fast computation is possible. If sufficient numbers of measurements are ensured at each node, then state estimation can be carried out at each bus. At Central computing station only the final estimated state variables are to be considered as sudeo-measurements. The other possibilities of this approach is under progress.

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