SANTIAGO PEÑATE VERA

PRACTICAL GRID MODELLING

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Contents

1	Introduction 9	
2	Notation 11	
3	General network model 13	
4	Magnitudes and units 15	
5	The branch element 21	
6	The bus and it's connected elements	29
7	Topology analysis and consolidation	31
8	Power flow 33	
9	Time series power flow 43	
10	Stochastic power flow 45	
11	State estimation 47	
12	Short-circuit 49	
	Bibliography 51	

List of Figures

3.1	Graph with 8 nodes and 7 branches. The graph contains two islands.	13
3.2	Two node system in 3 phases, split into three systems of two nodes	
	each. 14	
4.1	Voltage delay. 15	
4.2	Voltage delay in phasor representation in the complex plane. 15	
4.3	Delta connection scheme. 16	
4.4	Wye connection scheme. 16	
5.1	General branch model. 21	
5.2	Conversion from the Delta branch model to the Y branch model.	7
6.1	YISV Bus model. 29	
8.1	Circuit's admittance matrix and vectors reduction. Representation	
	for a six node circuit with two slack nodes. 34	
8.2	Jacobian based linear system to compute the voltage increments. 36	5

List of Tables

4.1	Electrical magnitudes and their units. 15	
4.2	Magnitudes and their real and imaginary complex components.	15
4.3	Electrical magnitudes and their per unit base. 17	
_	Equations of the generalized Π model. 21 Three-phase transformer impedances of the branch model. 24	
8.1	Bus types. 33	

1 Introduction

This book aims at explaining grid modelling from a practical perspective, providing state of the art models, algorithms as well as implementation hints and examples.

The goal is to provide a reference document for researchers and computer engineers in the field of electric systems modelling, so that the task of building your own simulator or extend the already available open source ones becomes feasible.

The book assumes that the reader has a basic understanding of matrices, vectors and complex numbers. The notation used in the book is intended to be clear and accessible as much as it is practically possible.

Most of the formulas presented will imply vectorized operations such as matrix multiplications or element-wise multiplications. This notation makes it much easier to implement the numerical methods, and in case of scripting languages such as python, enables the use of fast linear algebra libraries. This is key in the implementation of electrical systems simulators, since most articles and books provide the formulas using summations, which implementation in languages other than c or Fortran would make the simulator prohibitively slow in comparison with the vectorized version.

Vectorized formulas also allow a much simpler debug and maintenance of the produced software.

2 Notation

- $A \times B \rightarrow$ Matrix-Matrix dot product.
- $A \times b \rightarrow$ Matrix-Vector dot product.
- $a \times b^{\top} \rightarrow$ vector-vector dot product. The result is a number.
- A · B → Matrix-Matrix element wise multiplication. A and B must have the same dimensions.
 The result is a matrix.
- *a* · *b* → vector-vector element wise multiplication. The result is a vector.
- A* → Element-wise conjugate of the numbers composing A.
- $A^{-1} \rightarrow \text{Matrix inverse.}$
- $A^T \rightarrow$ Transposed matrix or vector.
- A⁻¹ × b → Solve the linear system where A is the coefficients matrix and b is the free terms vector. Never perform the inverse of A and then multiply it by b. Instead use a linear system
- A[rows,:] → From the matrix A, pick the rows which indices are contained in the vector rows.

- A[:,cols] → From the matrix A, pick the columns which indices are contained in the vector cols.
- A[rows, cols] → From the matrix A, pick the rows which indices are contained in the vector rows and the columns which indices are contained in the vector cols.
- b[rows] → From the vector b, pick the elements contained in the vector or list rows.
- diag(b) → Convert the vector b into a diagonal matrix.
- diag(A) → Extract the diagonal of the matrix A as a vector.
- $real(A) \rightarrow Extract$ the real part of A.
- $imag(A) \rightarrow Extract$ the imaginary part of A.
- $max(b) \rightarrow Maximum$ value of the vector b.
- $max(c,d) \rightarrow Pick$ the greater value between c and d.

The electrical grid can be assimilated to a graph. A graph is a mathematical object composed of nodes and edges (or branches). From a calculation point of view, a node is the place where the voltage is calculated given power and current injections, and a branch is the place where the current and power that flows through it are computed given the nodes voltage.

In the general network model, the neutral wire and the earth "wire" are embedded into the three-phase equivalent using Kron's reduction. See ¹ for a comprehensive explanation or simply ² for practical application.

A node has power injection or consumption devices attached to it. Examples are loads, generators, capacitor banks or any other device that injects or consumes power from the grid.

A branch might have devices attached that modify the flow through the branch. Such devices are known as FACTS (Flexible Alternating Current Transmission Systems).

A real life grid can be composed of several isolated groups of nodes (islands). It is impossible to perform calculations of several islands at once in the same numerical process. Hence, the maximal calculation object is the island circuit: a graph that does not contain further islands inside. In practice, a grid is split in its islands, each island is computed separately and the results are merged back to provide a consistent analysis interface of the whole grid.

¹ Florian Dorfler and Francesco Bullo. Kron reduction of graphs with applications to electrical networks. *IEEE Transactions on Circuits and Systems I: Regular Papers*, 60(1):150–163, 2013 ² William H Kersting. *Distribution system modeling and analysis*. CRC press, 2012

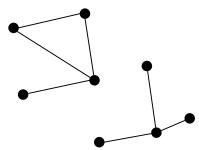


Figure 3.1: Graph with 8 nodes and 7 branches. The graph contains two islands

n-phase modelling

The electrical grid calculations are mostly done in what is called *positive sequence equivalent* (a single phase equivalent of a three-phase grid) In practice the positive sequence equivalent is applied to branches with one, two or three phases, this simplifies the modelling, but makes the realistic analysis of the grid much harder since the phases imbalance has been neglected.

In this book, three-phase models will be presented, but the calculations will be done on phase-by-phase basis. This is possible using the models presented by Vieira, Freitas and Morelato ³. In their work, the authors pick the diagonal of the elements admittance matrices to form single-phase grids, that are simulated independently. The magnetic coupling effects are included in the single-phase admittance matrices as shunt elements. This allows a very flexible

³ JCM Vieira, W Freitas, and A Morelato. Phase-decoupled method for three-phase power-flow analysis of unbalanced distribution systems. *IEE Proceedings-Generation, Transmission and Distribution*, 151(5):568–574, 2004

3.1

model of the grid while retaining the calculation accuracy of a full-blown n-phase admittance matrix.

So, because of the latter, the models will be introduced in threephases and whenever needed in positive sequence equivalents, but the numerical methods will consider only one phase at the time. In practice we will solve one numerical problem per phase. To simulate n phases, we will perform n numerical simulations, merging the results back into n-phase structures.

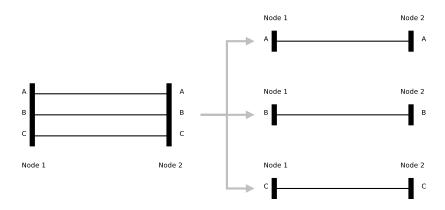


Figure 3.2: Two node system in 3 phases, split into three systems of two nodes each.

Further details about how to decouple the three-phase impedance matrices into three single phase impedances are presented in the section 5.4.

4 Magnitudes and units

As the electric energy is mostly distributed in alternating current, electrical magnitudes are waves that vary their polarity (positive and negative value) and amplitude in time. because of this, the electrical magnitudes are expressed by complex numbers to denote the position of the value in the two-dimensional plane amplitude-time.

The units in electrical grid modelling are represented in the tables 4.1 and 4.2.

Magnitude	Unit	Recommended user input unit
Voltage	V (Volt)	kV (kilo-Volt)
Current	A (Ampere)	kA (kilo-Ampere)
Power	VA (Volt-Ampere)	MVA (Mega-Volt-Ampere)
Active power	W (Watt)	MW (Mega-Watt)
Reactive power	VAr (Volt-Ampere-reactive)	MVAr (Mega-Volt-Ampere-reactive)
Impedance	Ω (Ohm)	Ω (Ohm) or per-unit
Admittance	S (Siemens)	S (Siemens) or per-unit

The figure 4.1 shows two voltage waves. The one represented with a dotted line is delayed an angle δ with respect to the reference voltage wave represented by the plain black line. Since both waves are periodical, both can be represented as "phasors" or vectors indicating the magnitude's value and angle in the complex rectangular plane as depicted in the figure 4.2. Figures 4.1 and 4.2 are equivalent representations.

Magnitude	Real part	Imaginary part	Relation
S (Power) V (Voltage) Expressed as	P (Active power) V_r (Real voltage)	Q (Reactive power) V_i (Imaginary voltage)	$S = P + jQ$ $V = V_r + jV_i$
r	V_m (Voltage module) δ (Voltage angle)		$V = V_m \cdot e^{\delta}$
I (Current) Z (Impedance) Y (Admittance)	I_r (Real current) R (Resistance) G (Conductance)	I_i (Imaginary current) X (Inductance) B (Susceptance)	$I = I_r + jI_i$ $Z = R + jX$ $Y = G + jB$

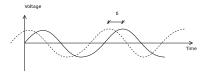


Figure 4.1: Voltage delay.

Table 4.1: Electrical magnitudes and their units.

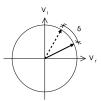


Figure 4.2: Voltage delay in phasor representation in the complex plane.

Table 4.2: Magnitudes and their real and imaginary complex components.

Components connection and their conversions

Let us assume a three-phase grid. The phases of the grid are denoted by the names of *A*, *B* and *C*. There are two main connection types that arise: Wye (like the letter "y") and Delta.

The wye and delta connections provide the ground to introduce the *phase* and *line* voltages. The phase to neutral voltage is called *phase voltage*, those are V_A , V_B and V_C . The phase to phase voltage is called *line voltage*, those are V_{AB} , V_{AC} and V_{BC} .

The delta connection is depicted in the figure 4.3. The delta connection has no neutral.

The wye connection is depicted in the figure 4.4. The star connection does have neutral (N). The wye connection is also known as star connection.

Delta to Wye ($\Delta \to Y$) The transformation of a three phase connected shunt element in Delta to it's Wye equivalent is:

$$Elm_{Wye} = D \times Elm_{Delta}$$
 (4.1)

Where *D* is:

$$D = \begin{bmatrix} 1 & -1 & 0 \\ 0 & 1 & -1 \\ -1 & 0 & 1 \end{bmatrix} \tag{4.2}$$

For instance, considering the impedances transformation from figures 4.3 and 4.4:

$$\begin{bmatrix} Z_A \\ Z_B \\ Z_C \end{bmatrix} = \begin{bmatrix} 1 & -1 & 0 \\ 0 & 1 & -1 \\ -1 & 0 & 1 \end{bmatrix} \times \begin{bmatrix} Z_{AB} \\ Z_{AC} \\ Z_{BC} \end{bmatrix}$$
(4.3)

Wye to Delta $(Y \to \Delta)$ The transformation of a three phase connected shunt element in Wye to it's Delta equivalent is:

$$Elm_{Delta} = D^{-1} \times Elm_{Wye} \tag{4.4}$$

Where D^{-1} is:

$$D^{-1} = \frac{1}{3} \begin{bmatrix} 1 & 0 & -1 \\ -1 & 1 & 0 \\ 0 & -1 & 1 \end{bmatrix}$$
 (4.5)

For instance, considering the impedances transformation from figures 4.3 and 4.4:

$$\begin{bmatrix} Z_{AB} \\ Z_{AC} \\ Z_{BC} \end{bmatrix} = \frac{1}{3} \begin{bmatrix} 1 & 0 & -1 \\ -1 & 1 & 0 \\ 0 & -1 & 1 \end{bmatrix} \times \begin{bmatrix} Z_A \\ Z_B \\ Z_C \end{bmatrix}$$
(4.6)

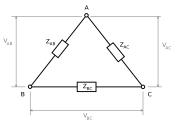


Figure 4.3: Delta connection scheme.

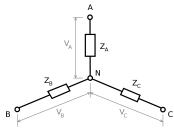


Figure 4.4: Wye connection scheme.

Per unit system

In an electrical grid there are multiple levels of voltage. This situation introduces discontinuities in the numerical methods used to

solve power flows and state estimations among others, producing an unstable convergence behaviour. To avoid this, the per unit system is introduced. A side effect of the per unit representation is to have a very convenient way of visualizing the grid magnitudes, all referenced to their base. In the per unit system, all the voltages are expressed in terms of their nominal value. In this case, all the grid voltage values are around one. For instance, a voltage value of 0.98 means that the voltage is 98 % of the nominal voltage value at that point.

For most exchange formats in computer programs, the element's magnitudes are expressed with a mix of actual units and per unit values. Regardless of this, a practical way of converting any electrical magnitude to its per unit equivalent is presented.

First, we must choose an arbitrary value of power base conversion. This value can be seen as the grid's nominal power, even though that concept is not related to any physical quantity, but it is rather a numerical artifice.

Magnitude	Base
V (Voltage) S (Power)	V_{Base} : terminal's nominal voltage. S_{Base} : Arbitrary value.
<i>I</i> (Current)<i>Z</i> (Impedance)<i>Y</i> (Admittance)	$I_{Base} = S_{Base} / V line_{Base} = S_{Base} / (V_{Base} \cdot \sqrt{3})$ $Z_{Base} = S_{Base} / V_{Base}^2$ $Y_{Base} = 1 / Z_{Base}$

The base power is most commonly chosen to be $S_{Base} = 100MVA$.

Table 4.3: Electrical magnitudes and their per unit base.

Sequence components simplification

Charles L. Fortescue presented in 1918 his famous article ¹ in which he describes how to represent a three-phase element in the so-called *sequence components*.

The main use of this technique is to reduce the amount of impedances needed to represent a line or transformer from usually nine, to three (or even two) if the element is considered to be balanced. An element is considered balance if the impedance in all it's phases is equal and the phase-to-phase coupling impedances are also equal. This is an assumption that is commonly made for transmission grids (very high voltage) and distribution grids in high voltage. This advance allowed the popularization of the single-line diagrams in which every line represents a a number of wires transmitting power in a balanced scheme.

Fortesue defined a transformation matrix A_s and it's inverse as:

$$A_{s} = \begin{bmatrix} 1 & 1 & 1 \\ 1 & a^{2} & a \\ 1 & a & a^{2} \end{bmatrix}$$
 (4.7)

$$A_s^{-1} = \frac{1}{3} \begin{bmatrix} 1 & 1 & 1 \\ 1 & a & a^2 \\ 1 & a^2 & a \end{bmatrix}$$
 (4.8)

¹ Charles L Fortescue. Method of symmetrical co-ordinates applied to the solution of polyphase networks. *Transactions of the American Institute of Electrical Engineers*, 37(2):1027–1140, 1918 Where *a* is the transformation eigenvector:

$$a = 1^{120_{deg}} = 1 \cdot e^{j\frac{2}{3}\pi} = 1 \cdot \cos\left(\frac{2}{3}\pi\right) + 1j \cdot \sin\left(\frac{2}{3}\pi\right) \tag{4.9}$$

$$a^{2} = 1^{-120_{deg}} = 1 \cdot e^{-j\frac{2}{3}\pi} = 1 \cdot \cos\left(\frac{2}{3}\pi\right) - 1j \cdot \sin\left(\frac{2}{3}\pi\right)$$
 (4.10)

Then, any 3x3 impedance matrix representing the rectangular ABC three-phase impedance of an element (line, transformer, capacitor, etc.) can be transformed to a sequence equivalent using the formula:

$$Z_{seq} = A_s^{-1} \times Z_{ABC} \times A_s \tag{4.11}$$

Example Consider the following impedance matrix of a three-phase line. Example from ².

$$Z_{ABC} = \begin{bmatrix} 0.4576 + j1.0780 & 0.1560 + j0.5017 & 0.1535 + j0.3849 \\ 0.1560 + j0.5017 & 0.4666 + j1.0482 & 0.1580 + j0.4236 \\ 0.1535 + j0.3849 & 0.1580 + j0.4236 & 0.4615 + j1.0651 \end{bmatrix}$$

Using equation 4.11, we obtain the sequence impedance matrix:

$$Z_{seq} = \begin{bmatrix} 0.7735 + j1.9373 & 0.0256 + j0.0115 & 0.0321 + j0.0159 \\ -0.0321 + j0.0159 & 0.3061 + j0.6270 & 0.0723j0.0060 \\ 0.0256 + j0.0115 & -0.0723j0.0059 & 0.3061 + j0.6270 \end{bmatrix}$$

For the sequence matrix, the non diagonal elements are neglected. Using only the three diagonal elements as:

$$Z_0 = 0.7735 + j1.9373$$

 $Z_1 = 0.3061 + j0.6270$
 $Z_2 = 0.3061 + j0.6270$

Observe that Z_1 and Z_2 are identical (with the shown numerical precision). It is very common in utilities to store only Z_0 and Z_1 to define a line. The balanced element assumption is very common and should be carefully used.

Building Z_{ABC} from the sequence components

Once the complete 3x3 impedance matrix has been reduced to the sequence components and only those have been stored in the utility database, the obtaining of the full 3x3 matrix might be necessary to perform unbalanced calculations. Of course we will not be able to obtain the exact original Z_{ABC} from the reduced sequence components, but the approximation is fair.

The approximated full impedance matrix is obtained from the sequence components as:

$$Z_{ABC_{approx}} = \frac{1}{3} \begin{bmatrix} 2Z_1 + Z_0 & Z_0 - Z_1 & Z_0 - Z_1 \\ Z_0 - Z_1 & 2Z_1 + Z_0 & Z_0 - Z_1 \\ Z_0 - Z_1 & Z_0 - Z_1 & 2Z_1 + Z_0 \end{bmatrix}$$
(4.12)

² William H Kersting. *Distribution* system modeling and analysis. CRC press, 2012

Example We need to compute two values, before assembling the 3x3 matrix:

$$\frac{1}{3}(2 \cdot Z_1 + Z_0) = \frac{1}{3}(2 \cdot (0.3061 + j0.6270) + (0.7735 + j1.9373)) = 0.4619 + j1.0638$$

$$\frac{1}{3}(Z_0 - Z_1) = \frac{1}{3}((0.7735 + j1.9373) - (0.3061 + j0.6270)) = 0.1558 + j0.4368$$

The the approximated impedance matrix is:

$$Z_{ABC_{approx}} = \begin{bmatrix} 0.4619 + j1.0638 & 0.1558 + j0.4368 & 0.1558 + j0.4368 \\ 0.1558 + j0.4368 & 0.4619 + j1.0638 & 0.1558 + j0.4368 \\ 0.1558 + j0.4368 & 0.1558 + j0.4368 & 0.4619 + j1.0638 \end{bmatrix}$$

We observe that the calculation outcome is a fair approximation of the original Z_{ABC} and that the approximated matrix is symmetric, matching the balanced assumption, which the original impedance matrix did not fully comply with. Therefore, is the original Z_{ABC} was reduced assuming a balanced impedance distribution when that was not the case, if we build the approximated matrix, we will never know if it represents the reality.

5 The branch element

To the effect of most calculations run in operation of a electrical system, the lines, transformers, and any other element that connects two nodes are represented by the so-called Π model.

л П Model

The pi model is composed by a series admittance Y_{serie} and a shunt admittance Y_{sh} divided in two. The shunt admittances are connected at the sending and receiving terminals (primary and secondary). To accommodate the possibility of regulating the voltage at the sending and/or receiving terminals, two per-unit transformers are included as well. The per unit transformers are modelled with the *tap ratio* parameters α and β .

In the case of lines, the series admittance Y_{serie} is computed as the inverse of Z_{ABC} . From the line's calculation it is obtained the series impedance matrix Z_{ABC} and the shunt admittance matrix Y_{sh} .

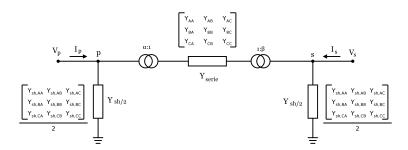


Figure 5.1: General branch model.

The generalized admittance matrix that corresponds to the $\boldsymbol{\Pi}$ model is:

$$\begin{bmatrix} I_p \\ I_s \end{bmatrix} = \begin{bmatrix} Y_{pp} & Y_{ps} \\ Y_{sp} & Y_{ss} \end{bmatrix} \times \begin{bmatrix} V_p \\ V_s \end{bmatrix}$$
 (5.1)

Where:

$$\begin{array}{c|cccc} Y_{pp} & Y_{ps} & Y_{sp} & Y_{ss} \\ \hline Y_{series} + \frac{Y_{sh}}{2} & -Y_{series} & -Y_{series} & Y_{series} + \frac{Y_{sh}}{2} \\ \alpha \beta & \alpha \beta & \beta \end{array}$$

Table 5.1: Equations of the generalized Π model.

The formula 5.1 will be used to compute the network admittance matrices.

The line element, utilizes the branch model as it has been formulated, only that the use of the tap ratio relations α and β is not necessary.

5.2.1 Overhead lines

5.2.2 Underground cables

Example Let's assume that we have already computed the series impedance and the shunt admittance of a three-phase line. The nominal voltage at the line terminals is 66kV and we choose the base power to be $S_{base} = 100MVA$.

$$Z_{ABC} = \left[\begin{array}{ccc} 0.4576 + j1.0780 & 0.1560 + j0.5017 & 0.1535 + j0.3849 \\ 0.1560 + j0.5017 & 0.4666 + j1.0482 & 0.1580 + j0.4236 \\ 0.1535 + j0.3849 & 0.1580 + j0.4236 & 0.4615 + j1.0651 \end{array} \right] \Omega$$

$$Y_{sh} = \begin{bmatrix} j5.6712 & -j1.8362 & -j0.7034 \\ -j1.8362 & j5.9774 & -j1.1690 \\ -j0.7034 & -j1.1690 & j5.3911 \end{bmatrix} \cdot 10^{-6}S$$

The first thing we need to do is to compute the base magnitudes:

$$Z_{base} = \frac{100MVA}{(66kV)^2} = 0.022956841 \quad \Omega$$

$$Y_{base} = \frac{1}{Z_{base}} = 43.56 \quad S$$

Then we must obtain the line series admittance matrix Y_{series} by inverting the 3x3 matrix Z_{ABC} . The we divide the resulting matrix by Y_{base} . Analogously we can invert the per unit impedance matrix:

$$Y_{series} = \left(\frac{Z_{ABC}}{Z_{base}}\right)^{-1} p.u.$$

$$Y_{series} = \begin{bmatrix} 0.0112 - j0.0232 & -0.0054 + j0.008 & -0.0020 + j0.0052 \\ -0.0054 + j0.008 & 0.0125 - j0.024 & -0.0033 + j0.0063 \\ -0.0020 + j0.0052 & -0.0033 + j0.0063 & 0.0100 - j0.0224 \end{bmatrix} p.u.$$

Dividing the shunt admittance by the base admittance we obtain the per unit shunt admittance:

$$Y_{sh} = \begin{bmatrix} j13.0193 & -j4.2153 & -j1.6148 \\ -j4.2153 & j13.7222 & -j2.6837 \\ -j1.6148 & -j2.6837 & j12.3763 \end{bmatrix} \cdot 10^{-8} \quad p.u.$$

Now we need to find the branch model admittances Y_{pp} , Y_{ps} , Y_{sp} and Y_{ss} .

$$Y_{pp} = Y_{ss} = Y_{series} + Y_{sh}/2 = \begin{bmatrix} 0.0112 - j0.0232 & -0.0054 + j0.008 & -0.0020 + j0.0052 \\ -0.0054 + j0.008 & 0.0125 - j0.024 & -0.0033 + j0.0063 \\ -0.0020 + j0.0052 & -0.0033 + j0.0063 & 0.0100 - j0.0224 \end{bmatrix} p.u.$$

$$Y_{ps} = Y_{sp} = -Y_{series} = \begin{bmatrix} -0.0112 + j0.0232 & 0.0054 - j0.008 & 0.0020 - j0.0052 \\ 0.0054 - j0.008 & -0.0125 + j0.024 & 0.0033 - j0.0063 \\ 0.0020 - j0.0052 & 0.0033 - j0.0063 & -0.0100 + j0.0224 \end{bmatrix} p.u.$$

If we convert the three-phase matrices into positive sequence values we obtain:

$$Y_{pp} = Y_{ss} = 0.0148 - j0.0296$$

$$Y_{ps} = Y_{sp} = -0.0148 + j0.0296$$

$$\begin{bmatrix} I_p \\ I_s \end{bmatrix} = \begin{bmatrix} 0.0148 - j0.0296 & -0.0148 + j0.0296 \\ -0.0148 + j0.0296 & 0.0148 - j0.0296 \end{bmatrix} \times \begin{bmatrix} V_p \\ V_s \end{bmatrix}$$

Transformer

The transformer model implements the branch model as well, but the model admittances Y_{pp} , Y_{ps} , Y_{sp} and Y_{ss} vary depending on the connections types at the primary and at the secondary.

The most common transformer connections at the terminals are:

- Delta (Δ)
- Wye (*Y*)
- Grounded Wye (the neutral is grounded) (*Yg*)

Primary	Secondary	Y_{pp}	Y_{ss}	Y_{ps} and Y_{sp}
Yg	Yg	$\frac{1}{\alpha^2} \Upsilon_I$	$\frac{1}{\beta^2} Y_I$	$-\frac{1}{\alpha\beta}Y_I$
Yg	Y	$\frac{1}{3\alpha^2}Y_{II}$	$\frac{1}{3\beta^2}Y_{II}$	$-\frac{1}{3\alpha\beta}Y_{II}$
Yg	Δ	$\frac{1}{\alpha^2}Y_I$	$\frac{1}{\beta^2} Y_{II}$	$\frac{1}{\alpha\beta}Y_{III}$
Υ	Yg	$\frac{1}{3\alpha^2}Y_{II}$	$\frac{1}{3\beta^2}Y_{II}$	$-\frac{1}{3\alpha\beta}Y_{II}$
Υ	Υ	$\frac{1}{3\alpha^2}Y_{II}$	$\frac{1}{\beta^2}Y_{II}$	$\frac{1}{\alpha\beta}Y_{III}$
Δ	Δ	$\frac{1}{\alpha^2} Y_{II}$	$\frac{1}{\beta^2} Y_{II}$	$-\frac{1}{\alpha\beta}Y_{II}$
Δ	Yg	$\frac{1}{\alpha^2} Y_{II}$	$\frac{1}{\beta^2}Y_I$	$\frac{1}{\alpha\beta}Y_{III}$

Table 5.2: Three-phase transformer impedances of the branch model.

The table 5.2 lays out the branch admittances for every pair of connections. The source for this model is the excellent book by J.Arrillaga¹.

$$Y_{I} = \begin{bmatrix} y_{t} & 0 & 0 \\ 0 & y_{t} & 0 \\ 0 & 0 & y_{t} \end{bmatrix}$$
 (5.2)

$$Y_{II} = \begin{bmatrix} 2y_t & -y_t & -y_t \\ -y_t & 2y_t & -y_t \\ -y_t & -y_t & 2y_t \end{bmatrix}$$
 (5.3)

$$Y_{III} = \begin{bmatrix} -y_t & y_t & 0\\ 0 & -y_t & y_t\\ y_t & 0 & -y_t \end{bmatrix}$$
 (5.4)

 y_t is the transformer winding per unit admittance. It is given in values per units from the transformer specifications sheet. Usually it is given either directly as magnetizing resistance r_m and inductance x_m , in which case:

$$y_t = \frac{3}{r_l + jx_l} \tag{5.5}$$

Or it is given as the "short circuit study" values. This is a more complete case.

¹ Jos Arrillaga and CP Arnold. *Computer analysis of power systems*. Wiley Online Library, 1990

5.3.1 Transformer definition from the short circuit study

In order to get the series impedance and shunt admittance of the transformer to match the branch model, it is advised to transform the specification sheet values of the device into the desired values. The values to take from the specifications sheet are:

- S_n : Nominal power in MVA.
- U_{hv} : Voltage at the high-voltage side in kV.
- U_{lv} : Voltage at the low-voltage side in kV.
- *U_{sc}*: Short circuit voltage in %.
- P_{cu} : Copper losses in kW.
- *I*₀: No load current in %.
- GX_{hv1} : Reactance contribution to the HV side. Value from 0 to 1.
- GR_{hv1} : Resistance contribution to the HV side Value from 0 to 1.

Then, the series and shunt impedances are computed as follows:

Nominal impedance HV (Ohm):

$$Zn_{hr} = U_{hr}^2 / S_n$$
 (5.6)

Nominal impedance LV (Ohm):

$$Zn_{lv} = U_{lv}^2 / S_n \qquad (5.7)$$

Short circuit impedance (p.u.):

$$z_{sc} = U_{sc}/100$$
 (5.8)

Short circuit resistance (p.u.):

$$r_{sc} = \frac{P_{cu}/1000}{S_n}$$
 (5.9)

Short circuit reactance (p.u.):

$$x_{sc} = \sqrt{z_{sc}^2 - r_{sc}^2}$$
 (5.10)

HV resistance (p.u.):

$$r_{cu,hv} = r_{sc} \cdot GR_{hv1} \qquad (5.11)$$

LV resistance (p.u.):

$$r_{cu,lv} = r_{sc} \cdot (1 - GR_{hv1})$$
 (5.12)

HV shunt reactance (p.u.):

$$xs_{hv} = x_{sc} \cdot GX_{hv1}$$
 (5.13)

LV shunt reactance (p.u.):

$$xs_{lv} = x_{sc} \cdot (1 - GX_{hv1})$$
 (5.14)

Shunt resistance (p.u.):

$$r_{fe} = \frac{Sn}{P_{fe}/1000}$$
 (5.15)

Magnetization impedance (p.u.):

$$z_m = \frac{1}{I_0/100} \tag{5.16}$$

Magnetization reactance

$$x_m = \frac{1}{\sqrt{\frac{1}{z_m^2} - \frac{1}{r_{fe}^2}}}$$
 (5.17)

If the content of the square root is negative, set the magnetization impedance to zero.

The final complex calculated parameters in per unit are:

Magnetizing impedance (or series impedance):

$$z_{series} = Z_m = r_{sc} + j \cdot x_{sc} \quad (5.18)$$

The series admittance is [p.u.]:

$$y_{series} = \frac{1}{z_{series}} \tag{5.19}$$

Leakage impedance (or shunt impedance):

$$Z_1 = r_{fe} + j \cdot x_m \tag{5.20}$$

$$y_{shunt} = 1/Z_l \tag{5.21}$$

$$y_t = \frac{3}{z_{series}}$$
 (5.22)

we divide the impedance by 3, to reflect the three phases.

Example Let's consider a distribution transformer with the following nameplate characteristics:

- Primary connection: Δ
- $p_{cu} = 6kW$
- Secondary connection: *Yg*
- $p_{fe} = 1.4kW$
- $S_n = 0.5MVA$
- $I_0 = 0.28\%$

• $U_{hv} = 20kV$

• $GR_{hv} = 0.5$

• $U_{lv} = 0.4kV$

• $U_{sc} = 6\%$

• $GX_{hv} = 0.5$

First we obtain the impedance value, from the short circuit study:

$$Zn_{hv}=800$$
 Ω $r_{fe}=357.1429$ $p.u.$ $r_{fe}=357.1429$ $p.u.$ $r_{fe}=357.1429$ $p.u.$ $r_{sc}=0.0600$ $p.u.$ $r_{sc}=0.0120$ $p.u.$ $r_{sc}=17.0103$ $p.u.$ $r_{cu,hv}=0.0060$ $p.u.$ $r_{cu,lv}=0.0060$ $p.u.$ $r_{cu,lv}=8.5052$ $p.u.$ $r_{cu,lv}=8.5052$ $p.u.$ $r_{cu,lv}=0.0000$ r_{cu} $r_{cu}=0.0000$ r_{cu} $r_{cu}=0.0000$ $r_{cu}=0.0000$

Once the we have obtained the y_t value, we start building the appropriate branch impedance matrices Y_{pp} , Y_{ps} , Y_{sp} and Y_{ss} .

$$Y_I = \left[\begin{array}{ccc} 0.0001 - j0.1764 & 0.0000 + j0.0000 & 0.0000 + j0.0000 \\ 0.0000 + j0.0000 & 0.0001 - j0.1764 & 0.0000 + j0.0000 \\ 0.0000 + j0.0000 & 0.0000 + j0.0000 & 0.0001 - j0.1764 \end{array} \right]$$

$$Y_{II} = \begin{bmatrix} 0.0002 - j0.3527 & -0.0001 + j0.1764 & -0.0001 + j0.1764 \\ -0.0001 + j0.1764 & 0.0002 - j0.3527 & -0.0001 + j0.1764 \\ -0.0001 + j0.1764 & -0.0001 + j0.1764 & 0.0002 - j0.3527 \end{bmatrix}$$

$$Y_{III} = \begin{bmatrix} -0.0001 + j0.1764 & 0.0001 - j0.1764 & 0.0000 + j0.0000 \\ 0.0000 + j0.0000 & -0.0001 + j0.1764 & 0.0001 - j0.1764 \\ 0.0001 - j0.1764 & 0.0000 + j0.0000 & -0.0001 + j0.1764 \end{bmatrix}$$

According to the transformer $\Delta \to Yg$ connection, we build the branch model admittances using the formulas for the table 5.2:

$$Y_{pp} = \begin{bmatrix} 0.0002 - j0.3527 & -0.0001 + j0.1764 & -0.0001 + j0.1764 \\ -0.0001 + j0.1764 & 0.0002 - j0.3527 & -0.0001 + j0.1764 \\ -0.0001 + j0.1764 & -0.0001 + j0.1764 & 0.0002 - j0.3527 \end{bmatrix}$$

$$Y_{ss} = \begin{bmatrix} 0.0001 - j0.1764 & 0.0000 + j0.0000 & 0.0000 + j0.0000 \\ 0.0000 + j0.0000 & 0.0001 - j0.1764 & 0.0000 + j0.0000 \\ 0.0000 + j0.0000 & 0.0000 + j0.0000 & 0.0001 - j0.1764 \end{bmatrix}$$

$$Y_{ps} = Y_{sp} =$$

$$\begin{bmatrix} -0.0001 + j0.1764 & 0.0001 - j0.1764 & 0.0000 + j0.0000 \\ 0.0000 + j0.0000 & -0.0001 + j0.1764 & 0.0001 - j0.1764 \\ 0.0001 - j0.1764 & 0.0000 + j0.0000 & -0.0001 + j0.1764 \end{bmatrix}$$

There is no computational need to assemble the 6x6 element impedance matrix in a computer program. The branch model matrices are added to the full circuit matrix instead.

5.4 Voltage regulator

Since all the electrical models are in per-unit values, the voltage regulator is modelled in the exact same way as a transformer.

Decoupling the three-phase model. Delta branch to Y branch transformation

We have presented a three-phase branch model with 3x3 matrices. However we indicated at the beginning of this book that we would not perform the three-phase calculations using the coupled matrices (which would lead to larger linear systems to solve and the impossibility to solve multi-phase grids). This is why we are going to introduce a technique to decouple the phases while keeping the phase-phase influences (the non diagonal elements of the branch element matrix.)

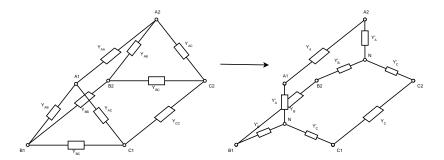


Figure 5.2: Conversion from the Delta branch model to the Y branch model.

The three-phase series admittance matrix can be immediately split into three single-phase systems each of them represented by one of the admittances of the diagonal of the three-phase matrix. However by doing this we would be neglecting the phase coupling

that is present in the 3x3 matrix.

$$Y_{series,\Delta} = \begin{bmatrix} Y_{AA} & Y_{AB} & Y_{AC} \\ Y_{BA} & Y_{BB} & Y_{BC} \\ Y_{CA} & Y_{CB} & Y_{CC} \end{bmatrix} \longrightarrow Y_{AA}, Y_{BB}, Y_{CC}, couplings?$$

In order to take into account the phase couplings we need to perform the delta to star conversion of the magnitudes.

$$\begin{bmatrix} Y'_A \\ Y'_B \\ Y'_C \end{bmatrix} = \begin{bmatrix} 1 & -1 & 0 \\ 0 & 1 & -1 \\ -1 & 0 & 1 \end{bmatrix} \times \begin{bmatrix} Y_{AB} + Y_{AC} \\ Y_{BC} + Y_{BA} \\ Y_{CA} + Y_{CB} \end{bmatrix}$$
(5.23)

 Y'_A , Y'_B and Y'_C are the shunt magnitudes that appear by embedding the phase-phase coupling admittances by performing the delta-star transformation.

Now we can perform the same operation on the shunt admittance matrix:

$$Y_{shunt,\Delta} = \begin{bmatrix} Y_{sh,AA} & Y_{sh,AB} & Y_{sh,AC} \\ Y_{sh,BA} & Y_{sh,BB} & Y_{sh,BC} \\ Y_{sh,CA} & Y_{sh,CB} & Y_{sh,CC} \end{bmatrix} \longrightarrow Y_{sh,AA}, Y_{sh,BB}, Y_{sh,CC}, couplings?$$

$$\begin{bmatrix} Y'_{sh,A} \\ Y'_{sh,B} \\ Y'_{sh,C} \end{bmatrix} = \begin{bmatrix} 1 & -1 & 0 \\ 0 & 1 & -1 \\ -1 & 0 & 1 \end{bmatrix} \times \begin{bmatrix} Y_{sh,AB} + Y_{sh,AC} \\ Y_{sh,BC} + Y_{sh,BA} \\ Y_{sh,CA} + Y_{sh,CB} \end{bmatrix}$$
(5.24)

Finally we can formulate the three single-phase systems:

• System A

$$Y_{series} = Y_{AA}$$

$$Y_{shunt} = Y_{sh,AA} + Y'_A + Y'_{sh,A}$$

• System B

$$Y_{series} = Y_{BB}$$

 $Y_{shunt} = Y_{sh,BB} + Y'_B + Y'_{sh,B}$

• System C

$$Y_{series} = Y_{CC}$$

$$Y_{shunt} = Y_{sh,C} + Y'_{C} + Y'_{sh,C}$$

In Matrix form:

$$\begin{bmatrix} Y_{series,decoupled,A} \\ Y_{series,decoupled,B} \\ Y_{series,decoupled,C} \end{bmatrix} = \begin{bmatrix} Y_{AA} \\ Y_{BB} \\ Y_{CC} \end{bmatrix}$$
 (5.25)

$$\begin{bmatrix} Y_{sh,decoupled,A} \\ Y_{sh,decoupled,B} \\ Y_{sh,decoupled,C} \end{bmatrix} = \begin{bmatrix} Y_{sh,A} \\ Y_{sh,B} \\ Y_{sh,C} \end{bmatrix} + \begin{bmatrix} 1 & -1 & 0 \\ 0 & 1 & -1 \\ -1 & 0 & 1 \end{bmatrix} \times \begin{bmatrix} Y_{AB} + Y_{AC} + Y_{sh,AB} + Y_{sh,AC} \\ Y_{BC} + Y_{BA} + Y_{sh,BC} + Y_{sh,BA} \\ Y_{CA} + Y_{CB} + Y_{sh,CA} + Y_{sh,CB} \end{bmatrix}$$

$$(5.26)$$

6 The bus and it's connected elements

The bus model is based upon the traditionally called ZIP bus model to indicate that the same model includes an impedance (Z), a current (I) and a power (P) value. To be consistent with our nomenclature, here we are going to refer to it as the YISV model since at the node there is voltage(V), power (S), admittance (Y) and current (I) potentially connected.

The admittance, power and currents connected to the bus are derived from the elements connected to it (loads, generators, capacitor banks, etc.)

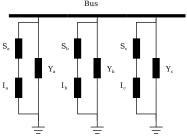


Figure 6.1: YISV Bus model.

- 6.1 The substation
- 6.2 Load
- 6.3 Voltage controlled generator
- 6.4 Battery
- 6.5 Capacitor banks

7 Topology analysis and consolidation

- 7.1 Islands detection
- 7.2 Calculation of the bus connected phases
- 7.3 Calculation of the admittance matrices
- 7.4 Calculation of the voltage, power and current vectors

8 Power flow

The power flow problem consists in finding the node voltages that correspond to the injected current and power values. The fundamental equation to be solved is:

$$S = V \times (Y \times V - I)^* \tag{8.1}$$

Where:

- *S*: Vector of power injections.
- *I*: Vector of current injections.
- *V*: Vector of voltages.
- Y: Admittance matrix.

Types of buses

At each node, we can consider the following magnitudes to exist:

- |V|: Voltage module.
- δ : Voltage angle.
- *P*: Active power injection/consumption.
- *Q*: Reactive power injection/consumption.

In some nodes we know the active power injection (generation nodes), in some nodes we just know the consumption power (load nodes), and in some nodes we artificially define the voltage (slack nodes). The power flow problem is formulated such that in every node two variables are set and the other two are computed. Thus we can define the node types as:

Bus type	V	δ	P	Q
PQ	Calc.	Calc.	Set	Set
PV	Set	Calc.	Set	Calc.
VD	Set	Set	Calc.	Calc.

There must be at least one VD or slack node in order to be able to compute the power flow problem.

Note that the equation 8.1 is non-linear when solving for the voltage (V). However if there are no power injections and only current injections, the equation becomes linear. Because of the non-linearity, we will be using iterative methods to solve this equation.

Table 8.1: Bus types.

Z-Matrix

The Z-Matrix method is a simple iterative method that provides good results, especially in distribution grids where the voltage variations are not large.

The Z-Matrix method requires the Kron reduction of the slack nodes of the circuit. This reduction is explained below:

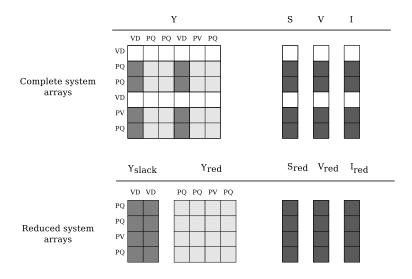


Figure 8.1: Circuit's admittance matrix and vectors reduction. Representation for a six node circuit with two slack nodes.

$$Y_{red} = Y[pqpv, pqpv] (8.2)$$

$$Y_{slack} = Y[pqpv, vd] \tag{8.3}$$

$$V_{red} = V[pqpv] (8.4)$$

$$V_{slack} = V[vd] (8.5)$$

$$S_{red} = S[pqpv] \tag{8.6}$$

$$I_{red} = I[pqpv] \tag{8.7}$$

Once we have all the reduced magnitudes, it means that we have removed the slack (VD) nodes from the circuit, but we must keep their influence in the rest of the nodes in the form of current injections.

First we copy V_{red} into another variable V_{prev} .

$$V_{prev} = V_{red} (8.8)$$

The current injections appearing by the removal of the slack nodes are:

$$I_{slack} = Y_{slack} \times V_{slack} \tag{8.9}$$

The voltages that arise from the slack current injections are:

$$C_k = Y_{red} \times I_{slack} \tag{8.10}$$

The total injected currents including the power injections are:

$$I_k = \frac{S_{red}}{V_{prev}} + I_{red} \tag{8.11}$$

The next step is to compute the nodes voltage due to all the power and current injections:

$$V_k = Y_{red}^{-1} \times I_k - C_k \tag{8.12}$$

Now the error is the infinite norm of the voltage difference.

$$error = ||V_{prev} - V_k||_{\infty} = max(abs(V_{prev} - V_k))$$
(8.13)

The infinite norm of the voltage difference between the previous iteration and the current iteration is a poor convergence criteria, but it is usually the only working criteria in practice for this method.

Next, we copy the voltage solution to the previous voltage solution vector:

$$V_{nrev} = V_k \tag{8.14}$$

Repeat equations 8.11 to 8.14 until the error is less or equal to a given tolerance not too strict like 1×10^{-3} .

Finally, since we have found the voltages for the PQ and PV nodes only, we should return a voltage vector with the voltages for all the nodes. Because of this, we copy the voltage solution for the reduced system into a copy of the original voltage vector.

$$V[pqpv] = V_k \tag{8.15}$$

Jacobian based power flow

8.3

The derivative based methods are more accurate than the derivative free methods. They are usually more robust at the cost of being more computationally expensive. In this section we'll start by introducing the Jacobian matrix. It is the derivative of the power flow equation [8.1] with respect to a set of voltage values. The section continues by introducing the methods that use this matrix to solve the power flow problem.

The Jacobian linear system ($J \times \Delta x = F$), allows the computation of the voltage increment (Δx) to apply to the voltage solution (V) in order to achieve a voltage that solves the power flow equation 8.1.

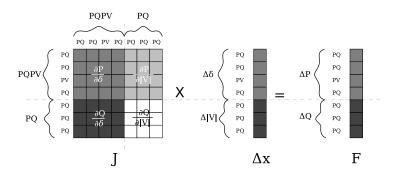


Figure 8.2: Jacobian based linear system to compute the voltage increments.

This process is iterative and recursive, hence to have a initial voltage solution in the vicinity of the final solution is crucial. Otherwise the method will converge to a non physical solution. A common good-enough voltage initial solution is to set all the voltages according to the slack voltage value.

.3.1 Building the Jacobian matrix

The Jacobian matrix (J) is the derivative of the fundamental power flow equation 8.1, with respect to the voltage magnitudes to be solved |V| and δ . It is one of the most expensive parts of the derivative based power flow methods.

In most literature, the reader will find the Jacobian presented with non matrix formulas which are hard to understand in their relation to the Jacobian matrix formation. Ray D. Zimmerman came up with a fantastic way of building the Jacobian matrix ¹ by using lineal algebra operations over the complex matrices and vectors that represent the circuit. The formulation presents a huge advantage: No use of cosine and sine operations for the derivatives computation. The calculation of trigonometric functions is a recursive operation and should be avoided whenever possible in high performance computing. The formulation is presented as follows.

First we compute the total current injection. It is a sparse diagonal matrix, so do not use a dense matrix format.

$$I_{diag} = diag(Y \times V - I) \tag{8.16}$$

Then we convert the buses voltage solution V into a square diagonal matrix. Use a sparse format for this.

$$V_{diag} = diag(V) (8.17)$$

Compute the buses voltage normalized solution V into a square diagonal matrix. This is acomplished by dividing each complex voltage by its absolute value. Use a sparse format for this.

$$E_{diag} = diag(V/|V|) \tag{8.18}$$

Now, Compute the derivative of the power injections S with respect to the voltage module |V|:

$$\frac{\partial S}{\partial |V|} = V_{diag} \times (Y \times E_{diag})^* + I_{diag}^* \times E_{diag}$$
 (8.19)

¹ Ray D Zimmerman. Ac power flows, generalized opf costs and their derivatives using complex matrix notation, 2010

Compute the derivative of the power injections S with respect to the voltage angle δ :

$$\frac{\partial S}{\partial \delta} = 1j \cdot V_{diag} \times (I_{diag} - Y \times V_{diag})^*$$
(8.20)

Finally, assemble the Jacobian matrix J. The Jacobian matrix is sparse and only contains real values. If the circuit contains npq buses of type PQ and npv buses of type PV, the Jacobian matrix is a square matrix with 2npq + npv rows and the same number of columns. See the figure 8.2.

$$J = \begin{bmatrix} real \left(\frac{\partial S}{\partial \delta}[pqpv, pqpv] \right) & real \left(\frac{\partial S}{\partial |V|}[pqpv, pq] \right) \\ imag \left(\frac{\partial S}{\partial \delta}[pq, pqpv] \right) & imag \left(\frac{\partial S}{\partial |V|}[pq, pq] \right) \end{bmatrix}$$
(8.21)

Observe that the Jacobian matrix is composed of four subsets of the previously computed derivatives $\frac{\partial S}{\partial \delta}$ and $\frac{\partial S}{\partial |V|}$, where combinations of the PQ and PV node indices are selected.

pqpv is a vector that contains the indices of the PQ and PV type indices in sequential order.

pq is a vector that contains the indices of the PQ type indices in sequential order.

Newton-Raphson

Newton-Raphson is a recursive, iterative numerical method that minimizes the value of a function. In our case the value to minimize is the power mismatch:

$$s_{calc} = V \times (Y \times V - I_{bus})^* \tag{8.22}$$

$$\Delta S = S_{specified} - S_{calc} \tag{8.23}$$

Since the mismatch (ΔS) is a vector of complex values, we transform it into a vector (F) with the real part of ΔS for the PQ and PV bus indices and the imaginary part of ΔS for the PQ bus indices of the vector. Observe that the length of the vector is 2npq + npv, matching the Jacobian dimensions.

$$F = \begin{bmatrix} \Delta P \\ \Delta Q \end{bmatrix} \tag{8.24}$$

$$\Delta Q = imag(\Delta S[pq]) \tag{8.25}$$

$$\Delta P = real(\Delta S[pqpv]) \tag{8.26}$$

Then we define the error to minimize as the infinite norm of F. The infinite norm is the same as the maximum absolute value. If the error is less than the specified tolerance, we stop the iterations otherwise, continue until a certain number of iterations considered as the maximum (for example 20).

$$error = ||F||_{\infty} = max(abs(F))$$
 (8.27)

If the error is higher than the tolerance, we need to solve the voltages increment (Δx), for which we must compute the Jacobian first. Some simplifications of the method only compute the Jacobian one time, but for better accuracy, the Jacobian needs to be calculated on every iteration.

$$\Delta x = J^{-1} \times F \tag{8.28}$$

Notice that Δx is a vector with the voltage angles (δ) for the PQ and PV nodes followed by the voltage modules (|V|) for the PQ nodes. This matches the way we built the Jacobian (J) and the mismatch (F). Now we must assign those polar voltage angle and module increments to the rectangular voltage vector that we use to compute the Jacobian and the mismatch.

To accomplish this, we declare two polar voltage increment vectors $\Delta |V|$ and $\Delta \delta$, initialize them to zero, and we fill the corresponding values with the following:

$$\Delta \delta[pqpv[i]] = \Delta x[i] \quad \forall i \in 0..npqpv \tag{8.29}$$

Here we use the auxiliary vectors *pqpv* and *pq* which contain the indices of the pq and pv type buses, and the pq type buses respectively. *npqpv* and *npq* are the sizes of those two vectors.

$$\Delta |V|[pq[i]] = \Delta x[i + npqpv] \quad \forall i \in 0..npq \tag{8.30}$$

Then we need to add those values to the voltage module and angle vectors |V| and δ :

$$|V|^{(k+1} = |V|^{(k} + \Delta|V|^{(k)}$$
(8.31)

$$\delta^{(k+1)} = \delta^{(k)} + \Delta \delta^{(k)} \tag{8.32}$$

At last, we convert the polar vectors |V| and δ into a single complex voltage vector V.

$$V^{(k+1)} = |V|^{(k+1)} \cdot \left(\cos(\delta^{(k+1)}) + 1j \cdot \sin(\delta^{(k+1)})\right)$$
(8.33)

The method consists in repeating formulas 8.22 to 8.33 until the error is less or equal to the tolerance or a number of iterations is reached.

The algorithm is then:

- 1. Start.
- 2. Compute the mismatch function (*F*) using the initial voltage solution (*V*). Equation 8.24.
- 3. Compute the error. Equation 8.27.
- 4. While *error* > *tolerance* or *iterations* < *max_iterations*:
 - (a) Compute the Jacobian
 - (b) Solve the linear system. Equation 8.28.
 - (c) Assign Δx to V. Equations 8.29 to 8.33.
 - (d) Compute the mismatch function (F) using the latest voltage solution (V). Equation 8.24.
 - (e) Compute the error. Equation 8.27.
 - (f) iterations = iterations + 1
- 5. End.

Here we introduce the notion of the k_{th} iteration wich represents the current iteration and the $k_{th} + 1$ represents the next iteration.

Levenberg-Marquardt

The Levenbarg-Marquardt is a recursive and iterative technique that is usually not related to power flow, but rather to non-linear least squares problems. Nevertheless, it solves the exact same problem as Newton-Raphson, but in a much more robust manner. It is advised when Newton-Raphson does not converge, because it exhibits excellent convergence properties at the cost of a higher computational effort.

At each iteration we need to assemble the system matrix A as follows:

$$A = I^{\top} \times I + \lambda \cdot I \tag{8.34}$$

Where λ is computed only in the first iteration as:

$$\lambda = 0.001 \cdot max(diag(J \times J^{\top})) \tag{8.35}$$

The right hand side of the linear system to obtain the voltage increments is:

$$rhs = I^{\top} \times F \tag{8.36}$$

Solve the voltage increments vector Δx . As in Newton-Raphson, this vector has 2npq + npv elements and the structure is the same as the one explained in the Newton-Raphson method. See figure 8.2.

$$\Delta x = A^{-1} \times rhs \tag{8.37}$$

Compute the objective function to minimize f. It is a value.

$$f = \frac{1}{2} \cdot (F \times F^{\top}) \tag{8.38}$$

Calculate the decision function ρ . It is a value.

$$\rho = \frac{f^{(k-1) - f^{(k)}}}{\frac{1}{2} \cdot (\Delta x \times (\lambda \cdot \Delta x + rhs)^{\top})}$$
(8.39)

Now, based on the value of ρ we decide what to do in the next iteration.

If ρ is greater than zero, we mark a flag to update the Jacobian in the next iteration, and we update the voltage solution using Δx as described in equations Equations 8.29 to 8.33. We also need to modify λ :

$$\lambda = \lambda \cdot max(\frac{1}{3}, 1 - (2 \cdot \rho - 1)^3)$$
 (8.40)

And set a variable $\nu = 2$.

If ρ is less or equal to zero, then the values of Δx will not improve the solution and we need to take corrective actions. We mark a flag to not to update the Jacobian in the next iteration, we set $\lambda = \lambda \cdot \nu$ and $\nu = \nu \cdot 2$.

The complete algorithm is:

F is provided in the equation 8.24.

 $f^{(k-1)}$ is the value of f in the previous iteration. $f^{(k)}$ is the value of f computed in the current iteration. The value of $f^{(k-1)}$ in the first iteration should be a very large number i.e. 1×10^9 .

- 1. Start.
- 2. Compute the mismatch function (*F*) using the initial voltage solution (V). Equation 8.24.
- 3. Compute the error. Equation 8.27.
- 4. While *error* > *tolerance* or *iterations* < *max_iterations*:
 - (a) If the computation of the Jacobian is enabled, compute the Jacobian and the system matrix A using equation 8.34.
 - (b) Compute the linear system right hand side using equation 8.36.
 - (c) Solve the linear system to obtain Δx . Equation 8.37.
 - (d) Compute the objective function f using equation 8.38.
 - (e) Compute the decision value ρ using equation 8.39.
 - (f) If $\rho > 0$:
 - i. Assign Δx to V. Equations 8.29 to 8.33.
 - ii. Update λ using equation 8.40.
 - iii. Set $\nu = 2$.
 - iv. Set the Jacobian update flag to true.
 - (g) If $\rho \leq 0$:
 - i. Assign Δx to V. Equations 8.29 to 8.33.
 - ii. Update $\lambda = \nu \cdot \lambda$.
 - iii. Update $\nu = 2 \cdot \nu$.
 - iv. Set the Jacobian update flag to false.
 - (h) Compute the mismatch function (*F*) using the latest voltage solution (V). Equation 8.24.
 - (i) Compute the error. Equation 8.27.
 - (j) iterations = iterations + 1
- 5. End.
- Holomorphic embedding
 - Post voltage solution: Compute the power flows

Ironically, what is called the power flow problem does not compute the power flows in the branches of a grid. It computes the node voltages. In this section we explain how to compute the current and power that flows through the grid branches using a given voltage solution vector.

9 Time series power flow

10 Stochastic power flow

10.1 Cumulative Density Function (CDF)

10.2 Monte Carlo

10.3 Latin Hypercube

11 State estimation

12 Short-circuit

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