# Automation Engineering 

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## Foreword

This book will examine the foundations of automation engineering in a world increasingly focused on the development and implementation of automation. It will provide the reader with an understanding of the key principles and components of automation engineering and how these principles can then be combined and implemented in a real industrial system to provide safe, economically viable, and efficient systems. Application will focus on a wide range of different systems including chemical, electrical, and mechanical systems.

At the end of each chapter, questions testing the reader's understanding of the material presented as well as providing places for extension and deeper understanding of the topics at hand are given.

In order to clearly differentiate between regular text and computer-related symbols, the font "Courier New" is used for all computer-related symbols.

This book is meant to be used as the course material for an introductory, bachelor's-level course in automation engineering. It is possible to present different permutations and combinations of the material depending on what topics are of interest to the particular students.

Computer files and related material can be downloaded from the book website (please insert the appropriate book website once the book has been published).

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# Chapter 1: Introduction to Automation <br> <br> Engineering 

 <br> <br> Engineering}

Automation engineering is an important component of modern industrial systems that focuses on the development, analysis, optimisation, and implementation of complex systems to provide safe, economically viable, and efficient processes. Automation engineering seeks to eliminate as much as possible human intervention into the process. However, it should be noted that this does not mean that humans are not required to monitor and assist with the running of the process; it simply implies that the mundane, often repetitive, tasks are delegated to computer systems that are better suited for performing such work.

In order to understand automation engineering, it is important to briefly review its long history, its main principles, and its foundations.

## Section 1.1: The History of Automation Engineering

Ever since humans developed the need to implement complex tasks, the desire for making them faster and easier was also present. All such endeavours have sought to harness the power of nature using ingenious methods to provide the desired outcomes. Often, these devices have had practical or military implications.

One of the first to develop an interest in automating processes were the ancient Greeks, who developed a wide range of devices. Since these devices acted on their own, they were often called automata (singular: automaton; from the Greek $\alpha$ v̇ó $\mu \alpha \tau \tau^{1}{ }^{1}$, which means acting of one's own will). These automata performed a wide variety of tasks and were first described by Homer
 tripods. The first known device with feedback control is the water clock developed by Ctesibius (gr: K $\eta \boldsymbol{\sigma} \dot{i} \beta$ ıos; fl. 285 to 222 BC ) that was able to accurately maintain the time. In fact, this water clock remained the most accurate time-keeping device until the invention of the pendulum clock

[^0]in AD 1656 by Christiaan Huygens. Later, a primitive steam engine, called an æolipile, was developed by Hero of Alexandria (gr: "Hp $\omega v$ ó ${ }^{A} \lambda \lambda \varepsilon \xi \alpha v \delta \rho \varepsilon$ ́s, c. AD 10 to 70). Examples of these devices are shown in Figure 1. As well, objects that assisted in the computation of heavenly bodies (essentially the first computers) were also developed. The most famous is the Antikythera mechanism, a type of astronomical clock based on a gear-driven apparatus. This tradition of developing mechanical automated objects was continued long into the Middle Ages both in Europe and in the Middle East, with such work as the Book of Ingenious Devices (ar: كتاب الحيل (Kitab alHiyal) or pe: كتاب ترفندها (Ketab tarfandha)) by the brothers Banu Musa published in A.D. 850, which describe various automata, including primitive control methods, automatic fountains, mechanical music machines, and water dispensers. Similarly, in the court rooms around the world, various automata in the shape of singing animals were being developed and maintained. Famous examples can be found in the (now destroyed) palaces of the Khanbaliq of the Chinese Yuan dynasty and the court of Robert II, Count of Artois.

The interest in the development of automata continued into the Renaissance with the development of life-size automata, such as The Flute Player by the French engineer Jacques de Vaucanson (1709-1782) in 1737.

With the rediscovery of steam engines, the ability to develop large and complex automated system become a reality sparking the first Industrial Revolution (1760 to 1840). One of the first such examples was the Jacquard loom that could be programmed to automatically weave different designs using punch cards. The development of advanced systems required methods of controlling them, so that explosions and damage were minimised. The first dedicated device for controlling a steam engine, called the Watt governor, was developed by James Watt (1736-1819). The Watt governor, shown in Figure 2, regulates the amount of fuel entering the system so that the speed of the engine remains in a desired range. The development of the first Watt governor was followed by a rash of patents trying to improve various aspects, including one by William Siemens (1823-1883). It was not until 1868, when James Clerk Maxwell (1831-1879) provided a mathematical description of the governor in his aptly named paper On Governors that a rigorous mathematical foundation for the development of methods for controlling a process was available. As the systems became more complex in the succeeding decades and centuries, there came an ever-increasing need for understanding these intricate systems in order that they be properly run so as to avoid unsafe operating conditions.


Figure 1: Automation engineering in the time of the ancient Greeks: (left) Æolipile (steam engine) and (right) Automated temple door opening device


Figure 2: Watt governor (a: fly ball, b: arms, c: spindle, d: sleeve, e: ball crank lever, and f: throttle)

The demand for automation continued with the onset of the second Industrial Revolution (1870-1915), which focused on the development of efficient manufacturing methods (production lines, Taylorism, and similar ideas) coupled with the development of electricity and the first electrical devices.

By the 1950s, a new industrial revolution, often called the third Industrial Revolution or the Digital Revolution, had started. This revolution focused on the implementation and use of complex electrical circuits that can be used to quickly and efficiently perform complex processes. With the development of these circuits, it became easy and cost-effective to implement automation in a wide range of different fields. From the perspective of automation engineering, the key event was the development of programmable logic controller (PLCs) that could be used to implement advanced control methods in an industrial setting. The first PLCs were developed by Bedford Associates from Bedford, Massachusetts, USA based on a white paper written by the engineer Edward R. Clark in 1968 for General Motors (GM). One of the key people working on this project was Richard E. "Dick" Morley (1932-2017), who is often considered the father of the PLC. Other important work was performed by Odo Josef Struger (1931-1998) of Allen-Bradley in the period 1958-1960. At the same time, there was an explosion of interest in a theoretical perspective on automation engineering, especially in the areas of control and process optimisation. This research by such people as Andrei Kolmagorov (ru: Андре́й Никола́евич Колмого́ров, 1903 - 1987), Rudolf Kálmán (hu: Kálmán Rudolf Emil, 1930 - 2016), Richard E. Bellman (1920-1984) and others lead to a strong foundation for the subsequent development and implementation of advanced control methods in industry.

Within the context of the third Industrial Revolution, the concept of robots was also considered. The word robot itself was first used by the Czech author Karel Čapek (1890-1938) in his 1920 drama R.U.R. (cz: Rossumovi Univerzální Roboti or en: Rossum's Universal Robots) as a word for artificial humanoid servants that provided cheap labour. Karel credited his brother painter Josef Čapek (1887-1945) with inventing this word. The word robot stems from the Czech root robota, that means serfdom, which ultimately comes from the proto-Indo-European word, * $h_{3} e r b^{h-}$, that means "to change or evolve status" from which the English word arbitrate is also derived.

More recently, with the growth of interconnectedness and the development of smart technologies, some have proposed that a new industrial revolution is dawning. This revolution has
been called the fourth Industrial Revolution, or Industry 4.0, which focuses on the development of self-functioning, interconnected systems in an increasingly globalised world. Its main drivers are ever increasing automation and digitalisation of the industrial plant combined with globalisation and customisation of the supply chains.

## Section 1.2: The Key Concepts in Automation Engineering

Automation engineering can be applied to a wide range of different cases. Instead of considering each situation separately, automation engineering has developed a set of abstract concepts that allow the ideas to be applied to any relevant situation.

The basic concept in automation engineering is a process or system. ${ }^{2}$ Figure 3 shows a typical system with some important components. In automation engineering, as shown in Figure 3, a system consists of inputs, denoted by $\boldsymbol{u}$, and outputs, $\boldsymbol{y}$. Inputs represent those variables, whose change will lead to a change in the system. Inputs can be divided into two types: manipulative and disturbance. A manipulative input is an input, whose value can be changed through some device, for example, the flow rate in a valve can be manipulated by opening or closing a valve. A disturbance input is an input, whose value cannot be changed (at least in context of the given situation), for example, the ambient temperature cannot be readily modified. Similarly, outputs can be divided into two categories: observable or measurable and unobservable or unmeasurable. Observable or measurable outputs are those variables whose value can be measured or inferred using some device, for example, the temperature of a fluid can be measured or determined using a thermometer. Unobservable or unmeasurable outputs are those variables that cannot be measured (at least in context of the given situation), for example, the density of a complex mixture consisting of multiple phases and components may not be easy to measure. Finally, we can consider the states of the system, which are the internal variables that describe the behaviour of the system. In many cases, the states of a system are equivalent to the outputs and can be categorised using a similar terminology. Often, not all the states can be observed.

[^1]

Figure 3: A system in automation engineering
It goes without saying that if we wish to understand the system, we require a model of the system. A model of the system is a mathematical representation of the relationship between the inputs, outputs, and states. The complexity of the required model depends on the purposes that the model will be used. Modelling a process is a complex endeavour that requires insight into the process and the ability to handle large data sets quickly and effortlessly.

In order to complete the picture regarding the system, it is necessary to extend our view to include parts that allow us to interact with the system or influence its behaviour. Such a view is provided in Figure 4, which shows the key components and their interactions. It is important to briefly consider the impact of these in order to understand how the system can be influenced and how it will react. In this view, the eight key components can be described as (the number is the same as in Figure 4):

1) Actual Process: This represents the process under consideration. Normally, the actual process is unknown. Instead, a model of the actual process is used.
2) Sensors: The sensors provide the ability to measure the process and understand how the variables are changing.
3) Actuators: The actuators allow the value of a variable to be changed. If we cannot change the value of a variable, then it is hard to use it in an automation strategy. When deciding on which actuators to use, it is important to consider such factors as the variable being manipulated, the automation requirements (e.g. required accuracy, tolerance, or precision), and the type of service required (e.g. continuous, discrete, or emergency).
4) Automation Devices: The automation devices are the controllers and related components that are used to automate the process. Most of the time the automation hardware consists of computers and other digital devices, such as programmable logic controllers (PLCs), that implement the required functions. The design of the automation hardware (and software) requires knowledge of the limitations and requirements of the system.
5) Environment: The environment represents everything that surrounds the system to be automated that can have an impact on the overall performance of the system. This influence can be caused either by changes in other processes that interact with the system of interest
or by direct environmental changes, such as for example, changes in the ambient temperature. The system will exchange mass, energy, and information with its environment.
6) Automation Objectives: The automation objectives play a very important role in the development and implementation of the final automation system. Poorly defined or unclear objectives can make achieving the project difficult if not impossible. Furthermore, the objectives often need to be translated from the language of business into actually implementable objectives on a system. This translation can cause additional uncertainties and lack of clarity.
7) Operators: Although the human component is often minimised or ignored when designing automation systems, it is in fact very important. Many complicated automation systems have failed due to a lack of proper consideration of the operators. In general, the operators need to have the required information easily available (no fancy graphics are needed) and they can enter the required information into the system quickly and efficiently. Appropriate feedback and safety checking of the entered values must be performed to avoid confusion. The operators interact with the process using a human-machine interface (HMI). An HMI provides two key functions. It allows the operators to see the important process values and to manipulate as necessary the process values. Manipulating the values implies that the operators can change at what value the process operates, for example, changing the flow rate in a pipe. Finally, when designing the HMI, it is important to consider any safety features, such as logic, that limits which values can be entered by the operators. This prevents mistakes both accidental, such as mistyped numbers or the wrong information in a given field, and malicious, such as changes introduced by illegal access to the system, from having an impact on the process.
8) Disturbances: Disturbances are everything that can impact the system, but whose presence cannot be directly controlled. Disturbances can originate in the environment (for example, the ambient temperature) or in the devices themselves (for example, measurement noise in sensors). One of the objectives of the automation system is to minimise, as much as possible, the impact of these disturbances on the process.

The final aspect that needs to be considered is safety. The automation system that has been designed should allow the process to operate in a safe manner without any unexpected behaviour, from serious errors to a system failure. The system should also be robust, that is, minor changes
in the process should not cause the whole system to fail catastrophically. A robust system can handle small changes in the process conditions and still attain the required goals.


Figure 4: General structure of an automated system

## Section 1.3: Automation-Engineering Framework

When solving an automation engineering problem, the following steps should be considered:

1) Modelling of the Process, which involves developing an appropriate model of the process.
2) Analysis of the Process, which involves using the model to analyse how the process will behave under different conditions. This can be performed using either mathematical analysis or simulations.
3) Design of the Automation Strategy, which based on the process properties and required behaviour, provides an appropriate automation strategy that attains the required automation objectives.
4) Validation of the Proposed Automation Strategy, which validates the proposed automation strategy using the process model. This determines if the proposed strategy can, in fact, achieve all the desired automation objectives. Should it be found that the proposed strategy is lacking, then the strategy needs to be refined and retested. This implies that there may need to iterate until the final automation strategy is found.
5) Implementation and Commissioning of the Proposed Automation Strategy, which involves the implementation of the strategy in the real process. Naturally, the implementation on the real system may lead to changes in the strategy. This implies that before commissioning, the automation strategy should be tested on the actual process in as realistic conditions as possible. In certain cases, this may not be feasible and advanced simulations, using hardware-in-the-loop methods can be implemented to provide a realistic simulation of the system.

## Section 1.4: The Automation-Engineering Pyramid

The automation-engineering pyramid is an overall description of the way in which different automation strategies can be organised and structured. It examines two key components: speed of response, or how often the system is expected to respond to changes, and process details. Figure 5 shows the general pyramid that consists of 6 different levels (from top to bottom):

- Level 5 - Enterprise Resource Planning (ERP) level: This level focuses on the abstract analysis of the overall company strategies for dealing with the current market conditions. This level runs at a very long-time horizon, often in terms of years. At this level, the focus is on market analysis, strategic investment and personnel planning, and corporate governance.
- Level 4 - Manufacturing Execution Level: This level considers the specific strategies that allow for the overall plant/process to run. This includes such details as production planning, production data acquisition, organisation of delivery orders, and deadline monitoring. The time horizon for actions can reach into months. Often, at this level, various software systems, such as a manufacturing execution system (MES) or a management information system (MIS), are used.
- Level 3 - Process Control Level: This level focuses on short-term production planning, quality control, and maintenance planning with an action horizon of at most days. Often,
software, such as a supervisory control and data acquisition (SCADA) programme, will be used.
- Level 2 - Control Level: This level considers how to run the algorithms for controlling the process using the available process information on the time horizon of minutes or even seconds. Various hardware, such as programmable logic controllers (PLCs) or industrial PCs (IPCs) are used to compute the required actions. It is on this level that this book focuses.
- Level 1 - Field Level: This level focuses on the sensors and actuators. Information is actively exchanged with Level 2, often using fieldbuses. The time horizon is often in the range of seconds or faster.
- Level 0 - Process Level: This level represents the actual process that is running in real time.

It can be noted that the process description becomes more abstract as we go from bottom to top. While Level 0 is very concrete and describes everything in great detail, Level 5 is very abstract and only focuses on describing the process inputs and outputs to give an overall picture of the process.


Figure 5: A description of the components of the automation-engineering pyramid

## Section 1.5: Chapter Problems

Problems at the end of the chapter consist of three different types: (a) Basic Concepts (True/False), which seek to test the reader's comprehension of the key concepts in the chapter; (b) Short Exercises, which seek to test the reader's ability to compute the required parameters for a simple data set using simple or no technological aids. This section also includes proofs of theorems; and (c) Computational Exercises, which require not only a solid comprehension of the basic material, but also the use of appropriate software to easily manipulate the given data sets.

## Section 1.5.1: Basic Concepts

Determine if the following statements are true or false and state why this is the case.

1) A system consists of inputs, outputs, and states.
2) Inputs are variables that influence the system.
3) All outputs can always be measured.
4) A disturbance is a variable whose value can be easily changed as desired.
5) The state of a system describes the internal behaviour of the system.
6) Actuators are used to manipulate disturbances.
7) PLCs are commonly used to automate the process.
8) The environment has minimal impact on an automation system.
9) HMIs should be easy to read and understand.
10) An automation system that explodes periodically is a well-designed system.
11) An automation system should fail the second the system deviates from the expected conditions.
12) A system exchanges mass, energy, and information with the environment.
13)Disturbances can affect sensors, actuators, and control devices.
13) Validating a proposed automation strategy using appropriate models of the system is a good strategy to consider.
15)Before commissioning of the automation strategy, it should be tested on the actual system in as real conditions as possible.
14) The enterprise resource planning level focuses on controlling the process in fine detail and making decisions every millisecond.
15) The process control level often uses SCADA systems to implement its tasks.
16) In the field level, sensors gather the information and transfer it using fieldbuses to the control level.
17) A process variable that cannot be measured should be used to control the process.
18) Safety is always an unimportant topic in automation engineering.

## Section 1.5.2: Short Exercises

These questions should be solved using only a simple, nonprogrammable, nongraphical calculator combined with pen and paper.
21)How is that information can be created and destroyed, but matter and energy cannot be? Provide some examples of such cases.
22) You have been assigned the task of designing an automation system for a traffic light system. Explain how you would apply the automation engineering framework to this problem.
23) You have been given the task of designing a large, multi-unit chemical plant. Explain how the automation-engineering pyramid could apply to this problem.
24) You have been given the task of designing a self-driving car. Explain how you would implement the automation engineering framework to this problem. Would you consider safety and robustness to be significant factors?

## Chapter 2: Instrumentation and Signals

The foundational component of any automation system is the instrumentation, that is, the sensors, actuators, and the computer hardware which together produce a stream of values, often called a signal, that can be used for subsequent processing. Before we can look at the sensors, actuators, and control concepts, it is helpful to understand the types of signals and how they can be generated.

## Section 2.1: Types of Signals

In automation engineering, signals can be classified using two domains: time and value. Each domain has two options: continuous and discrete. In general, a continuous signal can take any value within the set of (positive) real numbers, while a discrete signal can only take certain, specified values (for example, only natural numbers).

In the time domain, a signal is said to be continuous, if there exists a signal value for any time $t$, that is, the signal can be written as a continuous function of $t$. An example of a continuoustime signal would be the outdoor temperature, which contains a value for each possible time instance.

A signal is said to be discrete in the time domain, if it is only defined at certain values $t_{k}$, where $k \in \mathbb{Z}$ (or $\mathbb{N}$ ). Normally, it is assumed that the values are available with a specified constant sampling rate, $t_{s}$, so that $t_{k}=k t_{s}$. Obviously, as the sampling time decreases, then the signal approaches that of a continuous signal.

Similar to the time domain, the value domain of the signal can be classified into continuous or discrete. A continuous value for a signal implies that the signal can take any real number (subject to any physical constraints, for example, always positive or in the range [0, 1]). A continuous valued signal can be written as a continuous function that in general depends on time and any additional variables.

A discrete-valued signal can only take specific values. Often the discrete values are partitioned into equidistant bands. In automation systems, a common discrete-valued signal is a binary signal that can only take two values (conventionally denoted by 0 or 1 ). Such a signal is common when dealing with alarms that are triggered when a certain condition occurs, for example, if the pressure in the reactor surpasses a given value, then the signal value is set to 1 and the alarm
is triggered. Such signals work by assigning one value (say 0 ) to the normal state and the other value to the alarm state (say 1). Binary signals are often displayed using what is called a timing diagram. In a timing diagram, the binary signal is plotted as a function of time. Multiple different binary signals are normally placed on separate $y$-axis, but a common time $x$-axis. Figure 6 shows such a typical timing diagram.


Figure 6: A timing diagram for two binary signals A and B
Based on the classification in the time and value domains, it is possible to have four different types of signals, which are shown in Figure 7. By convention, a signal that is continuous in both the time and value domains is called an analogue signal, while a signal that is discrete in both the time and value domains is called a digital signal.

Since most real processes require and produce continuous, analogue signals, but computers perform their computations using discrete, digital signals, there is a need to understand how signals can be converted between the two forms. Converting from analogue to digital signals is shown in Figure 8, and consists of three components: sampler, quantiser, and encoder. The sampler measures (samples) the value of the analogue signal on a fixed frequency to convert the signal into the discrete time domain. Next, the quantiser coverts this sampled signal into the available closest value (quantum) to create a digital signal. The encoder simply encodes the digital signal into a
given digital representation that can be used by the computer. This process is often denoted as an analogue-to-digital (A/D) converter.


Figure 7: Continuous and discrete signals


Figure 8: Analogue-to-digital conversion
When quantising a continuous signal, its values are compared against fixed (equidistant) quantisation levels. If the value of the continuous signal is between two decision thresholds, the lower value is usually selected. For example, if a continuous signal has a value of 0.44 with quantisation levels at 0.25 and 0.5 , the quantised (sampled) signal will be set to 0.25 . The selection of appropriate quantization levels is important, since this selection affects the accuracy of the mapping of the process, for example, if the steps are too far apart, important information can be lost.

When going in the other direction, it is normal to assume that the value of the signal will not be changed and only the time component needs to be made analogue. This is normally accomplished using a hold, which holds the value of signal until a new value is received. The most
common hold is the zero-order hold, which simply holds the last value received until a new value is received. A more accurate hold is the first-order or linear hold, which uses an interpolation between the previous two data points to obtain a linearly varying value over the sampling period. This process is called digital-to-analogue (D/A) conversion.

## Section 2.2: Sensors

A sensor is a device that can detect changes in a variable and upon calibration display these changes in a manner that can be understood by others, often with reference to some absolute scale.

A sensor is characterised by two properties: accuracy and precision or reproducibility. Accuracy measures the ability of a sensor to give the "true" value, which is usually determined based on some standard. The difference between the true and measured values is often called bias. Precision or reproducibility measures the variability of the sensor when measuring the same value. Ideally, it is desired that the values reported by a sensor be tightly located about the mean value, that is, the variance of the sensor values should be small. It should be noted that an inaccurate sensor may be very precise with a tight distribution about an incorrect value. Another issue to consider is the range of the sensor. The range of a sensor is defined as the difference between the largest and smallest values that the sensor can measure. The accuracy and precision of a sensor is often a function of the range of values that the sensor is meant to measure. The larger the range the less precise the values will be. Similarly, the smaller the range, the more precise the values will be. This implies that when dealing with processes that have values covering a large range it may be necessary to install multiple sensors which are accurate in a limited region and then using only the appropriate sensor value.

Sensors need to be calibrated before being used or to check that they are behaving as expected. Calibration involves using standards with accurate and well-defined values to compare against the measured value given by the sensor. Plotting the measured and true values against each other will allow the calibration of the sensor to be determined. A typical calibration curve is shown in Figure 9. There are two parameters of interest here: the $y$-intercept, which gives the bias, and the slope of the line (or lack of linearity). The slope of the line and the general distribution of the points suggest whether or not an appropriate calibration curve was used. Ideally, the line should be linear and the slope equal to exactly one.


Figure 9: Typical Calibration Curve
A single sensor depending on its calibration and physical arrangement can be used to measure different physical variables, for example, a differential pressure cell can measure both flow rates and level.

Selecting an appropriate sensor depends on the following criteria:

1) Measurement Range (Span): The required measurement range for the process variable should lie within the instrument's range.
2) Performance: Depending on the specific application, different factors, such as accuracy, precision, and speed of response, will need to be considered.
3) Reliability: How well does the sensor work in the given operating conditions, for example, if the sensor must be placed in harsh operating conditions, can it handle them and for how long?
4) Materials of Construction: Depending on the application, the required materials of construction for the sensor may be different, for example, a temperature sensor in a blast furnace will require a different material than a temperature sensor in the living room of a house.
5) Invasive or Noninvasive: An invasive sensor comes in direct contact with the object being measured, for example, inserting a probe into a liquid to measure the temperature. If an invasive sensor comes into contact with the process, it can influence the process or be influenced by the process itself. Thus, invasive methods can have issues with long-term accuracy due to fouling or corrosion of the probe surface. On the other hand, noninvasive sensors do not come in contact with the process. In such cases, the process is not disturbed,
but the measurement may be less accurate. However, noninvasive sensors are generally easier to use and can be easily retrofitted into an already built environment.

## Section 2.2.1: Pressure Sensor

In general, most industrial pressure sensors are constructed using a transducer that can convert the force per given area (pressure) into an electric signal. Nonelectronic pressure sensors, often called manometers, do not produce an electric signal and often have a calibrated faceplate that allows the value to be read. For this reason, the signals produced cannot, in most cases, be used for industrial automation. A pressure sensor can either measure an absolute pressure or a pressure difference. A simplified schematic of these two possibilities is shown in Figure 10. Most pressure sensors measure a pressure difference, which is often expressed as deviations from atmospheric pressure. As shown in Figure 10 (a), this is accomplished by leaving one of the two taps or sides of a pressure sensor open to the atmosphere. Such a reading is often called gauge pressure. A negative gauge pressure is expressed by using the term vacuum, for example, a reading of 10 kPa vacuum, would imply that the pressure of 10 kPa below atmospheric pressure has been achieved. Instead of using the varying atmospheric pressure, it is possible to fix the pressure value on one side of the sesnsor and use the resulting system to measure the system pressure. As shown in Figure 10 (b), such a set up will allow the measurement of absolute pressure to be achieved.


Figure 10: Measurement set-up for pressure sensors: (a) Measuring differential pressure and (b) Measuring absolute pressure (a: flexible membrane, $P_{a}$ : ambient pressure, $P_{m}$ : to-be-measured pressure, and $P_{f}$. fixed pressure)

Most transducer-based pressure sensors use some type of electric circuit (for example, a Wien bridge) to measure the strain induced by the pressure on the system. Common strain gauges include piezoresistive, capacitive, electromagnetic, piezoelectric, and optical. Figure 11 shows a typical high-pressure transducer-based pressure sensor. Similarly, manometers are also based on the effect of pressure on some system property. They include the common hydrostatic manometers,
which basically measure the difference in pressure between the two taps, and mechanical manometers, which measure the effect of pressure on the strain of the material. Mechanical manometers have the advantage that they do not interact strongly with the fluid and can provide very sensitive readings. On the other hand, compared to the hydrostatic manometers, they can be more expensive.


Figure 11: High-pressure transducer-based pressure sensor (a: measuring diaphragm, b: strain gauge, c: reference hole to the atmosphere, d : true gauge diaphragm, and e: area for the resistor for temperature correction and internal electronic amplifiers)

## Section 2.2.2: Liquid-Level Sensors

Determining the liquid level in tanks or other similar containers is a very common need in many chemical plants. Liquid level can be determined using many different methods, including differential pressure cells, floats, and various radio-based methods. Of these, the most common approach is to use a differential pressure cell, which measures the pressure gradient between the top of the liquid and the bottom of the liquid. Since pressure is proportional to the height of the liquid in the tank, it is relatively easy to calibrate and determine the height. On the other hand, since density depends on the temperature, this approach will not work in cases where there can be wide fluctuations in temperature (for example, a boiling liquid). A float measurement device works on a similar principle, of pressure difference, but records the values differently. Figure 12 shows an example of float to measure the level in a toilet for controlling the flow of water into the toilet tank. This shows a relatively simple example of automation that can be implemented with an appropriately selected sensor. Finally, various radio-based methods, for example, ultrasonic pulse
generating devices, can be used to determine the surface level. However, in order to get an accurate estimate, it is required that the surface be relatively flat and consistent. Froths and other particles can impact the accuracy and precision of the measurements.


Figure 12: Level measurement and control using a float (a: trip lever, b: refill tube, c: float, d: overflow tube, e: ballcock, f: lift chain, and $g$ : flush valve)

## Section 2.2.3: Flow Sensors

Flow sensors are used to measure the speed at which a liquid or gas is moving. There are three main types of flow sensors: mechanical flow sensors, pressure-based flow sensors, and electromagnetic flow sensors (including optical and ultrasonic flow sensors). Depending on the fluid present, each of the three types will have different accuracies and characteristics.

Mechanical flow sensors are based on the idea of timing how long it takes the liquid or gas to fill some known unit of volume. Most mechanical flow sensors have some type of wheel or paddle that is turned by the flowing medium inducing an electric signal, which is then calibrated to give a flow rate. In general, mechanical flow sensors are good with simple fluids, for example, water, in a single phase over a limited range of flow rates. Suspended particles in the fluid as well as multiple phases can cause the mechanical flow sensor to give incorrect readings.

Pressure-based flow sensors measure the pressure difference caused by some constriction in flow to determine the flow rate. Common pressure-based flow sensors include Venturi tubes, orifice-plate differential pressure cells, and Pitot tubes. As with mechanical flow sensors, these flow sensors tend to work best with uniphase flow of simple fluids without any suspended particles. Figure 13 shows the basic operating principles of a Venturi tube.


Figure 13: Venturi tube (a: datum, b: U-tube manometer, c : manometer fluid, d: inlet, e: outlet, f: main pipe, g: converging cone; h: throat, i: diverging cone, $D_{1}$ : diameter of the main pipe, $D_{2}$ : diameter of the throat, $Z_{1}$ : reference height $1, Z_{2}$ : reference height 2 , and $h$ : height difference in the manometer)

Electromagnetic flow sensors use various electromagnetic waves (including light) to measure the flow of the fluid. Although not strictly speaking an electromagnetic flow sensor, ultrasonic flow sensors are based on a similar principle and can be included here. Magnetic flow sensors measure the changes induced by a flowing fluid in a local electric field. In many cases, the flowing fluid should be conductive and the surrounding pipe a nonconductor. Optical flow sensors use the time it takes for small, suspended particles inside a flowing gas to cross two laser beams. Based on this time, the bulk flow rate can be computed. Finally, ultrasonic flow sensors use the Doppler effect to measure the flow rate. An advantage of the ultrasonic flow sensors is that they can be nonintrusive, that is, the measurement can be done without coming in contact with either the pipe or the fluid. On the other hand, these sensors require the speed of sound for the flowing fluid to be known in order to calibrate the results correctly.

In most industrial cases, pressure-based flow meters are used, since they are simple to use, robust, and easy to maintain. On the other hand, when dealing with exotic or extreme fluids, then more complex methods may be required.

## Section 2.2.4: Temperature Sensors

Temperature sensors are used to measure the temperature of a stream, be it liquid or gas. The most common industrial temperature sensor is a thermocouple, where temperature changes induce a voltage (called the Seebeck effect) in the two different touching conductors. It is possible to calibrate this change with temperature and hence use it. The construction of the thermocouple is shown in Figure 14. Determining an appropriate span and calibration for a thermocouple is very important in order to allow for maximal benefit. Inappropriate calibration can lead to accuracy issues with the sensor. The most common types of thermocouples are shown in Table 1.


Figure 14: Thermocouple (a: metal 1, b: metal 2, c: measurement point, d: reference location, $T_{1}$ : to-be-measured temperature, $T_{2}$ : reference temperature, and V : voltmeter)

Table 1: Thermocouple Types

| Thermocouple <br> Type | Temperature Range <br> $\left({ }^{\circ} \mathrm{C}\right)$ | Material of <br> Construction | Comments |
| :---: | :---: | :--- | :--- |
| K | -200 to 1350 | chromel-alumel | Cheap, not too precise |
| E | -110 to 140 (narrow) | chromel-constantan <br> -50 to 740 (wide) | Good for cryogenic use |
| J | -40 to 750 | iron-constantan | More sensitive than K |
| B and R | 50 to 1800 | platinum-rhodium | Good for high temperature |
|  |  | alloys | use, expensive |
|  |  |  |  |


| Thermocouple <br> Type | Temperature Range <br> $\left({ }^{\circ} \mathrm{C}\right)$ | Material of <br> Construction | Comments |
| :---: | :---: | :---: | :--- |
| C, D, and G | 0 to 2320 | tungsten/rhenium <br> alloys | Cannot be used in <br> presence of oxygen, <br> expensive. |
| Chromel- <br> gold/iron | -273.15 to 25 | chromel-gold-iron | Cryogenic applications |

## Section 2.2.5: Concentration, Density, Moisture, and Other Physical Property Sensors

There exist various online sensors that can quickly provide readings for many different types of physical properties. Unfortunately, most of these sensors are relatively limited in their accuracy or precision. This often results from the extreme or nonideal conditions in which such systems are used. Often the laboratory measurements are more accurate and trusted. Most of these sensors use various photonic, magnetic, or sonic pulse methods to determine concentrations or densities. Consider, for example, a moisture sensor for water in soil. At present, there are 4 main ways to measure the moisture: tensiometers, which measure the water tension in the soil; electrical resistance blocks, which measure the resistance of a ceramic block in contact with the soil; electrical conductivity probes, which measure the conductivity of the soil between 2 plates; and dielectric sensors, which measure the dielectric constant of the soil. All methods require regular maintenance and are sensitive to ambient conditions, for example, placing more fertiliser will change the electrical conductivity of the moisture and hence disrupt the sensor.

## Section 2.3: Actuators

An actuator is a device that can change the amount of some variable that enters a system. An actuator is characterised by three properties: accuracy, precision or reproducibility, and performance. Accuracy measures the ability of an actuator to give the "true" value, which is usually determined based on some standard. The difference between the true and measured values is often called bias. Precision or reproducibility measures the variability of the actuator when delivering the same value. Ideally, it is desired that the value delivered by an actuator be close to the desired true value. The final issue to consider is the performance or how long does it take for
the actuator to respond to a change. In general, it is desired that an actuator respond quickly to changes and deliver the final desired value in a short period of time. Most actuators are designed so that $0 \%$ corresponds to no value and $100 \%$ corresponds to the maximal value. A common problem with actuators is that their response is nonlinear. One way to resolve this problem is to use a characterisation function that will convert the desired linear values into the actuator's nonlinear values.

Actuators need to be calibrated before being used or to check that they are behaving as expected. Calibration involves using standards with accurate and well-defined values to compare against the measured value by the sensor. However, the exact relationships and behaviour of the calibration will depend on the specific actuator.

There are three common actuators: valves, pumps, and variable current actuators.

## Section 2.3.1: Valves

Valves are one of the most common actuators seen in a plant. They allow the flow rate of a liquid or gas to be controlled. Since valves are so ubiquitous in plants, there is a vast amount of work done on understanding valves and how they impact the performance of automation. The most common type of control valve is a pneumatic control valve that uses air to change the flow rate. A typical control valve is shown in Figure 15. A pneumatic control valve consists of three components: the current-to-pressure converter (I/P converter), the valve itself, and a positioner. The I/P converter takes the electric 4-20 mA signal and converts it to an appropriate pressure signal that can provide the motive force required to move the piston. There exist two types of control valves: air-to-close and air-to-open. Air-to-open valves are also called fail close valves because in the event of a loss of air pressure, the valve will close. Similarly, air-to-close valves are called fail open valves because in the event of a loss of air pressure, the valve will remain open. The choice of air-to-open or air-to-close valves is based on the outcome of a process hazard review.

For an air-to-close valve as shown in Figure 15, the compressed air enters the top of the valve actuator, exerts force against the diaphragm, and moves the diaphragm until the resistance force from the spring is equal to the force on the diaphragm from the compressed air. An increase in air pressure will tend to push the diaphragm down, while a decrease in air pressure will result in the spring forcing the diaphragm up to a new equilibrium point. The valve stem is attached to the diaphragm, and as the stem moves up and down it changes the position of a tapered plug (or
"trim") relative to a seat. As the stem moves, it changes the cross-sectional area available for flow, and thus the resistance to flow. For an air-to-close valve, an increase in air pressure will push the plug towards the seat, thus reducing the area available for flow, increasing the resistance to flow, and reducing the flow rate though the valve. For an air-to-open valve, the air enters on the bottom of the diaphragm, so an increase in air pressure will raise the diaphragm, thus opening the valve.


Figure 15: Pneumatically Actuated Control Valve (air-to-close) (a: supply air connection, b: mechanical stop, c: diaphragm, d: upper chamber, d: lower chamber, f: spring, g : housing, h : local position indicator, i : stem, j : flow direction, and k : transducer)

Many modern valves have an additional element called a positioner that seeks to overcome any potential errors in the valve. A positioner basically compares the current valve location against the reference value and will change the air supply to allow for the difference to be zero.

The behaviour of a valve is normally specified based on the percentage of the total distance that the valve is opened or closed. This eliminates the need to know the exact flow rates. Therefore, the flow rates of a valve are often stated in terms of percent open (often abbreviated as \%open).

Properly selecting a valve for automation (or control) is very important. There are two variables to consider: sizing and dynamic performance.

## Section 2.3.1.1: Valve Sizing

Control valves must be specified, just like piping, heat exchangers, and other process equipment. The sizing of a control valve determines the range of flow rates over which the valve can produce a hydrodynamically stable flow.

A valve that is too small will not permit enough flow when it is fully open, which is defined in terms of the needs for process regulation. A control valve must permit significantly greater flow than the steady-state requirement in order to be able to provide reasonable performance. A valve that is too large will have a sufficiently high maximum flow but will provide poor regulation when the flow is low. For a number of reasons, valves are not very precise instruments, and the relative errors tend to be greatest when the valve is almost closed.

A valve that is open by less than $10 \%$ is generally considered effectively closed, and a valve that is more than $90 \%$ open is generally considered effectively fully open.

To determine whether a control valve is undersized or oversized, examine the range of controller output values used when the valve is in service. If the valve spends a significant fraction of the time fully open, then it can be considered undersized. If it spends a significant fraction of the time less than $10 \%$ open, or if the flow through the valve reaches a maximum before the valve is fully open, then the valve is oversized.

## Section 2.3.1.2: Dynamic Performance of Valves

In most flow control applications, it is desirable for the flow through the valve to be a linear function of valve position. If a plot of valve \%open as a function of steady-state flow rate produces a straight line, then the valve is said to be linear. There exist two types of valves which will give a nonlinear plot: quick-opening valves and equal-percentage valves, whose plots are shown in Figure 16. A quick-opening valve, as its name suggests, will quickly open and reach the maximal flow rate, while an equal-percentage valve will reach the maximal flow rate more slowly. Since this behaviour of the valve is determined by the manufacturer, it is often called the inherent valve characteristic. If the valve is nonlinear, then the control performance can be improved by including a characterisation block that converts the linear flow rates desired into the corresponding percentage, rather than assuming a linear relationship.

Furthermore, since a valve is a mechanical system, friction will cause two additional types of nonlinearities to be present in the system: static and dynamic. Static nonlinearity refers to the nonlinearity introduced by static friction when initiating changes in the valve position, while dynamic nonlinearity results from the dynamic friction in the valve, usually between the valve stem and the seal. Static friction between the stem and seal results in the valve not moving when there is a small change in the control signal (air pressure) input to the valve. Static friction, or "stiction," can significantly affect controller performance.

To find static and dynamic nonlinearities in a control valve, step the valve from $0 \%$ open to $100 \%$ open and then back to $0 \%$ in small steps and record the steady-state flow at each step. Plot the data points on a graph of steady-state flow versus \%open, and use different lines for opening and closing. The resulting graph will be similar to that shown in Figure 17, where the ideal valve behaviour is shown as a dashed line. Static nonlinearity consists of the deadband region over which the flow rate does not change even though the signal to the valve increases and the slip-jump behaviour, while dynamic nonlinearity consists of hysteresis, which is the gap between the two lines. Slip-jump behaviour is caused by the static friction that must first be overcome before the valve changes its position. Since the dynamic friction is much smaller than the static friction, the valve will overshoot the position defined by the force as soon as it starts to move. This will make the curve look like a staircase. Hysteresis arises from the difference between the measured flow rate when opening and closing a valve. This effect can be explained by the different effect of dynamic friction on the moving valve. When the valve is being opened, due to dynamic friction, it will open less than desired. On the other hand, when it is being closed, the valve will be more open than specified, that is, it will close less than desired.


Figure 16: Inherent valve characteristics


Figure 17: Phase plot for the typical behaviour of a valve with stiction (after (Shoukat Choudhury, Thornhill, \& Shah, 2005)). The arrows show the direction in which the values were changed.

## Section 2.3.2: Pumps

Another actuator that can control the flow rate in a stream is a pump. A pump is a mechanical device that takes a fluid, most often a liquid, from a storage tank and moves it to some other location. The main types of pumps are centrifugal pumps, positive-displacement pumps, and axial-flow pumps. In centrifugal pumps, the flow direction changes by $90^{\circ}$ as it moves over the impeller, while in an axial-flow pump the flow direction is not changed. In positive-displacement pumps, the fluid is trapped in a fixed volume and forced (or displaced) into the discharge pipe. The traditional hand pump is a good example of a positive-displacement pump. Of these three types, the most common type is a centrifugal pump. A positive-displacement pump is often used if flows are small or extreme precision is required. Figure 18 shows a schematic diagram of a centrifugal pump, while Figure 19 shows a positive-displacement pump.

Compared to valves, pumps tend to provide better control of the flow rate and have fewer nonlinear characteristics. However, they have much higher energy consumption than valves. The main considerations for a pump, as for valves, are sizing and performance.


Figure 18: Centrifugal Pump (a: inflow, b: impeller, c: shaft, and d: outflow)


Figure 19: Positive-Displacement Pump (a: inflow, b: packing, c: piston rod, d: stuffing-box bushing, e: liner, f: piston, g : working fluid, h: valve, and i: outflow)

## Section 2.3.2.1: Pump Sizing

Control pumps must be specified, just like piping, heat exchangers, and other process equipment. The sizing of a control pump determines the range of flow rates over which the pump can produce a hydrodynamically stable flow.

An undersized pump will not permit enough flow when it is fully operating, which is defined in terms of the needs for process regulation. A control pump must permit significantly greater flow than the steady-state requirement in order to be able to provide reasonable performance. An oversized pump will have a sufficiently high maximum flow, but will provide poor regulation when the flow is low.

Pumps operating for most of the time at less than about $10 \%$ are said to be oversized, while those operating at more than $90 \%$ are said to be undersized.

## Section 2.3.2.2: Dynamic Performance of Pumps

Unlike for a valve, the performance of a control pump is easier to quantify. In general, pumps are linear. The only significant issue is that with certain types of pumps a deadband may be present at low flow rates, that is, there may be no observed flow. This can be attributed to the fact that the pump cannot overcome the effects of gravity and friction, and hence, produce a flow.

The behaviour of a pump is characterised by its pump characteristic curve, which is often supplied by the manufacturer of the pump. A typical pump characteristic curve for a centrifugal
pump is shown in Figure 20. The head, H, normally expressed in units of length, such as metres or feet, shows how high a given column of liquid could be lifted by a pump. It represents the effective pressure gradient that the pump can overcome. The efficiency, $\eta$, represents how much work put into the pump is converted into lifting the liquid. As in many engineering applications, the higher the efficiency, the better it is. For centrifugal pumps, the net positive suction head (NPSH) is the minimum head (pressure) at the inlet before cavitation occurs. Cavitation is defined as the boiling of a liquid in a pump, which is evidently a very undesirable event. Therefore, the head at the inlet must be greater than the specified value.


Figure 20: Typical pump characteristic curve for a centrifugal pump

## Section 2.3.3: Variable Current Devices

The final actuator of interest is a variable current device that can modulate (vary) the current entering a device. Since these devices are mechanical, they tend not to have any issues with nonlinearities or undesired behaviour. As with all equipment, sizing can be a problem that needs to be appropriate resolved.

## Section 2.4: Programmable Logic Computer (PLCs)

Programmable logic computers (PLCs) are small computers that have been made robust and rugged. They are often used in industry for controlling processes, such as assembly lines, robots, or complex chemical processes. PLCs allow the collection of the different signals and their subsequent processing to reach a decision or action that may need to be taken. Given the computational power, they can also perform relatively advanced logical and mathematical functions that can be used to control the process.

As shown in Figure 21, a typical PLC consists of 6 key components:

1) Inputs: These allow information from outside the PLC to be incorporated and used. Most often, they are electrical signals coming from sensors or switches.
2) Power Supply: This provides the power required to run the PLC and operate all the circuitry. Internally, most PLCs use a 5-V standard. However, the power supply most often provides $230 \mathrm{~V} \mathrm{AC}, 120 \mathrm{~V} \mathrm{AC}$, or 24 V DC. Furthermore, the power supplies are often built as a replaceable module, so that depending on the application, the appropriate power can be provided, for example, in Europe one could use 230 V , but in North America 120 V .
3) Central Processing Unit (CPU): The CPU is the brain of the PLC that performs all the required instructions, calculations, operations, and control functions.
4) Memory: The PLC must also contain memory or the ability to store relevant information for future use. The amount of memory available depends on the PLC and the programming requirements. Some PLCs can have additional memory cards inserted, so that they have more available memory. There are two main types of memory:
a. Read-Only Memory (ROM): ROM is the permanent storage for the operating system and system data. Since a true ROM cannot be changed, in practice, an erasable programmable ROM (EPROM) is used, so that it is possible to update the operating system for the PLC.
b. Random-Access Memory (RAM): RAM is used to store all other information required by the PLC including any programmes and variables. Accessing RAM is very fast. However, when power is lost, the information in RAM is also lost. Therefore, a PLC will always have an additional battery to maintain power to the RAM, so that the information that is contained in the RAM is not lost during a power outage.
5) Communications Unit: The communications unit allows for the PLC to interact with other devices to exchange information using different types of protocols. Often, this exchange of information involves sending new or updated programmes to the PLC. In general, the communications unit will often include the ability to communicate with an operator panel, printers, networks, or other computers.
6) Outputs: These allow the PLC to exchange information with other devices to cause them to take an action. Such devices include motors, valves, pumps, and alarms.

In a PLC, communication between the components occurs using groups of copper wires called buses. A bus consists of a bundle of wires that allow for the transmission of binary information, for example, if the bus contains eight wires, then it is possible to transmit up to 8 bits of information per bus. A typical PLC consists of four buses:

1) Data Bus: The data bus is used to transfer information between the CPU, memory, and I/O.
2) Address Bus: The address bus is used to transfer the memory addresses from which the data will be fetched or to which data will be written.
3) Control Bus: The control bus is used to synchronise and control the traffic circuits.
4) System Bus: The system bus is used for input-output communications.


Figure 21: Layout of a PLC
Since a PLC operates more or less similarly to that of a computer, this means that in order for a PLC to do anything useful, it must be programmed. The basic idea for a PLC is to monitor and control a process. A process in this context is anything that requires monitoring and control and can range from a simple unit that requires its output to be maintained at a prespecified value to a complex, highly interacting system like a room with multiple heating ovens, lights, ventilations systems, and windows, wherein the temperature, carbon-monoxide levels, and humidity must be monitored and maintained at safe levels. In PLC terms, a process is said to be in a given operating
mode $^{3}$, when the process is operating in some specified mode, for example, when the pump is on, off, or operating. Finally, the PLC requires a user programme that takes the inputs, outputs, and internal information to make decisions about the process.

Before we look at how a PLC operates, it can be useful to note the different operating modes (states) that a PLC itself can be in. In general, the operating modes for a PLC are programming, stop, error, diagnostic, and run. In programming mode, a PLC is being programmed, usually using an external device. In stop mode, the PLC is stopped and will only perform some basic operations. In error mode, the PLC has encountered some sort of problem and has stopped working. In diagnostic mode, the PLC runs without necessarily activating any inputs or outputs, allowing for the validity of the programme to be determined. Often, test signals are used in place of the real inputs and outputs. In run mode, a PLC is actively working and performing the requested actions. Each PLC manufacturer may call these operating modes different names and not all of them may be present for a given PLC.

Naturally, the most important mode from the perspective of automation is the run mode. Thus, its behaviour will be examined in greater detail. In run mode, the PLC performs the same four operations in a repeating cycle:

1) Internal Processing, where the PLC checks its own state and determines if it is ready for operation. Should the response from hardware or communication units be lacking, the PLC can give notice of these events by setting a flag, which is an internal Boolean address (or visual indicator) that can be checked by the user to determine the presence of an error state. Normally, the PLC will continue operation, unless the error is serious, in which case, it will suspend operation. In this step, software-related events are also performed. These include such things as updating clocks, changing PLC modes, and setting watchdog times to zero. A watchdog is a timer that is used to prevent a programme from taking too long to execute, for example, it could be stated that if a programme does not terminate in one second, then it could be stuck inside an unending while-loop and an error state will be returned.

[^2]2) Reading Inputs: Next, all the input statuses are copied into memory. This means that the PLC will only use the values from memory rather than checking to make sure that the most recent value is available. This means that at any given point the programme will use the same values during the same scan time. Furthermore, reading values from memory is faster than reading them each and every time from the input.
3) Executing Programmes: After the inputs have been read, the programmes are executed in the order in which the code has been written. The order of execution can be changed by changing the associated priority of the given code or using conditional statements and subroutines. It should be noted that at this point only internal variables and output addresses are updated in memory; the physical outputs are not changed at this point.
4) Updating Outputs: Once the programmes have been finished, the output memory is written so that the state and values of the outputs can be updated. This will then complete a single cycle and the PLC will return to the first step, that is, internal processing.
Since it can be seen that the PLC in run mode performs these four operations repeatedly, the question becomes what is the best approach for how the PLC should repeat these operations. By convention, a single pass through the four operations is called a scan and the scan time or cycle time is the amount of time required to perform a single scan. In practice, the scan time can vary between scans due to the presence of different events or conditions requiring the execution of more or less code. In order to accomplish the repetitions effectively, a programme can be associated with a task, whose execution type can be specified. There are three common execution types:

1) Cyclic Execution: In cyclic execution, the time between scans is fixed. Obviously, the time must be set so that the PLC has the time to complete all the required code. Naturally, certain tasks, such as counting or timing, should always be run on cyclic execution.
2) Freewheeling Execution: In freewheeling execution, as soon as one scan has been completed, then the next scan is started. This is the fastest way of running a task, since there is no waiting between scans.
3) Event-Driven Execution: In event-driven execution, a task is only executed when a given Boolean condition is fulfilled. Event-driven execution is useful for emergency stop routines, start-up routines, and other extraordinary events.
Furthermore, the nature of the task must be specified: can the task be interrupted by another task? If the task can be interrupted, then it is called pre-emptive; if it cannot be, nonpre-emptive. The
difference between these two types of tasks is shown in Figure 22. Finally, the priority of a task must be specified ranging from high to low (the exact details depend on the specific PLC and standard used).


Figure 22: Pre-emptive and nonpre-emptive tasks

## Section 2.5: Communication Devices

The last type of instrumentation is the communication devices which allow all the actuators, sensors, and control logic to communicate with each other. The following equipment is often found:

1) Analogue-to-Digital Converters (A/D Converter), which convert the analogue signal received from the sensor into a digital signal. At this point, quantisation (or resolution) can be an issue. If an insufficient number of levels (decimals) are present, then the resulting data may not carry as much information as before.
2) Digital-to-Analogue Converters (D/A Converter), which convert a digital signal received from the computer/software into an analogue signal that can be used by the actuator. Different methods exist for implementing this conversion including a zero-order hold, where the previous value is maintained until a change occurs, or a first-order hold, where some average of the previous values is used until a new value is available. In most implementations, a zero-order hold is used due to its simplicity and sufficiency.
3) Control Software, which can reside either on the programmable logic computer or as a separate software programme on a computer.
4) Data Historian, which stores all the values for future retrieval. Selecting an appropriate sampling time, or how fast the data are recorded, can determine the usefulness of the stored data for future applications.
5) Network Cables, Switches, and Accessories, which physically connect all the equipment and allow for the strategy to be implemented.

The main issue with the design of the communication units is the available bandwidth. The faster the data is sampled and the more computations that need to be performed, the larger the bandwidth and computational power that will be needed.

In automation, signals can be encoded using many different standards. The two most common standards are the current-based standard of 4 to 20 mA and the pressure-based standard of 3 to 15 psig. ${ }^{4}$ It should be noted that both of these standards do not start at zero, since a value of zero is ambiguous: is the device not working properly or is the value actually zero. By using a live zero, that is, the value of zero corresponds to some nonzero current or pressure means that it is possible to distinguish between the case of a zero value and a faulty device. Furthermore, a live zero allows part of the remaining current or pressure to be used for operating the device without needing additional power supply. This means that the device can be used in remote areas without its own power supply. The lower limit for the live zero was historically determined as the smallest measurable value. In most cases, the upper limit was chosen so that the lower and upper limits are in the ratio 1:5.

## Section 2.6: Chapter Problems

Problems at the end of the chapter consist of three different types: (a) Basic Concepts (True/False), which seek to test the reader's comprehension of the key concepts in the chapter; (b) Short Exercises, which seek to test the reader's ability to compute the required parameters for a simple data set using simple or no technological aids. This section also includes proofs of theorems; and (c) Computational Exercises, which require not only a solid comprehension of the basic material, but also the use of appropriate software to easily manipulate the given data sets.

[^3]
## Section 2.6.1: Basic Concepts

Determine if the following statements are true or false and state why this is the case.

1) An analogue signal is continuous in both the time and value domains.
2) A binary signal is an example of a digital signal.
3) A signal can only be discretised in the value domain.
4) A precise sensor will always give a value close to the true value.
5) If the observer sensor value is 3 kg with a standard deviation of 0.5 kg and the true value is 10 kg , then we can say that the sensor is precise and accurate.
6) A manometer measures the pressure difference using a transducer.
7) A differential pressure cell can measure the level of a liquid in a tank.
8) Venturi tubes are an example of a pressure-based flow sensor.
9) The Doppler effect can be used to measure flow rates.
10) The Seebeck effect allows us to measure pressure changes.
11) A J-type thermocouple can be used to measure the temperature of molten silica which is at least $1000^{\circ} \mathrm{C}$.
12) It is not possible to use thermocouples to measure temperatures below $0^{\circ} \mathrm{C}$.
13) Actuators should be selected so that they are used in the range 20 to $60 \%$.
14) Valves are a type of actuator that restricts flows.
15) An air-to-close valve will remain open should the air supply fail.
16) A quick-opening valve is useful for rapid dosing of a liquid.
17) Slip-jump in valves results from dynamic friction of the moving parts.
18) The efficiency of a pump represents the pressure gradient that it can produce at a given flow rate.
19) A PLC consists of inputs, power supply, CPU, memory, outputs, and communication devices.
20) A bus in a PLC is a location in memory used to store information about where different variables are stored.
21) A PLC in programming mode is being programmed from an external device.
22) A flag in a PLC is a Boolean variable that shows the presence of an error state.
23) A watchdog in a PLC is a variable that prevents the PLC from being interrupted by external users as it is running.
24) A PLC scan is the amount of time it takes for the PLC to read the inputs.
25)Freewheeling execution occurs when the PLC executes a task solely on demand from an external event.
25) A pre-emptive task can never be interrupted.
26) An analogue-to-digital converter is found in all computer-based automation solutions.
27) A data historian stores the data collected from a process.
28) Live zero implies that when the signal value is zero then the current has a value of 4 mA .
29) A PLC normally has a battery to provide power in case of a power failure.

## Section 2.6.2: Short Questions

31) Consider that you have been assigned the task of designing an automation system for a door to monitor who is present at the door and allow the occupant of the house to open the door if necessary. List what sensors, actuators, and other devices would be needed to accomplish this task.
32) Consider the task of monitoring the temperature of a glass furnace. What considerations should you take into account? Which sensors would you use?
33) Consider the task of pumping a mixture of sand, water, oil, air, and various particles. What would you need to consider when designing the pump? What kind of sensors would you consider? Do you think you can achieve highly accurate results?
34) Consider the data shown in Table 2. Create the valve characterisation curve using the data. Determine what type of valve this is and how reproducible the values are. Are there any static or dynamic nonlinearities present? How can you determine this?

Table 2: Data for creating the valve characterisation curve for Question 34)

| \%open | Flow Rate, $\dot{m}(\mathrm{~kg} / \mathrm{min})$ |  |  |
| :---: | :---: | :---: | :---: |
|  | Run 1 | Run 2 | Run 3 |
| 0 | 0 | 0 | 0 |
| 10 | 0.5 | 0.5 | 0.4 |
| 20 | 2.3 | 2.3 | 2.3 |
| 30 | 4 | 4 | 4 |
| 40 | 5.6 | 5.5 | 5.5 |
|  |  |  |  |


| \%open | Flow Rate, $\dot{m}(\mathrm{~kg} / \mathrm{min})$ |  |  |
| :---: | :---: | :---: | :---: |
|  | Run 1 | Run 2 | Run 3 |
| 50 | 6.8 | 6.8 | 6.8 |
| 60 | 8 | 8 | 7.9 |
| 70 | 8.9 | 8.9 | 8.9 |
| 80 | 9.7 | 9.7 | 9.7 |
| 90 | 10.4 | 10.5 | 10.5 |
|  |  |  |  |


| \%open | Flow Rate, $\dot{m}(\mathrm{~kg} / \mathrm{min})$ |  |  |
| :---: | :---: | :---: | :---: |
|  | Run 1 | Run 2 | Run 3 |
| 100 | 11 | 11 | 11 |
| 90 | 10.7 | 10.7 | 10.7 |
| 80 | 10 | 10 | 10 |
| 70 | 9.2 | 9.2 | 9.2 |
| 60 | 8.3 | 8.3 | 8.4 |
| 50 | 7.2 | 7.2 | 7.2 |


| \%open | Flow Rate, $\dot{m}(\mathrm{~kg} / \mathrm{min})$ |  |  |
| :---: | :---: | :---: | :---: |
|  | Run 1 | Run 2 | Run 3 |
| 40 | 6 | 6 | 6 |
| 30 | 4.6 | 4.6 | 4.6 |
| 20 | 3 | 3 | 3 |
| 10 | 1.2 | 1.2 | 1.2 |
| 0 | 0 | 0 | 0 |

35) Consider the sensor data shown in Table 3. It is desired to determine if the sensor values are properly calibrated against the measured values. Determine if the calibration is correct. Are the values reliable?

Table 3: Sensor Calibration Data for Question 35)

| Measured Height <br> (m) | Sensor Value (m) |  |
| :---: | :---: | :---: |
|  | Run 1 | Run 2 |
| 0.00 | 0.02 | 0.03 |
| 0.10 | 0.115 | 0.105 |
| 0.15 | 0.149 | 0.152 |
| 0.20 | 0.229 | 0.215 |
| 0.25 | 0.248 | 0.251 |


| Measured Height (m) | Sensor Value (m) |  |
| :---: | :---: | :---: |
|  | Run 1 | Run 2 |
| 0.30 | 0.321 | 0.312 |
| 0.35 | 0.348 | 0.349 |
| 0.40 | 0.412 | 0.392 |
| 0.45 | 0.452 | 0.457 |
| 0.50 | 0.512 | 0.493 |

## Chapter 3: Mathematical Representation of a

## Process

In order to understand and provide useful information about a process, it is necessary to understand how different processes and systems can be represented. In practice, there exist two main types of representations: mathematical and schematic. A mathematical representation focuses on providing an abstract description of the process that provides information about the process. A good mathematical representation of the system can provide a deep understanding of how the system works and how it will behave in the future. On the other hand, a schematic representation focuses on the relationships between the different components and how they relate to each other. It is primarily a visual approach that allows the actual process to be represented on a piece of paper.

Common mathematical representations include state-space models, transfer function models, and automata.

## Section 3.1: Laplace and Z-Transforms

Before considering the mathematical models themselves, it is helpful to review two very common transformations between the time and frequency domains: the Laplace and $z$-transforms. If the time domain is continuous, then the Laplace transform is used, while if the time domain is discrete, then the $\boldsymbol{z}$-transform is used.

## Section 3.1.1: Laplace Transform

The Laplace transform converts a function from the time domain into the frequency domain. This transform allows for a simple algebraic solution of complex differential equations. Due to this feature, Laplace transforms are widely used in automation engineering to understand process behaviour and obtain solutions to various control problems.

The Laplace transform is defined as

$$
\begin{equation*}
F(s)=\int_{0}^{\infty} f(t) e^{-s t} d t \tag{1}
\end{equation*}
$$

where $F$ is the Laplace transformed function, $f$ the original function, and $s$ the Laplace variable. Conventionally, the Laplace transform is denoted by a Faktur $L$, $\mathfrak{Z}$ (U+1D50F), that is, $F(s)=\mathfrak{R}(f(t))$. Converting from the frequency domain to the time domain is conventionally shown using the inverse of the Laplace transform, $\mathfrak{L}^{-1}$. Table 4 presents a summary of the most common Laplace transforms. The following are some useful properties of the Laplace transform:

1) Linearity: $\mathcal{L}(f+g)=\mathcal{L}(f)+\mathcal{L}(g)$.
2) Superposition: $\mathfrak{L}(\alpha f)=\alpha \mathfrak{L}(f)$.
3) Convolution: $\mathfrak{L}(f * g)=\mathfrak{L}\left(\int_{0}^{t} f(\tau) g(t-\tau) d \tau\right)=F(s) G(s)$.
4) Shifting Properties: The following rules can be useful in solving problems involving Laplace transforms:

$$
\begin{align*}
& \mathfrak{L}(f(t-a) u(t-a))=e^{-a s} F(s)  \tag{2}\\
& \mathfrak{L}(g(t) u(t-a))=e^{-a s} \mathfrak{L}(g(t+a))  \tag{3}\\
& \mathfrak{L}(\delta(t-a))=e^{-a s}  \tag{4}\\
& \mathfrak{L}(f(t) \delta(t-a))=f(a) e^{-a s} \tag{5}
\end{align*}
$$

5) Final Value Theorem: Assuming that the poles (roots of the denominator) lie in the lefthand plane ${ }^{5}$, then

$$
\begin{equation*}
\lim _{t \rightarrow \infty} f(t)=\lim _{s \rightarrow 0} s F(s) \tag{6}
\end{equation*}
$$

6) Initial Value Theorem: The value in the time domain at the starting point $t=0$ is given by

$$
\begin{equation*}
\lim _{t \rightarrow 0} f(t)=\lim _{s \rightarrow \infty} s F(s) \tag{7}
\end{equation*}
$$

Table 4: Table of Common Laplace Transforms

| Case | Time Domain <br> $\boldsymbol{f}(\boldsymbol{t})$ | Frequency Domain <br> $\boldsymbol{F}(\boldsymbol{s})$ |
| :--- | :---: | :---: |
| Dirac Delta or Impulse <br> Function, $\delta$ | $\delta(t-a)= \begin{cases}0 & t \neq a \\ \infty & t=a\end{cases}$ | $e^{-a s}$ |

[^4]| Case | Time Domain <br> $\boldsymbol{f}(\boldsymbol{t})$ | Frequency Domain <br> $\boldsymbol{F}(\boldsymbol{s})$ |
| :--- | :---: | :---: |
| Unit Step Function, $u$ | $u(t)= \begin{cases}0 & t \leq 0 \\ 1 & t>0\end{cases}$ | $s^{-1}$ |
| Polynomials | $\frac{t^{n-1}}{(n-1)!}$ | $\frac{1}{s^{n}}$ |
| Exponential | $e^{-a t}$ | $\frac{(s+a)^{-1}}{(s+a)^{2}+\omega^{2}}$ |
| Cosine | $e^{-a t} \cos (\omega t)$ | $\frac{e^{-a t} \sin (\omega t)}{(s+a)^{2}+\omega^{2}}$ |
| Sine | $\frac{d^{n} f}{d t^{n}}=f^{(n)}$ | $s^{n} F(s)-\sum_{k=1}^{n} s^{k-1} f^{(n-k)}(0) 6$ |
| Derivative | $\int f(t) d t$ | $\frac{1}{s} F(s)$ |
| Integration | $f(t-a) u(t-a)$ | $e^{-a s} F(s)$ |
| Time Shift (or Time Delay) |  |  |

Often, when we are given an equation in the Laplace domain, it may be necessary to convert it into the time domain. This can be performed by using the inverse Laplace transform, $\mathfrak{R}^{-1}$, such that $\mathfrak{L}^{-1}(\mathcal{L}(f(t))=f(t)$. In most automation-engineering problems, the general case reduces to finding the time-domain function corresponding to some rational function of $s$. In such cases, partial fractioning (see Appendix I for the details) is required to break up the original fraction into its constituent parts so that the known functions given in Table 5 can be used.

Table 5: Useful Inverse Laplace Table

| $\boldsymbol{L}(f(\boldsymbol{t}))$ | $\boldsymbol{f}(\boldsymbol{t})$ |
| :---: | :---: |
| $\frac{A}{C s+D}$ | $\frac{A}{C} e^{\frac{-D}{C} t}$ |

${ }^{6}$ If it is assumed that the system is initial at steady-state and using deviational variables, then all derivatives will be zero and this equation will reduce to the first term $s^{n} F(s)$.

| $\mathfrak{L}(f(t))$ | $f(t)$ |
| :---: | :---: |
| $\frac{A}{(C s+D)^{n}}$ | $\frac{A}{(n-1)!C^{n}} t^{n-1} e^{\frac{-D}{C} t}$ |
| $\frac{C s+E}{\alpha s^{2}+\beta s+\gamma}$ <br> (irreducible quadratic) | $\begin{gathered} \left(\frac{E-\frac{C \beta}{2 \alpha}}{\sqrt{\alpha \rho}}\right) e^{\frac{-\beta}{2 \alpha} t} \sin \left(t \sqrt{\frac{\rho}{\alpha}}\right)+\left(\frac{C}{\alpha}\right) e^{\frac{-\beta}{2 \alpha} t} \cos \left(t \sqrt{\frac{\rho}{\alpha}}\right) \\ \text { with } \rho=\gamma-\frac{\beta^{2}}{4 \alpha} \geq 0 \end{gathered}$ |
| $\begin{gathered} \frac{C s+D}{\left(\alpha s^{2}+\beta s+\gamma\right)^{n}} \\ \text { (irreducible } \\ \text { quadratic) } \end{gathered}$ | $\begin{aligned} & \left(\frac{C}{\alpha^{n}(n-2)!}\right)\left[\left(t^{n-2} e^{\frac{-\beta}{2 \alpha} t}\right) \otimes\left(e^{\frac{-\beta}{2 \alpha} t} \cos \left(t \sqrt{\frac{\rho}{\alpha}}\right)\right)^{\otimes_{n}}\right]+ \\ & \left(\frac{2 \alpha D-C \beta}{2 \alpha^{n+1}}\left(\frac{\alpha}{\rho}\right)^{n / 2}\right)\left(e^{\frac{-\beta}{2 \alpha} t} \sin \left(t \sqrt{\frac{\rho}{\alpha}}\right)\right)^{\otimes_{n}} \end{aligned}$ <br> with $n \neq 1, \rho=\gamma-\frac{\beta^{2}}{4 \alpha} \geq 0, \otimes$ is convolution, and $(f)^{\otimes^{n}}$ is defined as the convolution of $f$, $n$ times, i.e. $\underbrace{f \otimes f \otimes f \cdots f}_{n \text { times }}$. |

## Example 1: Laplace Transform

Compute the Laplace transform for the function

$$
\begin{equation*}
y_{t}=t^{5}+e^{-5 t} \cos (7 t) \tag{8}
\end{equation*}
$$

## Solution

Since the Laplace transform is linear, we can determine the Laplace forms of each part separately and then combine them together. From Table 4, we see that

$$
\begin{equation*}
\mathfrak{L}\left(\frac{t^{n-1}}{(n-1)!}\right)=\frac{1}{s^{n}} \tag{9}
\end{equation*}
$$

Setting $n-1=5$, which is the exponent of $t^{5}$ and noting that we will need to multiply both sides by the factorial $(n-1)$ !, will give a Laplace form of

$$
\begin{equation*}
\frac{5!}{s^{6}} \tag{10}
\end{equation*}
$$

Similarly, for the second term, from Table 4, we have that

$$
\begin{equation*}
\mathfrak{L}\left(e^{-a t} \cos (\omega t)\right)=\frac{s+a}{(s+a)^{2}+\omega^{2}} \tag{11}
\end{equation*}
$$

Comparing with the form that we have, we see that $a=5$ and $\omega=7$. Thus, the Laplace transform is

$$
\begin{equation*}
\frac{s+5}{(s+5)^{2}+7^{2}} \tag{12}
\end{equation*}
$$

Combining the two parts together gives

$$
\begin{equation*}
\mathfrak{L}\left(t^{5}+e^{-5 t} \cos (7 t)\right)=\frac{5!}{s^{6}}+\frac{s+5}{(s+5)^{2}+7^{2}} \tag{13}
\end{equation*}
$$

which is the Laplace transform of $y_{t}$.

## Example 2: Inverse Laplace Transform

Compute the time-domain representation of the following partial-fractioned Laplace function

$$
\begin{equation*}
\frac{5}{10 s+1}+\frac{8}{(2 s+1)^{5}} \tag{14}
\end{equation*}
$$

## Solution

The solution will be found by treating each fraction separately and using the information in Table 5. For the first term, the general form of the transform can be found from Table 5 as

$$
\begin{equation*}
\mathfrak{L}^{-1}\left(\frac{A}{C s+D}\right)=\frac{A}{C} e^{\frac{-D}{C} t} \tag{15}
\end{equation*}
$$

Comparing the general form with our first term, we see that $A=5, C=10$, and $D=1$, which implies that the inverse Laplace form will be

$$
\begin{equation*}
\mathfrak{L}^{-1}\left(\frac{5}{10 s+1}\right)=\frac{5}{10} e^{\frac{-1}{10} t}=0.5 e^{-0.1 t} \tag{16}
\end{equation*}
$$

For the second term, the general form can be written as

$$
\begin{equation*}
\mathfrak{L}^{-1}\left(\frac{A}{(C s+D)^{n}}\right)=\frac{A}{(n-1)!C^{n}} t^{n-1} e^{\frac{-D}{C} t} \tag{17}
\end{equation*}
$$

Comparing the general form with our second term, we see that $A=8, C=2, D=1$, and $n=5$, which implies that the inverse Laplace form will be

$$
\begin{equation*}
\mathfrak{L}^{-1}\left(\frac{8}{(2 s+1)^{5}}\right)=\frac{8}{(5-1)!2^{5}} t^{5-1} e^{\frac{-1}{2} t}=\frac{1}{96} t^{4} e^{-0.5 t} \tag{18}
\end{equation*}
$$

Combining the two terms together gives

$$
\begin{equation*}
\mathfrak{L}^{-1}\left(\frac{5}{10 s+1}+\frac{8}{(2 s+1)^{5}}\right)=0.5 e^{-0.1 t}+\frac{1}{96} t^{4} e^{-0.5 t} \tag{19}
\end{equation*}
$$

This gives us the time-domain representation of the original Laplace function.

## Section 3.1.2: Z-Transform

The $\boldsymbol{z}$-transform converts a discrete function from the time domain into the frequency domain. This transform allows for a simple algebraic solution of complex difference equations. Due to this feature, $z$-transforms are widely used in automation engineering to understand process behaviour and obtain solutions to various control problems.

The $z$-transform is defined as

$$
\begin{equation*}
F(z)=\sum_{n=0}^{\infty} f_{n} z^{-n} \tag{20}
\end{equation*}
$$

where $F$ is the transformed function and $f$ the original discrete function. Conventionally, the $z$ transform is denoted by a script $Z,(\mathrm{U}+1 \mathrm{D} 4 \mathrm{~B} 5)$. Table 6 presents a summary of the most common $z$-transforms. The following are some useful properties:

1) Linearity: $Z(f+g)=Z(f)+Z(g)$.
2) Superposition: $Z(\alpha f)=\alpha Z(f)$.
3) Final Value Theorem: Assuming that the poles (roots of the denominator) lie inside the unit circle, then ${ }^{7}$

$$
\begin{equation*}
\lim _{k \rightarrow \infty} f_{k}=\lim _{z \rightarrow 1}(z-1) F(z) \tag{21}
\end{equation*}
$$

4) Initial Value Theorem: The value in the time domain at the starting point $k=0$ is given by

$$
\begin{equation*}
\lim _{k \rightarrow 0} f_{k}=\lim _{z \rightarrow \infty} z F(z) \tag{22}
\end{equation*}
$$

[^5]Table 6: Table of Common $z$-Transforms ( $T_{s}$ is the sampling time)

| Case | Continuous Time Domain $f(t)$ | Discrete-Time <br> Domain $\left(f\left(k T_{s}\right)\right)$ | Frequency Domain $F(s)$ |
| :---: | :---: | :---: | :---: |
| Dirac Delta, $\delta$ | $\delta(t-a)= \begin{cases}0 & t \neq a \\ \infty & t=a\end{cases}$ | $\delta_{k-a}=\left\{\begin{array}{cc}0 & k \neq a \\ \infty & k=a\end{array}\right.$ | $z^{-a}$ |
| Step Function, <br> $u$ | $u(t)= \begin{cases}0 & t \leq 0 \\ 1 & t>0\end{cases}$ | $u_{k}= \begin{cases}0 & k \leq 0 \\ 1 & k>0\end{cases}$ | $\frac{1}{1-z^{-1}}$ |
| Exponential | $e^{a t}$ | $e^{a k T_{s}}$ | $\frac{1}{1-e^{a T_{s}} z^{-1}}$ |
| Exponential + <br> Cosine | $e^{a t} \cos (\omega t)$ | $e^{a k T_{s}} \cos \left(\omega k T_{s}\right)$ | $\frac{1-e^{a T_{s}} \cos \left(\omega T_{s}\right) z^{-1}}{1-2 e^{a T_{s}} \cos \left(\omega T_{s}\right) z^{-1}+e^{2 a T_{s}} z^{-2}}$ |
| Exponential + <br> Sine | $e^{a t} \sin (\omega t)$ | $e^{a k T_{s}} \sin \left(\omega k T_{s}\right)$ | $\frac{e^{a T_{s}} \sin \left(\omega T_{s}\right) z^{-1}}{1-2 e^{a T_{s}} \cos \left(\omega T_{s}\right) z^{-1}+e^{2 a T_{s}} z^{-2}}$ |
| General Power Series |  | $a^{k}$ | $\frac{1}{1-a z^{-1}}$ |
| First Difference (Derivative) | $\frac{d f}{d t}$ | $f_{k}-f_{k-1}$ | $\left(1-z^{-1}\right) F(z)$ |
| Time Shift (Delay) | $f(t-a) u(t-a)$ | $f_{k-a} u_{k-a}$ | $z^{-a} F(z)$ |

As can be seen from Table 6, many of the forms involve $z^{-1}$. For this reason, automation engineering, it is common to treat $z^{-1}$ as the variable, which is called the backshift operator. Converting between the two representations is rather easy as it involves multiplying by the highest power present in the equation.

The inverse operation of finding the time-domain representation given the $z$-transform is performed using the inverse $z$-transform. Since most automation-engineering examples consider rational functions of $z$, this will require partial fractioning in order to split the rational function into its constituent parts (see Appendix I for details). Once the constituent parts have been obtained, we can then use Table 7 to find the corresponding time-domain representation. Instead of partial-
fractioning, it may be possible to perform long division to obtain the individual values. However, this can be quite long and difficult to do.

Table 7: Useful Inverse $Z$-Transform Table ( $A$ and $D$ can be complex)

| $Z(f(k))$ | $y_{k}$ |
| :---: | :---: |
| $\frac{A z}{C z+D}=\frac{A}{C+D z^{-1}}$ | $y_{k}=\frac{A}{C}\left(-\frac{D}{C}\right)^{k}$ |
| $\frac{A z}{(C z+D)^{n}}$ | $y_{k}=\frac{A}{C^{n}}\left(\frac{\prod_{j=1}^{n-1}(k-j+1)}{\alpha^{n-1}(n-1)!}\right) \alpha^{k}, \alpha=-\frac{D}{C}$ |
| $\frac{A}{\left(C+D z^{-1}\right)^{n}}, n \in \mathbb{Z}$ | $y_{k}=\frac{A}{C^{n}}\binom{k+n-1}{n-1} \alpha^{k}, \alpha=-\frac{D}{C}$ |

Example 3: Z-Transform
Compute the $z$-transform of the following function

$$
\begin{equation*}
y_{k}=5^{k-2}, k \geq 2 \tag{23}
\end{equation*}
$$

## Solution

From Table 6, we see that an exponential function $a^{k}$ has the $z$-transform

$$
\begin{equation*}
\frac{1}{1-a z^{-1}} \tag{24}
\end{equation*}
$$

In our case, this implies that $a=5$. However, we can note that the values are delayed by 2 samples. Therefore, we will need to also use the delay formula from Table 6 to obtain the final result. Thus, the $z$-transform is

$$
\begin{equation*}
\mathcal{Z}\left(5^{k-2}\right)=\frac{z^{-2}}{1-5 z^{-1}} \tag{25}
\end{equation*}
$$

Example 4: Inverse z-Transform
Compute for the following frequency-domain function

$$
\begin{equation*}
\frac{2}{1+5 z^{-1}} \tag{26}
\end{equation*}
$$

the corresponding time-domain representation.

## Solution

From Table 7, we can see that this represents the first case with $A=2, C=1$, and $D=5$. This implies that the time-domain representation is

$$
\begin{equation*}
\mathcal{Z}^{-1}\left(\frac{2}{1+5 z^{-1}}\right)=\frac{A}{C}\left(-\frac{D}{C}\right)^{k}=\frac{2}{1}\left(-\frac{5}{1}\right)^{k}=2(-5)^{k} \tag{27}
\end{equation*}
$$

Finally, we can note that there is a relationship between the Laplace transform of the continuous time domain and the $z$-transform of the discrete time domain. If we set

$$
\begin{equation*}
z=e^{s T} \tag{28}
\end{equation*}
$$

This means that results obtained in the continuous domain will have a transformed representation in the discrete domain. Of note, the imaginary axis of the continuous domain is mapped onto the boundary of the unit circle. This relationship will appear in our further analysis.

## Section 3.2: Time- and Frequency-Based Models

Time- and frequency-based models are one of the most commonly encountered mathematical representations of a process. Time-based models focus on the behaviour of the system with respect to time, that is, how the system evolves or changes over time given specific inputs and states. However, since solutions can only often be obtained by integrating the complex differential equations, recourse is often made to frequency-based models where it can be easier to understand how the process will behave when the inputs change.

Before looking into the different types of time- and frequency-based models, it can be useful to consider some of the terms that can be used to classify the different types of models:

1) Linear versus nonlinear: A model $f(t)$ is said to be linear with respect to $t$ if the following two statements hold:
a. Principle of Superposition or additivity: $f\left(t_{1}+t_{2}\right)=f\left(t_{1}\right)+f\left(t_{2}\right)$, and
b. Principle of Homogeneity: $f\left(\alpha t_{1}\right)=\alpha f\left(t_{1}\right)$.

If these two statements do not hold, then the model is said to be nonlinear. Linearity is a very useful property that allows for a system to be easily analysed using known, wellestablished theoretical methods. Nonlinear models can be linearised, so that in a given region they are well described by the linear equivalent model. A model can be linear with respect to one variable, but not another.
2) Time-invariant versus time-variant: A model is said to be time-invariant if the parameters in the model are constant with respect to time. In a time-varying system, the model parameters can depend on the time period.
3) Lumped parameter versus distributed parameter: A model is said to be a lumped parameter system if the model does not depend on the location, that is, there are no space derivatives present. A model that depends on the location, that is, it contains space derivatives, is called a distributed parameter system. For example, if the temperature $T$ is a function of the $x$ - or $y$-direction, that is, we have a derivative of the form $\frac{\partial^{2} T}{\partial x^{2}}$ in the model, then the model is a distributed-parameter system. If the temperature only depends on the time, then we have a lumped-parameter system. The analysis of distributedparameter systems is often much more difficult than that of a lumped-parameter system. A distributed-parameter system can be reduced to a lumped-parameter model if it is assumed that the variables are homogenously distributed within the space, and hence, all the space derivatives are zero.
4) Causal versus noncausal: A system is said to be causal if the future values only depend on the current and past values. In a noncausal system, the future values depend on the current, past, and future values. A noncausal system is not physically realisable, since the future values will never be known.
5) System with memory (dynamic) versus system without memory (static): A system is said to have memory if the future values depend on both the past and present. On the other hand, a system is said to be without memory or memoryless if future values only depend on the current value.

## Section 3.2.1: Time- and Frequency-Domain Representations

In automation engineering, there are two common representations of a system: state-space and transfer-function models.

The state-space model focuses on the relationship between states, inputs, and outputs in the time domain. The general state-space model is given as

$$
\begin{align*}
\frac{d \vec{x}}{d t} & =\vec{f}(\vec{x}, \vec{u})  \tag{29}\\
\vec{y} & =\vec{g}(\vec{x}, \vec{u})
\end{align*}
$$

where $x$ is the state variable, $u$ is the input variable, $t$ is the time, $y$ is an output, $f$ and $g$ are some functions, and an arrow above denotes a vector. The state variable is a variable that describes the current location of the system from which the system's future behaviour can be determined. A state variable will often appear in some equation as a derivative with respect to time. The input variable, $u$, is a variable that describes the properties of a stream entering a system. Traditionally, the number of inputs is denoted by $m$, the number of states by $n$, and the number of outputs by $p$. A system where $p=m=1$ is said to be univariate or single-input, single-output (SISO). If $p$ and $m$ are greater than 1 , then the system is said to be multivariate or multi-input, multi-output (MIMO). A system is said to be multi-input, single-output (MISO) if $p=1$ and $m>1$.

The general state-space model is often reduced to a linear form

$$
\begin{align*}
\frac{d \vec{x}}{d t} & =\mathcal{A} \vec{x}+\mathcal{B} \vec{u}  \tag{30}\\
\vec{y} & =\mathcal{C} \vec{x}+\mathcal{D} \vec{u}
\end{align*}
$$

where $\mathcal{A}$ is the $n \times n$ state matrix, $\mathscr{B}$ is the $n \times m$ input matrix, $\mathcal{C}$ is the $p \times n$ output matrix, and $\mathcal{D}$ is the $p \times m$ feed-through matrix.

On the other hand, the transfer-function representation focuses solely on the relationship between the inputs and outputs in the Laplace domain and allows easier analysis of the system than for a state-space model. The general transfer function representation can be written as

$$
\begin{equation*}
\vec{Y}(s)=\mathcal{G}(s) \vec{U}(s) \tag{31}
\end{equation*}
$$

where $G$ is a matrix containing the transfer function representation of the model, that is

$$
\vec{Y}(s)=\left[\begin{array}{c}
Y_{1}(s)  \tag{32}\\
\vdots \\
Y_{p}(s)
\end{array}\right], \vec{U}(s)=\left[\begin{array}{c}
U_{1}(s) \\
\vdots \\
U_{m}(s)
\end{array}\right], \mathcal{G}(s)=\left[\begin{array}{ccc}
G_{11}(s) & \cdots & G_{1 m}(s) \\
\vdots & & \vdots \\
G_{p 1}(s) & \cdots & G_{p m}(s)
\end{array}\right]
$$

It is often assumed that each transfer function can be written as

$$
\begin{equation*}
G(s)=\frac{N(s)}{D(s)} e^{-\theta s} \tag{33}
\end{equation*}
$$

where $N$ and $D$ are polynomials of $s$ and $\theta$ is the deadtime or time delay in the system. Time delay arises in real systems due to the transport phenomena, measurement delays, and approximations introduced when linearising a system. The order of a transfer function is equal to the highest power of the $D$-polynomial.

Given the simple form of a transfer-function representation, much of the analysis in control is performed using it.

## Section 3.2.2: Converting Between Representations

The state-space model can be converted to a transfer function model, by applying the Laplace transform to Equation (30) and re-arranging to give ${ }^{8}$

$$
\begin{gather*}
s \vec{X}=\mathcal{A} \vec{X}+\mathcal{B} \vec{U} \\
\vec{Y}=\mathcal{C} \vec{X}+\mathcal{D} \vec{U}  \tag{34}\\
\vec{X}=(s \mathcal{I}-\mathcal{A})^{-1} \mathcal{B} \vec{U}  \tag{35}\\
\vec{Y}=\mathcal{C} \vec{X}+\mathcal{D} \vec{U} \\
G(s)=\frac{\vec{Y}}{\vec{U}}=C(s I-A)^{-1} B+D \tag{36}
\end{gather*}
$$

where $\mathcal{J}$ is the $n \times n$ identity matrix and $s$ is the Laplace transform variable.
Example 5: Numeric Example of Obtaining a Transfer Function
Consider the example

$$
\left\{\begin{array}{l}
5 \frac{d x}{d t}+3 x=u  \tag{37}\\
y=x
\end{array}\right.
$$

Taking the Laplace transform of Equation (37) gives

$$
\left\{\begin{array}{l}
5 s X(s)+3 X(s)=U(s)  \tag{38}\\
Y(s)=X(s)
\end{array}\right.
$$

Substituting $Y$ for $X$ in the first equation of Equation (38) gives

$$
\begin{equation*}
5 s Y(s)+3 Y(s)=U(s) \tag{39}
\end{equation*}
$$

Solving for $Y / U$ gives

[^6]\[

$$
\begin{align*}
(5 s+3) Y(s) & =U(s) \\
\frac{Y(s)}{U(s)} & =\frac{1}{5 s+3} \tag{40}
\end{align*}
$$
\]

Therefore, the transfer function is

$$
\begin{equation*}
G(s)=\frac{Y(s)}{U(s)}=\frac{1}{5 s+3} \tag{41}
\end{equation*}
$$

## Example 6: General Univariate Case

Consider the following $N^{\text {th }}$-order ordinary differential equation

$$
\left\{\begin{align*}
\frac{d^{n} \tilde{x}}{d t^{n}}+a_{n-1} \frac{d^{n-1} \tilde{x}}{d t^{n-1}}+\cdots+a_{0} \tilde{x} & =b_{n-1} \frac{d^{n-1} \tilde{u}}{d t^{n-1}}+\cdots+b_{0} \tilde{u}  \tag{42}\\
\tilde{y} & =\tilde{x}
\end{align*}\right.
$$

such that at $t=0$, all the derivatives are equal to zero and $\tilde{x}=\tilde{y}=0$. Taking the Laplace transform of Equation (42) and simplifying gives

$$
\begin{gather*}
\left\{\begin{array}{c}
s^{n} X(s)+a_{n-1} s^{n-1} X(s)+\cdots+a_{0} X(s)=b_{n-1} s^{n-1} U(s)+\cdots+b_{0} U(s) \\
Y(s)=X(s)
\end{array}\right.  \tag{43}\\
s^{n} Y(s)+a_{n-1} s^{n-1} Y(s)+\cdots+a_{0} Y(s)=b_{n-1} s^{n-1} U(s)+\cdots+b_{0} U(s)
\end{gather*}
$$

Collecting like terms and re-arranging gives

$$
\begin{align*}
& \left(s^{n}+a_{n-1} s^{n-1}+\cdots+a_{0}\right) Y(s)=\left(b_{n-1} s^{n-1}+\cdots+b_{0}\right) U(s) \\
& G(s)=\frac{Y(s)}{U(s)}=\frac{\left(b_{n-1} s^{n-1}+\cdots+b_{0}\right)}{\left(s^{n}+a_{n-1} s^{n-1}+\cdots+a_{0}\right)} \tag{44}
\end{align*}
$$

Example 7: Multivariate Example
Consider the following differential equation

$$
\left\{\begin{align*}
\frac{d \tilde{x}}{d t} & =-a \tilde{x}+b_{1} \tilde{u}_{1}+b_{2} \tilde{u}_{2}  \tag{45}\\
\tilde{y} & =\tilde{x}
\end{align*}\right.
$$

and determine the transfer functions for the system. Note that there are two inputs $(m=2)$ and one output ( $p=1$ ).

Solution

$$
\left.\begin{array}{l}
\left\{\begin{array}{c}
s X(s)=-a X(s)+b_{1} U_{1}(s)+b_{2} U_{2}(s) \\
Y(s)=X(s)
\end{array}\right. \\
s Y(s)+a Y(s)=b_{1} U_{1}(s)+b_{2} U_{2}(s)  \tag{46}\\
(s+a) Y(s)=b_{1} U_{1}(s)+b_{2} U_{2}(s)
\end{array}\right\} \begin{aligned}
& Y(s)=\frac{b_{1}}{s+a} U_{1}(s)+\frac{b_{2}}{s+a} U_{2}(s)
\end{aligned}
$$

Rewriting this into matrix form gives

$$
Y(s)=\left[\begin{array}{ll}
\frac{b_{1}}{s+a} & \frac{b_{2}}{s+a}
\end{array}\right]\left[\begin{array}{l}
U_{1}(s)  \tag{47}\\
U_{2}(s)
\end{array}\right]
$$

Note that this can be written separately, since linear functions are being considered. As well, note that the principle of superposition holds, so that we can study each transfer function separately and then combine the results together.

The reverse operation of converting a transfer function into a state-space representation does not yield a unique solution or realisation. Consider the following transfer function, where $p<n$ :

$$
\begin{equation*}
G(s)=\frac{\beta_{p} s^{p}+\beta_{p-1} s^{p-1}+\cdots+\beta_{1} s+\beta_{0}}{s^{n}+\alpha_{n-1} s^{n-1}+\cdots+\alpha_{1} s+\alpha_{0}} \tag{48}
\end{equation*}
$$

then the controllable canonical realisation is

$$
\begin{align*}
& \mathcal{A}=\left[\begin{array}{ccccc}
-\alpha_{n-1} & -\alpha_{n-2} & \cdots & -\alpha_{1} & -\alpha_{0} \\
1 & 0 & \cdots & 0 & 0 \\
0 & 1 & \ddots & \vdots & \vdots \\
\vdots & \ddots & \ddots & 0 & \vdots \\
0 & \cdots & 0 & 1 & 0
\end{array}\right]_{n \times n} \quad \mathcal{B}=\left[\begin{array}{c}
1 \\
0 \\
\vdots \\
\vdots \\
0
\end{array}\right]_{n \times 1}  \tag{49}\\
& \mathcal{C}=\left[\begin{array}{lllllll}
0 & \cdots & 0 & \beta_{p} & \beta_{p-1} & \cdots & \beta_{1}
\end{array} \beta_{0}\right]_{1 \times n} \mathcal{D}=0_{1 \times 1}
\end{align*}
$$

while the observable canonical realisation is

$$
\begin{align*}
& \mathcal{A}=\left[\begin{array}{ccccc}
-\alpha_{n-1} & -\alpha_{n-2} & \cdots & -\alpha_{1} & -\alpha_{0} \\
1 & 0 & \cdots & 0 & 0 \\
0 & 1 & \ddots & \vdots & \vdots \\
\vdots & \ddots & \ddots & 0 & \vdots \\
0 & \cdots & 0 & 1 & 0
\end{array}\right]_{n \times n} \quad \mathcal{C}=\left[\begin{array}{c}
1 \\
0 \\
\vdots \\
\vdots \\
0
\end{array}\right]_{n \times 1}^{T}  \tag{50}\\
& \mathcal{B}
\end{align*}{\left[\begin{array}{lllllll}
0 & \cdots & 0 & \beta_{p} & \beta_{p-1} & \cdots & \beta_{1}
\end{array} \beta_{0}\right.}_{]_{1 \times n}^{T} \mathcal{D}=0_{1 \times 1}}
$$

Note that for $\mathcal{C}$ in the controllable canonical realisation and for $\mathscr{B}$ in the observable canonical realisation, there will be $n-p-1$ zeroes. The fact that the two forms are related by taking the transpose is not coincidental.

## Example 8: Converting a Transfer Function into its Controllable Canonical Realisation

Convert the transfer function

$$
\begin{equation*}
G(s)=\frac{s-2}{s^{2}-4 s+1} \tag{51}
\end{equation*}
$$

into its controllable canonical realisation.

## Solution

First, we need to make sure that the transfer function is the form given by Equation (48) and determine the values of the parameters. Since the equations have the same form, we note that $n=2$ (highest power in the denominator) and $p=1$ (highest power in the numerator), which implies that the transfer function satisfies the requirements. Thus, we can simply compare the parameters and obtain their values, that is, $\alpha_{2}=1, \alpha_{1}=-4, \alpha_{0}=1, \beta_{1}=1$, and $\beta_{0}=-2$. Thus, using Equation (49), the controllable canonical realisation is

$$
\begin{align*}
& \mathcal{A}=\left[\begin{array}{cc}
4 & -1 \\
1 & 0
\end{array}\right] \mathcal{B}=\left[\begin{array}{l}
1 \\
0
\end{array}\right]  \tag{52}\\
& \mathcal{C}=\left[\begin{array}{ll}
1 & -2
\end{array}\right] \mathcal{D}=0_{1 \times 1}
\end{align*}
$$

Thus, we have defined the four matrices for one of the state-space representations of the given transfer function.

## Section 3.2.3: Discrete-Domain Models

In the discrete domain, the form and types of available models is similar. The linearised state-space representation can be written as

$$
\begin{align*}
\vec{x}_{k+1} & =\mathcal{A}_{d} \vec{x}_{k}+\mathcal{B}_{d} \vec{u}_{k}  \tag{53}\\
\vec{y}_{k} & =\mathcal{C}_{d} \vec{x}_{k}+\mathcal{D}_{d} \vec{u}_{k}
\end{align*}
$$

The subscript $d$ in Equation (53) can be dropped if it is clear that a discrete state-space representation is being considered. The discrete transfer function is the same as the continuous version except that all $s$ are replaced by either $z$ or $z^{-1}$. To convert between the discrete state-space and transfer function representations, Equation (36), can be used, mutatis mutandis.

With discrete transfer functions, it is common to explicitly include the unmeasured disturbance, denoted by $e_{t}$, in the final model. In such cases, the unmeasured disturbance signal is assumed to be a stochastic (random) Gaussian, white noise signal. A Gaussian, white noise signal implies that the values of this signal are normally distributed and independent of past or future values. The general discrete transfer function can be written as

$$
\begin{equation*}
y_{t}=G_{p}\left(z^{-1}, \vec{\theta}\right) u_{t}+G_{l}\left(z^{-1}, \vec{\theta}\right) e_{t} \tag{54}
\end{equation*}
$$

where $G_{p}$ is the process transfer function, $\theta$ the parameters, and $G_{l}$ the disturbance transfer function. Since in most applications, it is assumed that the transfer functions are rational functions of $z^{-1}$, the most common discrete transfer function equation can be written in a form called the prediction error model, which has the following form:

$$
\begin{equation*}
A\left(z^{-1}\right) y_{t}=\frac{B\left(z^{-1}\right)}{F\left(z^{-1}\right)} u_{t-k}+\frac{C\left(z^{-1}\right)}{D\left(z^{-1}\right)} e_{t} \tag{55}
\end{equation*}
$$

where $A\left(z^{-1}\right), C\left(z^{-1}\right), D\left(z^{-1}\right)$, and $F\left(z^{-1}\right)$ are polynomials in $z^{-1}$ of the form

$$
\begin{equation*}
1+\sum_{i=1}^{n_{a}} \theta_{i} z^{-i} \tag{56}
\end{equation*}
$$

where $n_{a}$ is the order of the polynomial and $\theta_{i}$ are the parameters, $B\left(z^{-1}\right)$, is a polynomial in $z^{-1}$ of the form

$$
\begin{equation*}
\sum_{i=1}^{n_{b}} \theta_{i} z^{-i} \tag{57}
\end{equation*}
$$

where $n_{b}$ is the order of the polynomial, and $k$ is the time delay in the system. In general, it is very rare for this system to be used directly. Instead, any of the following simplifications may be used:

1) Box-Jenkins Model: In this model, the $A\left(z^{-1}\right)$ polynomial is ignored. Thus, this model is given as

$$
\begin{equation*}
y_{t}=\frac{B\left(z^{-1}\right)}{F\left(z^{-1}\right)} u_{t-k}+\frac{C\left(z^{-1}\right)}{D\left(z^{-1}\right)} e_{t} \tag{58}
\end{equation*}
$$

In practice, this method is sufficient to fit an accurate model of the system.
2) Autoregressive Moving-Average Exogenous Model (ARMAX): In this model, the $D\left(z^{-1}\right)$ and $F\left(z^{-1}\right)$ polynomials are ignored, which gives

$$
\begin{equation*}
A\left(z^{-1}\right) y_{t}=B\left(z^{-1}\right) u_{t-k}+C\left(z^{-1}\right) e_{t} \tag{59}
\end{equation*}
$$

This model assumes that the denominator for both the input and the error is the same. However, this model has the beneficial property that the estimation of its parameters can be performed using least-squares analysis. A further simplification is to ignore the $C\left(z^{-1}\right)$ term. This gives an autoregressive exogenous model (ARX). This model has the form given by:

$$
\begin{equation*}
A\left(z^{-1}\right) y_{t}=B\left(z^{-1}\right) u_{t-k}+e_{t} \tag{60}
\end{equation*}
$$

3) Output-Error Model (OE): In this model, only the model for the input is fit to the data. The error terms are ignored. Thus, the model is given as

$$
\begin{equation*}
y_{t}=\frac{B\left(z^{-1}\right)}{F\left(z^{-1}\right)} u_{t-k}+e_{t} \tag{61}
\end{equation*}
$$

## Section 3.2.4: Converting Between Discrete and Continuous Models

It is possible to convert between continuous and discrete forms of a model under certain assumption regarding the discretisation. The most common assumption is that the input stays constant between sampling instances, that is, a zero-order hold is used to convert from the continuous (analogue) to discrete (digital) domains. In such case, it can be shown that the continuous state-space representation can be converted into the discrete time using the following formulae:

$$
\begin{align*}
& \mathcal{A}_{d}=e^{A \tau_{s}} \\
& \mathcal{B}_{d}=\left(\int_{\tau=0}^{\tau=\tau_{s}} e^{A \tau} d \tau\right) B  \tag{62}\\
& \mathcal{C}_{d}=C \\
& \mathcal{D}_{d}=D
\end{align*}
$$

where $\tau_{s}$ is the sampling time. It should be noted that all exponentiation is matrixwise exponentiation. The corresponding transfer function can then be obtained using Equation (36). For a simple, first order, continuous transfer function given as

$$
\begin{equation*}
Y(s)=\frac{K}{\tau_{p} s+1} U(s) \tag{63}
\end{equation*}
$$

the discrete, transfer function is given as

$$
\begin{equation*}
y_{k}=\frac{K\left(1-\mathrm{e}^{-\frac{\tau_{p}}{\tau_{s}}}\right)}{1-\left(1-\mathrm{e}^{-\frac{\tau_{p}}{\tau_{s}}}\right) z^{-1}} u_{k} \tag{64}
\end{equation*}
$$

If a system has time delay, then it can be converted using the following formula

$$
\begin{equation*}
k=\left\lfloor\frac{\theta}{\tau_{s}}\right\rfloor \tag{65}
\end{equation*}
$$

where $\theta$ is the continuous time delay and $\lfloor\bullet\rfloor$ is an appropriately selected rounding function converts to an integer value.

The general rule of thumb for selecting the sampling time is that

$$
\begin{equation*}
\tau_{s}=(0.1 \text { to } 0.2) \tau_{p}^{(\min )} \tag{66}
\end{equation*}
$$

where $\tau_{p}^{(\min )}$ is the smallest time constant present in the system.

## Section 3.2.5: Impulse Response Model

In the discrete domain, the infinite impulse response model is defined as

$$
\begin{equation*}
y_{k}=\sum_{i=0}^{\infty} h_{i} z^{-i} u_{k}=\sum_{i=0}^{\infty} h_{i} u_{k-i} \tag{67}
\end{equation*}
$$

where $h$ are the impulse coefficients. The values of $h$ can be obtained by either performing long division on the transfer function or partial fractioning the result to obtain the impulse coefficients. Since the values of $h$ can often quickly taper off, it is possible to convert the infinite impulse response model into a finite impulse response (FIR) model, that is,

$$
\begin{equation*}
y_{k}=\sum_{i=0}^{n} h_{i} z^{-i} u_{k}=\sum_{i=0}^{n} h_{i} u_{k-i} \tag{68}
\end{equation*}
$$

where $n$ is an integer representing the number of terms selected for the model.

## Section 3.2.6: Compact State-Space Representation

In many theoretical applications, the state-space representation is often written in a compact manner that resembles a transfer-function formulism but using matrices. For the standard statespace representation given by Equation (30), the compact or block state-space representation can be written as

$$
G=\left[\begin{array}{l|l}
\mathcal{A} & \mathcal{B}  \tag{69}\\
\hline \mathcal{C} & \mathcal{D}
\end{array}\right]
$$

Using this representation, it is possible to easily partition the matrices and show this in a compact manner, for example,

$$
G=\left[\begin{array}{cc|c}
\mathcal{A}_{11} & \mathcal{A}_{12} & \mathcal{B}_{1}  \tag{70}\\
\mathcal{A}_{21} & \mathcal{A}_{22} & \mathcal{B}_{2} \\
\hline \mathcal{C}_{1} & \mathcal{C}_{2} & \mathcal{D}
\end{array}\right]
$$

where the states have been split into four groups and the relationships between the different groups can easily be shown.

When using the compact representation, various short cuts can be used when combining two models. When adding two compact representations, that is the transfer functions are in parallel,

$$
G_{1}+G_{2}=\left[\begin{array}{cc|c}
\mathcal{A}_{1} & 0 & \mathcal{B}_{1}  \tag{71}\\
0 & \mathcal{A}_{2} & \mathcal{B}_{2} \\
\hline \mathcal{C}_{1} & \mathcal{C}_{2} & \mathcal{D}_{1}+\mathcal{D}_{2}
\end{array}\right]
$$

When multiplying two representations, that is the transfer functions are in series,

$$
G_{1} G_{2}=\left[\begin{array}{cc|c}
\mathcal{A}_{1} & \mathcal{B}_{1} \mathcal{C}_{2} & \mathcal{B}_{1} \mathcal{D}_{2}  \tag{72}\\
0 & \mathcal{A}_{2} & \mathcal{B}_{2} \\
\hline \mathcal{C}_{1} & \mathcal{D}_{1} \mathcal{C}_{2} & \mathcal{D}_{1} \mathcal{D}_{2}
\end{array}\right]
$$

The inverse of a compact state-space representation can be written as

$$
G^{-1}=\left[\begin{array}{c|c}
\mathcal{A}-\mathcal{B D}^{-1} \mathcal{C} & -\mathcal{B D}^{-1}  \tag{73}\\
\hline \mathcal{D}^{-1} \mathcal{C} & \mathcal{D}^{-1}
\end{array}\right]
$$

provided that $D$ is invertible. Finally, the transpose of this representation can be written as

$$
G^{T}=\left[\begin{array}{c|c}
\mathcal{A}^{T} & \mathcal{C}^{T}  \tag{74}\\
\hline \mathcal{B}^{T} & \mathcal{D}^{T}
\end{array}\right]
$$

A common operation with state-space models is to transform the order or meaning of the states. In such cases, the compact representation can be derived as follows. Assume that

$$
\begin{align*}
& \tilde{x}=\mathcal{T} x \\
& \tilde{u}=\mathcal{S} u  \tag{75}\\
& \tilde{y}=\mathcal{R} y
\end{align*}
$$

where $T, R$, and $S$ are appropriately sized nonsingular (invertible) matrices. In this case, the transformed compact representation can be written as

$$
\tilde{G}=\left[\begin{array}{c|c}
\mathcal{T} \mathcal{A T}^{-1} & \mathcal{T B S}^{-1}  \tag{76}\\
\mathcal{R C \mathcal { T } ^ { - 1 }} & \mathcal{R D S}^{-1}
\end{array}\right]
$$

## Section 3.3: Process Analysis

Having considered the different models, it would be useful to understand what information can be easily extracted from such models that would allow us to understand how the process behaves. Since it is easier to extract information from transfer functions, the focus will be on analysing them. It will be assumed that the transfer function can be written in the continuous domain as

$$
\begin{equation*}
G(s)=\frac{N(s)}{D(s)} e^{-\theta s} \tag{77}
\end{equation*}
$$

or in the discrete domain as

$$
\begin{equation*}
G(z)=\frac{N(z)}{D(z)} z^{-k} \tag{78}
\end{equation*}
$$

The order of a polynomial, denoted by $n$, is defined as the highest power present in a given polynomial, for example, $x^{4}+x^{2}$ is a fourth-order polynomial, since the highest power is 4 . The order of a transfer function is equal to the order of the denominator polynomial, $D$. A transfer function is said to be proper if the order of the numerator is less than or equal to the order of the denominator, that is, $n_{N} \leq n_{D}$. A transfer function is said to be strictly proper if the order of the numerator is less than the order of the denominator, that is, $n_{N}<n_{D}$. For most physical processes, a transfer function will be strictly proper.

A process is said to be causal if the time delay, $\theta$ (or $k$ ) is nonnegative; otherwise, the process is noncausal. In the discrete domain, causality is ensured by a proper transfer function. A
causal process does not depend on unknown future values. Again, all physical processes must be causal or else they would be able to predict the future.

The poles of a transfer function are defined as the roots of the denominator, $D$, that is, those values of $s($ or $z)$ such that $D=0$. The zeros of a transfer function are defined as the roots of the numerator, $N$.

A process is said to be at steady state if all derivatives are equal to zero, that is,

$$
\begin{equation*}
\frac{d \vec{x}}{d t}=0 \tag{79}
\end{equation*}
$$

If Equation (79) does not hold, then the process is said to be operating in transient mode. It should be noted that small deviations from zero do not necessarily mean that the process is now in a transient mode. In practice, it may not be possible to achieve an exact steady state, where all the derivatives are exactly equal to zero, both due to continual small fluctuations and measurement imprecisions. In such cases, it is common to define the settling time, $t_{s}$, of a process as the time it takes for the process to reach and stay within an envelope centred on the new steady-state value and has bounds that are $5 \%$ of the change in the process variable. Figure 23 shows how the settling time is computed.


Figure 23: Determining the Settling Time

For a process modelled by a continuous transfer function, there are three key parameters of interest: gain, time constant, and time delay (deadtime or process delay). The gain, $\boldsymbol{K}$, of the process represents the steady-state behaviour of the system for a unit change in the input. It is defined as

$$
\begin{equation*}
K=\frac{y(t \rightarrow \infty)}{u(t \rightarrow \infty)} \tag{80}
\end{equation*}
$$

where it is assumed that the input is bounded as $t \rightarrow \infty$. The gain is the same irrespective of the bounded input used. Using the final value theorem for a general transfer function and a step change in the input, it can be seen that

$$
\begin{align*}
K & =\lim _{s \rightarrow 0} s Y(s)=\lim _{s \rightarrow 0} \frac{s G(s)}{s}=\lim _{s \rightarrow 0} G(s) \\
& =\lim _{s \rightarrow 0} \frac{\left(b_{n-1} s^{n-1}+\cdots+b_{0}\right)}{\left(s^{n}+a_{n-1} s^{n-1}+\cdots+a_{0}\right)} e^{-\theta s}  \tag{81}\\
& =\frac{b_{0}}{a_{0}}
\end{align*}
$$

The ratio $b_{0} / a_{0}$ is called the gain, $K$, of the process.
The process time constant represents the transient or dynamic component of the system, that is, how quickly the system responds to changes in the input and reaches a new steady-state value. The larger the time constant the longer it takes to reach steady state. It can be obtained from the transfer function by factoring the denominator into the form

$$
\begin{equation*}
\frac{\left(b_{n-1} s^{n-1}+\cdots+b_{0}\right)}{\left(s^{n}+a_{n-1} s^{n-1}+\cdots+a_{0}\right)}=\frac{\left(b_{n-1} s^{n-1}+\cdots+b_{0}\right)}{\prod_{i=1}^{n}\left(\tau_{i} s+1\right)} \tag{82}
\end{equation*}
$$

which gives the time constants, $\tau$, of the process. If there are multiple time constants, then the largest time constant will determine the overall process response speed.

The final parameter of interest is the time delay, $\theta$, which measures how long it takes before the system responds to a change in the input. Time delays can arise from two different sources: physical and measurement. Physically, time delays are perceived times at which a system is not responding to changes in the input. They can arise if it takes time for an observable change to be seen, for example, heating a large tank would take some time before the temperature rises by an appreciable amount that can be measured. On the other hand, measurement time delays arise
due to the placement of sensors. In many systems it may not be possible to measure the variable immediately where it will act on the process. In this case, it will take some time before the variable actually affects the system, for example, the flow rate in a pipe could be measured at the beginning of the pipe and it could take some time before it will reach the process. In the frequency domain, the time delay is found as

$$
\begin{equation*}
e^{-\theta s} \tag{83}
\end{equation*}
$$

where $\theta$ is the time delay. In many applications, it may be necessary to convert the exponential form of the time delay into a polynomial expansion. This conversion can be accomplished using the Padé approximation, denoted as the $n / m$ Padé Approximation, where $n$ is the order of the polynomial in the numerator and $m$ is the order of the polynomial in the denominator. The $\mathbf{1} / \mathbf{1}$ Padé approximation is given as

$$
\begin{equation*}
e^{-\theta s}=\frac{1-\frac{\theta}{2} s}{1+\frac{\theta}{2} s} \tag{84}
\end{equation*}
$$

The $\mathbf{2} / \mathbf{2}$ Padé approximation is given as

$$
\begin{equation*}
e^{-\theta s}=\frac{1-\frac{\theta}{2} s+\frac{\theta}{12} s^{2}}{1+\frac{\theta}{2} s+\frac{\theta}{12} s^{2}} \tag{85}
\end{equation*}
$$

A table of the Pade Approximation for the exponential function up to $3 / 3$ are given in Table 8, from which the Padé Approximation for the deadtime can be easily derived.

Table 8: Padé Approximations for the Exponential Function, $\mathrm{e}^{\mathrm{z}}$

| $n$ <br> $m \downarrow$ | 0 | 1 | 2 | 3 |
| :--- | :--- | :--- | :--- | :--- |
| 0 | $\frac{1}{1}$ | $\frac{1}{1-z}$ | $\frac{1}{1-z+\frac{1}{2} z^{2}}$ | $\frac{1}{1-z+\frac{1}{2} z^{2}-\frac{1}{6} z^{3}}$ |
| 1 | $\frac{1+z}{1}$ | $\frac{1+\frac{1}{2} z}{1-\frac{1}{2} z}$ | $\frac{1+\frac{1}{3} z}{1-\frac{2}{3} z+\frac{1}{6} z^{2}}$ | $\frac{1+\frac{1}{4} z}{1-\frac{3}{4} z+\frac{1}{4} z^{2}-\frac{1}{24} z^{3}}$ |


| $n$ <br> $m \downarrow$ | 0 | 1 | 2 | 3 |
| :--- | :--- | :--- | :--- | :--- |
| 2 | $\frac{1+z+\frac{1}{2} z^{2}}{1}$ | $\frac{1+\frac{2}{3} z+\frac{1}{6} z^{2}}{1-\frac{1}{3} z}$ | $\frac{1+\frac{1}{2} z+\frac{1}{12} z^{2}}{1-\frac{1}{2} z+\frac{1}{12} z^{2}}$ | $\frac{1+\frac{2}{5} z+\frac{1}{20} z^{2}}{1-\frac{3}{5} z+\frac{3}{20} z^{2}-\frac{1}{60} z^{3}}$ |
| 3 | $\frac{1+z+\frac{1}{2} z^{2}+\frac{1}{6} z^{3}}{1}$ | $\frac{1+\frac{3}{4} z+\frac{1}{4} z^{2}+\frac{1}{24} z^{3}}{1-\frac{1}{4} z}$ | $\frac{1+\frac{3}{5} z+\frac{3}{20} z^{2}+\frac{1}{60} z^{3}}{1-\frac{2}{5} z+\frac{1}{20} z^{2}}$ | $\frac{1+\frac{1}{2} z+\frac{1}{10} z^{2}+\frac{1}{120} z^{3}}{1-\frac{1}{2} z+\frac{1}{10} z^{2}-\frac{1}{120} z^{3}}$ |

## Example 9: Extracting Information from a Transfer Function

Determine the gain, time constant, and time delay for the following transfer function

$$
\begin{equation*}
G(s)=\frac{Y(s)}{U(s)}=\frac{1}{5 s+3} e^{-5 s} \tag{86}
\end{equation*}
$$

## Solution

Re-arranging into the required form gives

$$
\begin{equation*}
G(s)=\frac{Y(s)}{U(s)}=\frac{1 / 3}{\frac{5}{3} s+1} e^{-5 s} \tag{87}
\end{equation*}
$$

from which, by inspection, the gain is $1 / 3$ and the time constant is $5 / 3$. The time delay is 5 .

## Section 3.3.1: Frequency-Domain Analysis

In frequency-domain analysis, the transfer function is solely used and analysed to determine its behaviour at different frequencies. Frequency-domain analysis is often presented graphically and was once used extensively before the advent of computers. Understanding the principles of frequency-domain analysis is still important, especially when designing various filters for electronic systems, such as microphones.

Frequency-domain methods are based on setting $s=j \omega$, where $j$ is the imaginary number $\sqrt{1}$ and $\omega$ is the frequency. This basically means that we are considering the response of the system to sinusoidal waves with different frequencies. The goal is to see how the process changes the amplitude or strength and the phase shift, that is, assuming that the original input has the form $\sin \omega t$
the response will have the form

$$
\begin{equation*}
A \sin (\omega t+\phi) \tag{89}
\end{equation*}
$$

where $A$ is the amplitude and $\phi$ is the phase shift. Consider a general transfer function, $G(s)$, then define the amplitude ratio, AR, (which corresponds to a normalised amplitude) as

$$
\begin{equation*}
\mathrm{AR}=\|G(j \omega)\|=\sqrt{\operatorname{Re}(G(j \omega))^{2}+\operatorname{Im}(G(j \omega))^{2}} \tag{90}
\end{equation*}
$$

where $R e$ is the real component of a complex number, Im is the imaginary component, and $\|\cdot\|$ is the modulus function. It should be noted that in many applications, the logarithm of the amplitude ratio is used. In most cases, a base-10 logarithm is used. Occasionally, a decibel logarithm will be used, that is, $\mathrm{AR}=20 \log \|G(j \omega)\|$. The phase angle, $\phi$, conventionally expressed in degrees, is defined as

$$
\begin{equation*}
\phi=\arctan \left(\frac{\operatorname{Im}(G(j \omega))}{\operatorname{Re}(G(j \omega))}\right) \tag{91}
\end{equation*}
$$

where $\arctan$ is the standard inverse tangent function defined on $]-90^{\circ}, 90^{\circ}$. It should be noted that if the denominator is negative, then it is necessary to add $180^{\circ}$ to the value obtained previously. This is because the arctan function does not work on the complete circle. ${ }^{10}$

It can be noted that if the original transfer function is a product of simpler transfer functions, that is,

$$
\begin{equation*}
G(s)=\prod G_{i}(s) \tag{92}
\end{equation*}
$$

then the amplitude ratio can be computed as

$$
\begin{equation*}
\mathrm{AR}=\prod \mathrm{AR}_{G_{i}(s)} \tag{93}
\end{equation*}
$$

which is the product of the individual amplitude ratios corresponding to each simpler transfer function. The phase angle can be computed as

$$
\begin{equation*}
\phi=\sum \phi_{G_{i}(s)} \tag{94}
\end{equation*}
$$

These two formulae can simplify the determination of the amplitude ratio and phase angles of complex transfer functions. This formulae result from the fact that all transfer functions can be

[^7]expressed in polar form as $G(j \omega)=\|G(j \omega)\| \mathrm{e}^{-\phi j}$, from which it is easy to see that combining multiple transfer functions as given by Equation (92), the given formulae will result.

The amplitude ratio and phase angle can be graphically presented in two different forms, either as a Bode plot or as a Nyquist plot. In a Bode plot, the $x$-axis is the frequency, $\omega$, and the $y$-axis is either the amplitude ratio or phase angle. A typical Bode plot is shown in Figure 24.


Figure 24: Bode Plot: The Argument-Ratio and Phase-Angle Plots
In a Nyquist plot, the real component of $G(j \omega)$ is plotted on the $x$-axis and the imaginary component of $G(j \omega)$ is plotted on the $y$-axis. A typical Nyquist plot is shown in Figure 25. Both graphs give similar information. In most applications, Bode plots are used, but a Nyquist plot can be useful to analysis the interactions between different systems.


Figure 25: A Nyquist Plot (for the same process as given by Bode plot in Figure 24)

## Section 3.3.2: Stability

Stability of a model refers to the behaviour of the model as $t \rightarrow \infty$ for the case where the input is bounded, that is, $\left|u_{t}\right| \leq L$, where $L$ is a constant. A model is said to be stable if as $t \rightarrow \infty$ for a bounded input, $y_{\infty} \rightarrow K$, where $K$ is a constant. Otherwise, a model is said to be unstable. Determining the stability of a process depends on the type of representation used and nature of the time domain. Table 9 summarises the key results regarding stability. A detailed examination of how the different types of stability manifest themselves can be found in Section 3.3.4.6.

For a state-space representation, stability is determined by examining the eigenvalues of the state matrix, $\mathcal{A}$. In the continuous time domain, the real part of all the eigenvalues must be less than zero, for the model to be stable; otherwise, the model is said to be unstable. In the discrete time domain, the magnitude of all the eigenvalues must be less than 1 , that is, the eigenvalues must lie inside the unit circle, for the process to be stable; otherwise, the model is said to be unstable.

For a transfer-function representation, stability is determined by examining the poles, or roots of the denominator, $D$. In the continuous time domain, the real part of all the poles must be less than zero, for the model to be stable; otherwise, the model is unstable. In the discrete time
domain, the magnitude of all the poles must be less than 1, that is, the poles must lie inside the unit circle, for the process to be stable; otherwise, the process is unstable. ${ }^{11}$

Finally, it can be noted that if the state-space representation is stable, then so will the transfer function representation be. However, the converse is not true, since each transfer function does not have a unique state-space representation and it is possible to construct an unstable statespace representation for a stable transfer function, by adding an unobservable unstable state.

Table 9: Summary of the Stability Conditions for Different Representations and Time Domains

| Representation |  | Analysis Metric, $p$ | Stable | Unstable |
| :---: | :---: | :---: | :---: | :---: |
|  | State-Space | Eigenvalues of the statespace matrix, $\mathcal{A}$ | $\operatorname{Re}(p)<0$ | $\operatorname{Re}(p) \geq 0$ |
|  | Transfer <br> Function | Poles of the transfer function (Roots of $D$ ) | $\operatorname{Re}(p)<0$ | $\operatorname{Re}(p) \geq 0$ |
| 河 | State-Space | Eigenvalues of the statespace matrix, $\mathcal{A}$ | $\\|p\\|<1$ | $\\|p\\| \geq 1$ |
| $\begin{aligned} & 0 \\ & 0 \\ & 0 \\ & 0 \\ & 0 \end{aligned}$ | Transfer <br> Function | Poles of the transfer function (Roots of $D$ ) | $\\|p\\|<1$ | $\\|p\\| \geq 1$ |

## Example 10: Determining the Stability of a Transfer Function

Determine if the following processes are stable:

$$
\begin{gather*}
G_{1}(s)=\frac{1}{5 s+3} e^{-5 s}  \tag{95}\\
G_{2}(s)=\frac{1}{(5 s+3)\left(s^{2}+4\right)} e^{-5 s}  \tag{96}\\
G_{3}(z)=\frac{z^{2}}{z^{4}+z^{3}+z^{2}+z-1} \tag{97}
\end{gather*}
$$

[^8]
## Solution

For $G_{1}$, setting $5 s+3$ equal to zero and solving, gives a pole of $-3 / 5$. According to the results in Table 9 for a continuous transfer function, the system is stable, since the pole is less than zero.

For $G_{2}$, setting the two terms to zero and solving, gives three poles of $-3 / 5$ and $\pm 2 i$. From Table 9 for a continuous transfer function, we conclude that the system is unstable since the we have two poles with a real component equal to zero and one pole whose value is less than zero.

For $G_{3}$, the setting the denominator equal to zero and solving, gives poles of -1.291 , $-0.1141 \pm 1.217 i$, and 0.519 . For a discrete transfer function, Table 9 states that in order for the system to be stable, the magnitude of the roots must be less than one. We see that we have at least one pole ( -1.291 ) whose magnitude will be greater than zero. This implies that the system will be unstable.

## Example 11: Determining the Stability of a State-Space Model

Determine if the given $\mathcal{A}$-matrices represent stable continuous processes:

$$
\begin{gather*}
\mathcal{A}_{1}=\left[\begin{array}{cc}
2 & 0 \\
3 & -2
\end{array}\right]  \tag{98}\\
\mathcal{A}_{2}=\left[\begin{array}{ccc}
-5 & 5 & 6 \\
0 & -3 & 2 \\
0 & 0 & -1
\end{array}\right]  \tag{99}\\
\mathcal{A}_{3}=\left[\begin{array}{cc}
2 & 1 \\
-1 & 2
\end{array}\right] \tag{100}
\end{gather*}
$$

Solution
For $\mathcal{A}_{1}$, since it is a triangular matrix, the eigenvalues can be directly found by looking at the main diagonal entries. This implies that the eigenvalues are 2 and -2 . According to the results in Table 9 for a continuous state-space model, the system is unstable, since one of the eigenvalues is greater than zero.

Likewise, for $\mathcal{A}_{2}$, we can find the eigenvalues by looking at the main diagonal entries. This gives $-5,-3$, and -1 . Since all of the eigenvalues are less than zero, from Table 9 for a continuous state-space model, we conclude that the system is stable.

For $\mathcal{A}_{3}$, we need to compute the eigenvalues as follows

$$
\begin{align*}
& \operatorname{det}\left(\lambda I-\mathcal{A}_{3}\right)=(\lambda-2)^{2}-(1)(-1) \\
& =\lambda^{2}-4 \lambda+5  \tag{101}\\
& =2 \pm i
\end{align*}
$$

Since the real part of the eigenvalues is greater than zero, from Table 9 that states that in order for the system to be stable, the real part of roots must be less than zero, we conclude that the system is unstable.

## Section 3.3.2.1: Routh Stability Analysis

Although the results presented in Table 9 are very useful when a numerical transfer function is available, determining the roots algebraically can be difficult, if not impossible for higher-order systems. For this reason, there exist special tests for stability that do not require the roots of the polynomial to be determined. One of the most common tests for a continuous-time transfer function is the Routh stability analysis. By simply considering the coefficients of the denominator polynomial, the stability of the system can be determined. Consider the characteristic polynomial (the polynomial of the denominator) as

$$
\begin{equation*}
a_{n} s^{n}+a_{n-1} s^{n-1}+a_{n-2} s^{n-2}+\ldots+a_{1} s+a_{0}=0 \tag{102}
\end{equation*}
$$

There are two conditions for the Routh stability test:

1) All coefficients must be present and have the same sign (either all strictly positive, that is, greater than zero or all strictly negative, that is less than zero). If this is not the case, the system is unstable.
2) If condition 1 is satisfied, then create the table shown in Table 10.
a) In the first row, place the coefficients corresponding to $a_{n}, a_{n-2}, \ldots$
b) In the second row, place the coefficients corresponding to $a_{n-1}, a_{n-3}, \ldots$
c) Using the values from the rows above and the current column and the column to the right, compute a determinant-like number using the formula provided in the table. For the $i^{\text {th }}$ row in the $j^{\text {th }}$ column $(i \in\{1,2, \ldots, n+1, j \in\{1,2, \ldots,\lceil 0.5 n\rceil\})$, the general formula is

$$
\begin{equation*}
\lambda_{i}^{(j)}=\frac{\lambda_{i+1}^{(j-2)} \lambda_{i}^{(j-1)}-\lambda_{i}^{(j-2)} \lambda_{i+1}^{(j-1)}}{\lambda_{1}^{j-1}} \tag{103}
\end{equation*}
$$

where $\lambda_{i}^{(1)}=a_{n-2 i+2}$ and $\lambda_{i}^{(2)}=a_{n-2 i+1}$ (the first two rows are the coefficients of the polynomial of interest). Any values that are not given are assumed to be zero.
d) Continue computing this value until the table has $n+1$ rows.
e) A polynomial is stable (all roots of the characteristics polynomial are negative) if all the values in the first column have the same sign, either strictly positive or strictly negative.

Table 10: Routh Stability Analysis Table

| Row | $\mathbf{1}$ | $\mathbf{2}$ | $\mathbf{3}$ | $\boldsymbol{\cdots}$ |
| :---: | :---: | :---: | :---: | :---: |
| $\mathbf{1}$ | $a_{n}$ | $a_{n-2}$ | $a_{n-4}$ | $\cdots$ |
| $\mathbf{2}$ | $a_{n-1}$ | $a_{n-3}$ | $a_{n-5}$ | $\cdots$ |
| $\mathbf{3}$ | $b_{1}=\frac{a_{n-1} a_{n-2}-a_{n} a_{n-3}}{a_{n-1}}$ | $b_{2}=\frac{a_{n-1} a_{n-4}-a_{n} a_{n-5}}{a_{n-1}}$ | $\ldots$ |  |
| $\mathbf{4}$ | $c_{1}=\frac{b_{1} a_{n-3}-b_{2} a_{n-1}}{b_{1}}$ | $\ldots$ |  |  |
| $\boldsymbol{n}+\mathbf{1}$ | $\vdots$ | $\vdots$ |  |  |
|  | $z_{1}$ |  |  |  |

## Example 12: Example of Routh Stability Analysis

Using Routh stability, determine the stability of the following two transfer functions:

1) $G(s)=\frac{5}{s^{6}+5 s^{4}-6 s^{3}+2 s^{2}+3 s+1}$
2) $G(s)=\frac{4}{s^{4}+3 s^{3}+s^{2}+2 s+1}$

## Solution

For the first transfer function, we can determine the stability by inspection. Since we are missing one of the powers $\left(s^{5}\right)$ and the coefficients are both negative and positive, we can conclude on the basis of condition (1) of the Routh stability analysis that this transfer function is unstable.

For the second transfer function, we see that condition (1) is satisfied, as all the powers are present and the coefficients have the same sign. Therefore, we will need to check condition (2) by constructing the Routh array. There will be 3 (always $n / 2$, rounded up) columns in the array. In the first row (denoted as $\mathbf{1}$ in Table 11), we place the even coefficients ordered in decreasing
powers. In the second row, we place the odd coefficients. In Rows 3, 4, and 5, we compute the values using the formula:

$$
\begin{gathered}
b_{1}=\frac{a_{n-1} a_{n-2}-a_{n} a_{n-3}}{a_{n-1}}=\frac{3(1)-1(2)}{3}=\frac{1}{3} \\
b_{2}=\frac{a_{n-1} a_{n-4}-a_{n} a_{n-5}}{a_{n-1}}=\frac{3(1)-1(0)}{3}=1 \\
c_{1}=\frac{(1 / 3) 2-3(1)}{1 / 3}=-7 \\
z_{1}=\frac{-7(1)-(1 / 3)(0)}{-7}=1
\end{gathered}
$$

It can be noted that any values that do not exist (such as $a_{n-5}$ ) are equal to zero. As stated by the definition of the condition, we will always have $n+1$ rows. The formulae used to compute the subsequent values are essentially determinants with the sign in the numerator flipped.

Table 11: Routh Array

| Row | $\mathbf{1}$ | $\mathbf{2}$ | $\mathbf{3}$ |
| :---: | :---: | :---: | :---: |
| $\mathbf{1}$ | 1 | 1 | 1 |
| $\mathbf{2}$ | 3 | 2 |  |
| $\mathbf{3}$ | $1 / 3$ | 1 |  |
| $\mathbf{4}$ | -7 |  |  |
| $\mathbf{5}$ | 1 |  |  |

From Table 11, we see that there is a sign change in Column 1, row 4, which implies that the system is unstable.

## Section 3.3.2.2: Jury Stability Analysis

Although the results presented in Table 9 are very useful when a numerical transfer function is available, determining the roots algebraically can be difficult, if not impossible for higher-order systems. For this reason, there exist special tests for stability that do not require the roots of the polynomial to be determined. One of the most common tests for a discrete-time transfer function is the Jury Stability Analysis. By simply considering the coefficients of the denominator
polynomial, the stability of the system can be determined. Like with the Routh stability test, a table is constructed. Consider that the polynomial of interest is given as

$$
\begin{equation*}
D(z)=a_{n} z^{n}+a_{n-1} z^{n-1}+\cdots+a_{1} z+a_{0} \tag{104}
\end{equation*}
$$

Construct the following table (similar to that shown in Table 12), where the first row contains all the coefficients starting from $a_{0}$ and ending with $a_{n}$. The second row is the first row reversed, that is, the first column contains $a_{n}$, and the last column contains $a_{0}$. The third row is computed using

$$
\begin{equation*}
b_{i}=a_{0} a_{i}-a_{n} a_{n-i} \tag{105}
\end{equation*}
$$

while the fourth row is the third row reversed. Note that only the first $n$ values are reversed. Effectively, we have now created a new reduced polynomial that needs to be analysed. Thus, the fifth row is computed using the same formula as for the third row given by Equation (105), but replacing $n$ by $n-1$ coefficients, that is,

$$
\begin{equation*}
c_{i}=b_{0} b_{i}-b_{n-1} b_{n-1-i} \tag{106}
\end{equation*}
$$

This procedure is then repeated until there is row with only 3 elements left (denoted as $q_{0}, q_{1}$, and $q_{2}$, that is, the table will have $2 n-3$ rows. In general, for the $(2 j+1)^{\text {th }}$ row $(j=1,2, \ldots, n-2)$, that is, each odd-numbered row, we can write the relationship as follows

$$
\begin{equation*}
d_{i}^{(2 j+1)}=d_{0}^{(2 j-1)} d_{i}^{(2 j-1)}-d_{n-j+1}^{(2 j-1)} d_{n-j-i+1}^{(2 j-1)} \tag{107}
\end{equation*}
$$

where $d_{i}^{(1)}=a_{i}$, that is, the first row contains the actual polynomial coefficients. The $(2 j+2)^{\text {th }}$ row, that is, the even row, will then be the $n-j+1$ coefficients written in reversed order. It can be noted that the odd rows are essentially a determinant of the above two rows.

The conditions for stability can then be stated as:

1) $D(1)>0$
2) $(-1)^{n} D(-1)>0$
3) $\left|a_{0}\right|>\left|a_{n}\right|$
4) $\left|b_{0}\right|>\left|b_{n-1}\right|$, and continuing in this manner until the last row is reach, where $\left|q_{0}\right|>\left|q_{2}\right|$.

If any of the above conditions fail, then the system will have at least one pole outside the unit circle and hence be unstable.

Table 12: Table for the Jury Stability Analysis

| Row | $\mathbf{0}$ | $\mathbf{1}$ | $\cdots$ | $n-1$ | $n$ |
| :---: | :---: | :---: | :---: | :---: | :---: |


| $\mathbf{1}$ | $a_{0}$ | $a_{1}$ | $\ldots$ | $a_{n-1}$ | $a_{n}$ |
| :---: | :---: | :---: | :---: | :---: | :---: |
| $\mathbf{2}$ | $a_{n}$ | $a_{n-1}$ | $\ldots$ | $a_{1}$ | $a_{0}$ |
| $\mathbf{3}$ | $b_{0}=a_{0} a_{0}-a_{n} a_{n}$ | $b_{1}=a_{0} a_{1}-a_{n} a_{n-1}$ | $\ldots$ | $b_{n-1}=a_{0} a_{n-1}-a_{n} a_{1}$ | 0 |
| $\mathbf{4}$ | $b_{n-1}$ | $b_{n-2}$ | $\ldots$ | $b_{0}$ | 0 |
| $\mathbf{5}$ |  |  |  |  |  |
| $\vdots$ | $\vdots$ | $\vdots$ | $\vdots$ | $\vdots$ | $\vdots$ |
| $\mathbf{2 n - 3}$ | $\boldsymbol{q}_{\mathbf{0}}$ | $\boldsymbol{q}_{\mathbf{1}}$ | $\ldots$ | $\ldots$ | $\mathbf{0}$ |

Example 13: Example of Jury Stability Analysis
Using Jury stability, determine the stability of the following two discrete transfer functions:

1) $G(z)=\frac{5}{z^{6}+5 z^{4}+6 z^{3}+2 z^{2}+3 z+1}$
2) $G(z)=\frac{4}{10 z^{4}+3 z^{3}+z^{2}+5 z+1}$

## Solution

For the first transfer function, we need to test the initial constraints before we consider creating the table. For condition $1, D(1)>0$, we get

$$
1+5+6+2+3+1=18>1
$$

which implies that this condition is satisfied. Condition $2,(-1)^{n} D(-1)>0$, gives

$$
(-1)^{6}\left[1(-1)^{6}+5(-1)^{4}+6(-1)^{3}+2(-1)^{2}+3(-1)^{1}+1\right]=0
$$

which is not satisfied. This implies that the first transfer function is not stable.
For the second transfer function, testing the three initial conditions gives
Condition (1): $\quad 10+3+1+5+1=20>0$ (satisfied)
Condition (2):

$$
(-1)^{4}\left[10(-1)^{4}+3(-1)^{3}+(-1)^{2}+5(-1)^{1}+1\right]=4>0
$$

Condition (3):

$$
\left|a_{n}\right|>\left|a_{0}\right| \Rightarrow|10|>|1|
$$

Since all the preliminary conditions have been satisfied, the Jury array can be constructed. This is shown in Table 13. The first row consists of the coefficients arranged in increasing powers, while the second row consists of the coefficients arranged in decreasing power. In Row 3, we compute the values using the formula

$$
\begin{equation*}
b_{i}=a_{0} a_{i}-a_{n} a_{n-i} \tag{108}
\end{equation*}
$$

which gives

$$
\begin{gathered}
b_{0}=a_{0} a_{0}-a_{4} a_{4-0}=1(1)-10(10)=-99 \\
b_{1}=a_{0} a_{1}-a_{4} a_{4-1}=(1) 5-10(3)=-25 \\
b_{2}=a_{0} a_{2}-a_{4} a_{4-2}=1(1)-10(1)=-9 \\
b_{3}=a_{0} a_{3}-a_{4} a_{4-3}=3(1)-10(5)=-47
\end{gathered}
$$

From here, we should check that the condition $\left|b_{0}\right|>\left|b_{n-1}\right|$ is satisfied. Since $|-99|>|-47|$, it holds and we proceed to the next step. We reverse the order of the remaining $n-1$ coefficients and write them in Row 4. Effectively, we have created a new polynomial of order $n-1$ that we need to test. Thus, in Row 5, we will use a modified form of Equation (108) to give

$$
\begin{equation*}
c_{i}=b_{0} b_{i}-b_{n-1} b_{n-i-1} \tag{109}
\end{equation*}
$$

Effectively, this is the same as Equation (108), but with the value of $n$ reduced by 1. Evaluating Equation (109) for the $n-2$ columns gives

$$
\begin{gathered}
c_{0}=b_{0} b_{0}-b_{3} b_{3}=(-99)^{2}-(-47)^{2}=7592 \\
c_{1}=b_{0} b_{1}-b_{3} b_{2}=-99(-25)-(-47)(-9)=2052 \\
c_{2}=b_{0} b_{2}-b_{3} b_{1}=-9(-99)-(-47)(-25)=-284
\end{gathered}
$$

Since we are left with three columns, we simply need to determine $\left|q_{0}\right|>\left|q_{2}\right|$, which in our case becomes $|7592|>|-284|$. Since it holds, we can conclude that the system is stable.

Table 13: Table for Jury stability

| Row | $\mathbf{0}$ | $\mathbf{1}$ | $\mathbf{2}$ | $\mathbf{3}$ | $\mathbf{4}$ |
| :---: | :---: | :---: | :---: | :---: | :---: |
| $\mathbf{1}$ | 1 | 5 | 1 | 3 | 10 |
| $\mathbf{2}$ | 10 | 3 | 1 | 5 | 1 |
| $\mathbf{3}$ | -99 | -25 | -9 | -47 |  |
| $\mathbf{4}$ | -47 | -9 | -25 | -99 |  |
| $\mathbf{5}$ | 7592 | 2052 | -284 |  |  |

## Section 3.3.2.3: Closed-Loop Stability Analysis

For closed-loop systems, stability can be analysed by considering the closed-loop transfer function, $G_{c l}$,

$$
\begin{equation*}
G_{c l}=\frac{G_{c} G_{p}}{1+G_{c} G_{p}} \tag{110}
\end{equation*}
$$

Since it has been shown that the poles of the system are the determining factor for stability, it therefore suffices to consider only the denominator of the closed-loop transfer function, that is, the term $1+G_{c} G_{p}$. It is possible to apply either the Routh or the Jury Stability Tests to determine under what conditions the system will be stable.

However, in some circumstances, it can be instructive to look at the frequency-domain plots, especially the Bode plot, to determine stability. In order to make the analysis simpler, we will rewrite the denominator to

$$
\begin{equation*}
G_{c} G_{p}=-1 \tag{111}
\end{equation*}
$$

Taking the magnitude (or modulus) of Equation (111) gives

$$
\begin{equation*}
\left\|G_{c} G_{p}\right\|=\|-1\| \tag{112}
\end{equation*}
$$

while the phase angle for -1 is $-180^{\circ}$, since the imaginary component is 0 and the real component is -1 , which after subtracting $-180^{\circ}$ gives the correct location. Therefore, plotting the Bode plot for the open-loop transfer function $G_{c} G_{p}$ will allow the stability of the system to be determined by examining the critical frequency, $\boldsymbol{\omega}_{\boldsymbol{c}}$, which is the frequency corresponding to a phase angle of $-180^{\circ}$. If the value of amplitude ratio at this point is greater than 1 (or 0 if a logarithmic basis is being used), then the process will be closed-loop unstable. An example is shown in Figure 26.


Figure 26: Bode Plot for Closed-Loop Stability Analysis
Furthermore, we can define two additional parameters of interest: the gain margin (GM) and the phase margin (PM). These two margins represent how much room we have before the system becomes unstable. On a Bode plot, instability implies an amplitude ratio above 1 (or 0 for the logarithmic case) or a phase angle below $-180^{\circ}$. These parameters are shown in Figure 27 for the Bode plot and in Figure 28 for the Nyquist plot. The phase margin is defined by adding $180^{\circ}$ to the phase angle when the gain crosses the value of 1 , which is denoted by $\omega_{g}$ and called the gain crossover, that is,

$$
\begin{equation*}
\mathrm{PM}=\phi\left(\omega_{g}\right)+180^{\circ} \tag{113}
\end{equation*}
$$

The gain margin is defined as the difference between 1 (or 0 for a logarithmic case) and the value when the phase angle crosses the phase angle of $-180^{\circ}$, which is denoted by $\omega_{p}$ and called the phase crossover, that is,

$$
\begin{equation*}
\mathrm{GM}=1-\mathrm{AR}\left(\omega_{p}\right) \tag{114}
\end{equation*}
$$

It is possible to design controllers by specifying the phase and gain margins.


Figure 27: Closed-loop stability using the Bode plot


Figure 28: Closed-loop stability using the Nyquist plot

## Section 3.3.3: Controllability and Observability

Since we are ultimately interested in automating our process, it is important to understand if our process can in fact be automated. For state-space representations, it is very common to talk about the controllability and observability of the process. The state equation or the pair $(\mathcal{A}, \mathcal{C})$ is observable if and only if for any unknown initial state, $x_{0}$, there exists a finite time $t$ (greater than
zero), such that the knowledge of the input, $u(t)$, and output, $y(t)$ in the interval $[0, t]$ is sufficient to determine the initial state, $x_{0}$. This can be accomplished if and only if

1) $\mathcal{O}=\left[\begin{array}{c}\mathcal{C} \\ \mathcal{C A} \\ \mathcal{C A}^{2} \\ \vdots \\ \mathcal{C A}^{n-1}\end{array}\right]_{n \times \times n}$ has full rank $n$, where $\mathcal{O}$ is the observability matrix.
2) The observability Grammian, $\mathcal{W}_{o}(t)_{n \times n}=\int_{0}^{t} e^{\mathcal{A}^{T} \tau} \mathcal{C}^{T} \mathcal{C} e^{\mathcal{A} \tau} d \tau$ is nonsingular for all $t>0$. The energy of the output signal can be given as $E_{y}(t)=x^{T}(0) \mathcal{W}_{c}(t) x(0)$.
3) The $n$-by-n matrix $\left[\frac{\mathcal{A}-\lambda_{i} \mathcal{I}}{\mathcal{C}}\right]$ has rank $n$ for all eigenvalues $\lambda_{i}$ of $\mathcal{A}$. This allows the individual states or eigenvalues of $\mathcal{A}$ to be classified.
A system is said to be detectable if all unstable states of a state-space representation can be observed. This can be determined by examining $\left[\frac{\mathcal{A}-\lambda_{i} \mathcal{I}}{\mathcal{C}}\right]$ for each of the unstable states and determining if it is full rank.

The state equation or the pair $(\mathcal{A}, \mathcal{B})$ are said to be controllable if and only if for any initial state, $x_{0}$, and final state, $x_{1}$, there exists an input, $u(t)$, that transfers $x_{0}$ to $x_{1}$ in finite time. This can be accomplished if and only if:

1) $\overline{\mathcal{C}}_{n \times n p}=\left[\begin{array}{lllll}\mathcal{B} & \mathcal{A B} & \mathcal{A}^{2} \mathcal{B} & \cdots & \mathcal{A}^{n-1} \mathcal{B}\end{array}\right]$ has full rank $n$, where $\overline{\mathcal{C}}$ is the controllability matrix.
2) The controllability Grammian, $\mathcal{W}_{c}(t)_{n \times n}=\int_{0}^{t} e^{\mathcal{A} \tau} \mathcal{B B}^{T} e^{\mathcal{A}^{T} \tau} d \tau$, is nonsingular for all $t>0$. $\mathcal{W}_{c}^{-1}$ is a measure of the energy required to control the process. A larger value implies that more energy (or effort) is required.
3) The $n$-by- $n$ matrix $\left[\mathcal{A}-\lambda_{i} \mathcal{J} \mid \mathcal{B}\right]$ has rank $n$ for all eigenvalues $\lambda_{i}$ of $\mathcal{A}$. This allows the individual states or eigenvalues of $\mathcal{A}$ to be classified.

A system is said to be stabilisable if all unstable states of a state-space representation can be controlled. This can be determined by examining $\left[\mathcal{A}-\lambda_{i} \mathcal{J} \mid \mathcal{B}\right]$ for each of the unstable states and determining if it is full rank.

It can be noted that controllability and observability are duals of each other, that is, if ( $\mathcal{A}$, $\mathcal{B})$ is controllable, then $\left(\mathcal{A}^{T}, \mathcal{B}^{T}\right)$ is observable. Similarly, if $(\mathcal{A}, \mathcal{C})$ is observable, then $\left(\mathcal{A}^{T}, \mathcal{C}^{T}\right)$ is controllable.

## Section 3.3.4: Analysis of Special Transfer Functions

In this section, the properties of different commonly encountered transfer functions will be considered. Since much of the analysis in automation focuses on using transfer functions, it is important to understand the behaviour of the common transfer functions. Time-domain responses of these transfer functions to a step input will also be considered, since it is important to recognise these transfer functions in their most common manifestations in a real process. Discrete time systems will only be briefly considered, since most discrete time analysis is still based on the underlying continuous-time systems.

## Section 3.3.4.1: Integrator

The integrator, as its name suggests, integrates a variable. It is a common model for level in a tank. Its Laplace transform is

$$
\begin{equation*}
G_{I}=\frac{M_{I}}{s} \tag{115}
\end{equation*}
$$

where $M_{I}$ is the gain. It is an unstable system, which will always increase even if the input is bounded. The step response of the system can be determined as

$$
\begin{align*}
Y & =G_{I} U \\
Y & =\frac{M_{I}}{s} \frac{M}{s}  \tag{116}\\
& =\frac{M_{I} M}{s^{2}}
\end{align*}
$$

From Table 4, the time domain representation for Equation (116) is

$$
\begin{equation*}
y_{t}=M_{I} M t \tag{117}
\end{equation*}
$$

This shows that the integrator will continual increase even if the input is bounded. A representative plot is shown in Figure 29. Its Bode plot can be determined as

$$
\begin{align*}
& G_{I}(j \omega)=\frac{M_{I}}{j \omega}=-\frac{M_{I} j}{\omega} \\
& \Rightarrow \mathrm{AR}=\sqrt{0^{2}+\left(-\frac{M_{I}}{\omega}\right)^{2}}=\frac{\left|M_{I}\right|}{\omega}  \tag{118}\\
& \phi=\arctan \left(\frac{-\frac{M_{I}}{\omega}}{0}\right)=-90^{\circ}
\end{align*}
$$

A representative Bode plot is shown in Figure 30. The key properties of the integrator are summarised in Table 14.

Table 14: Basic Information About an Integrator


Figure 29: Response of the Integrator $1 / s$ to a Unit-Step Input


Figure 30: Bode Plot for an Integrator

## Section 3.3.4.2: Lead Term

The lead term is a simple, first-order transfer function that is occasionally found in combination with other transfer function to model unusual or more complex initial dynamics in a process. Its Laplace transform is

$$
\begin{equation*}
G_{L}=K\left(\tau_{L} s+1\right) \tag{119}
\end{equation*}
$$

where $K$ is the gain and $\tau_{L}$ is a time constant. The step response of the system can be determined as

$$
\begin{align*}
Y & =G_{L} U \\
Y & =K\left(\tau_{L} s+1\right) \frac{M}{s}  \tag{120}\\
& =K M \frac{\left(\tau_{L} s+1\right)}{s}
\end{align*}
$$

In order to obtain a time-domain representation, it is necessary to perform a partial fraction decomposition of Equation (120) to give

$$
\begin{align*}
& Y=K M \frac{\left(\tau_{L} s+1\right)}{s}  \tag{121}\\
& Y=K M\left(\tau_{L}+\frac{1}{s}\right)
\end{align*}
$$

From Table 5, the time domain representation for Equation (121) is

$$
\begin{equation*}
y_{t}=K M\left(\tau_{L} \delta+u_{t}\right) \tag{122}
\end{equation*}
$$

where $\delta$ is the Dirac delta function and $u_{t}$ is the unit step function. This shows that the lead function is bounded. Its Bode plot can be determined as

$$
\begin{align*}
& G_{I}(j \omega)=K\left(\tau_{L} j \omega+1\right)=K+K \tau_{L} \omega j \\
& \Rightarrow \mathrm{AR}=\sqrt{K^{2}+\left(K \tau_{L} \omega\right)^{2}}=|K| \sqrt{1+\tau_{L}^{2} \omega^{2}}  \tag{123}\\
& \phi=\arctan \left(\frac{K \tau_{L} \omega}{K}\right)=\left\{\begin{array}{cc}
\arctan \tau_{L} \omega & K>0 \\
\arctan \tau_{L} \omega+180^{\circ} & K<0
\end{array}\right.
\end{align*}
$$

A representative Bode plot is shown in Figure 31 for all 4 possible combinations of $K$ and $\tau_{L}$. The key properties of the lead term are summarised in Table 15.

Table 15: Basic Information About a Lead Term

| Property |  | Value |
| :---: | :---: | :---: |
| Laplace Transform |  | $G_{L}=K\left(\tau_{L} s+1\right)$ |
| Step Response (Time Domain) |  | $y_{t}=K M\left(\tau_{L} \delta+u_{t}\right)$ |
| $\begin{aligned} & \frac{0}{0} \\ & \frac{0}{2} \\ & \frac{0}{0} \\ & 0 \end{aligned}$ | AR | $\|K\| \sqrt{1+\tau_{L}^{2} \omega^{2}}$ |
|  | $\phi$ | $\left\{\begin{array}{cc}\arctan \tau_{L} \omega & K>0 \\ \arctan \tau_{L} \omega+180^{\circ} & K<0\end{array}\right.$ |
| Stable |  | Yes |
| Comment |  | Negative values of $\tau_{L}$ (positive zeros) can induce an inverse response in a system, that is, the variable first decreases in value and increases to reach its new steady-state value (or vice versa). |



Figure 31: Bode Plot for a Lead Term (top row) $K>0$, (bottom row) $K<0$, (left) $\tau_{L}>0$, and (right) $\tau_{L}<0$

## Section 3.3.4.3: First-Order Transfer Function

A first-order system is one of the most common transfer functions encountered in automation engineering. It can be used to model any system ranging from simple heated tanks to complex multicomponent reactions. Its Laplace transform is

$$
\begin{equation*}
G_{F}=\frac{K}{\left(\tau_{p} s+1\right)} \tag{124}
\end{equation*}
$$

where $K$ is the gain and $\tau_{p}$ is the process time constant. It is often coupled with a deadtime term to give a first-order plus deadtime (FOPDT) process model, that is,

$$
\begin{equation*}
G_{F D}=\frac{K}{\left(\tau_{p} s+1\right)} e^{-\theta s} \tag{125}
\end{equation*}
$$

The step response of the pure, first-order system can be determined as

$$
\begin{align*}
Y & =G_{F} U \\
Y & =\frac{K}{\left(\tau_{p} s+1\right)} \frac{M}{s}  \tag{126}\\
& =\frac{K M}{s\left(\tau_{p} s+1\right)}
\end{align*}
$$

In order to obtain a time-domain representation, it is necessary to perform a partial fraction decomposition of Equation (126) to give

$$
\begin{align*}
& Y=\frac{K M}{s\left(\tau_{p} s+1\right)}  \tag{127}\\
& Y(s)=K M\left(\frac{1}{s}+\frac{-\tau}{\tau s+1}\right)
\end{align*}
$$

From Table 5, the time domain representation for Equation (127) is

$$
\begin{equation*}
y(t)=K M\left(1-e^{-\frac{t}{\tau}}\right) \tag{128}
\end{equation*}
$$

The process is stable if $\tau>0$, but unstable if $\tau<0$. Assuming $\tau>0$, then using the final value theorem, it can be shown that the new steady-state value will be

$$
\begin{align*}
\lim _{t \rightarrow \infty} y_{t} & =\lim _{s \rightarrow 0} s Y(s)=\lim _{s \rightarrow 0} s \frac{M G_{F}(s)}{s} \\
& =\lim _{s \rightarrow 0} M G_{F}(s)=\lim _{s \rightarrow 0} M G_{F}(0)  \tag{129}\\
& =K M
\end{align*}
$$

The time response of a stable first-order system is shown in Figure 32, as well as how to compute the key parameters from its graph. Its Bode plot can be determined as

$$
\begin{align*}
& G_{I}(j \omega)=\frac{K}{\left(\tau_{p} j \omega+1\right)}=\frac{K\left(1-\tau_{p} \omega j\right)}{1+\tau_{p}^{2} \omega^{2}} \\
& \Rightarrow \mathrm{AR}=\sqrt{\frac{K^{2}+\left(-K \tau_{p} \omega\right)^{2}}{\left(1+\tau_{p}^{2} \omega^{2}\right)^{2}}}=\frac{|K|}{\sqrt{1+\tau_{p}^{2} \omega^{2}}}  \tag{130}\\
& \phi=\arctan \left(\frac{-K \tau_{p} \omega}{K}\right)=\left\{\begin{array}{cc}
-\arctan \tau_{p} \omega & K>0 \\
180^{\circ}-\arctan \tau_{p} \omega & K<0
\end{array}\right.
\end{align*}
$$

In order to avoid a discontinuity when it pass through $0^{\circ}$, it is common to use negative angles when plotting this function, that is $-45^{\circ}$ rather than the equivalent $315^{\circ}$. A representative Bode plot is
shown in Figure 31 for all 4 possible combinations of $K$ and $\tau_{p}$. The key properties of the firstorder system are summarised in Table 16.

Table 16: Basic Information About a First-Order System

| Property |  | Value |
| :---: | :---: | :---: |
| Laplace Transform |  | $G_{F}=\frac{K}{\left(\tau_{p} s+1\right)}$ |
| Step Response (Time Domain) |  | $y(t)=K M\left(1-e^{-\frac{t}{\tau}}\right)$ |
| $\begin{aligned} & \stackrel{n}{0} \\ & \frac{0}{2} \\ & 0 \\ & 0 \\ & 0 \end{aligned}$ | AR | $\frac{\|K\|}{\sqrt{1+\tau_{p}^{2} \omega^{2}}}$ |
|  | $\phi$ | $\left\{\begin{array}{cc}-\arctan \tau_{p} \omega & K>0 \\ 180^{\circ}-\arctan \tau_{p} \omega & K<0\end{array}\right.$ |
| Stable |  | Yes, if $\tau_{p}>0$ |
| Comment |  | This is one of the most common transfer functions encountered in process control and is used to model many different applications, often by including a time delay term. |



Figure 32: Step Response of a Stable, First-Order System to a Step Response


Figure 33: Bode Plot for a First-Order System (top row) $K>0$, (bottom row) $K<0$, (left) $\tau_{p}>0$, and (right) $\tau_{p}<0$

## Section 3.3.4.4: Second-Order System

A second-order system is another commonly encountered transfer function that is often used to model oscillations or periodic behaviour in a process. Its Laplace transform is

$$
\begin{equation*}
G_{I I}=\frac{K}{\left(\tau_{p}^{2} s^{2}+2 \zeta \tau_{p} s+1\right)} \tag{131}
\end{equation*}
$$

where $K$ is the gain, $\tau_{p}$ is the process time constant, and $\zeta$ is the damping coefficient. This basic transfer function can be augmented by adding a lead term to model various behaviours to give

$$
\begin{equation*}
G_{I I}=\frac{K\left(\tau_{L} s+1\right)}{\left(\tau_{p}^{2} s^{2}+2 \zeta \tau_{p} s+1\right)} \tag{132}
\end{equation*}
$$

As well, it can be coupled with a deadtime term to give a second-order plus deadtime (SOPDT) process model, that is,

$$
\begin{equation*}
G_{I I D}=\frac{K}{\left(\tau_{p}^{2} s^{2}+2 \zeta \tau_{p} s+1\right)} e^{-\theta s} \tag{133}
\end{equation*}
$$

The poles of the second-order transfer function can be written as

$$
\begin{align*}
s & =\frac{-2 \zeta \tau_{p} \pm \sqrt{4 \zeta^{2} \tau_{p}^{2}-4 \tau_{p}^{2}}}{\tau_{p}^{2}} \\
& =\frac{-2 \zeta \pm 2 \sqrt{\zeta^{2}-1}}{\tau_{p}} \tag{134}
\end{align*}
$$

Depending on the value of $\zeta$, three different cases can be determined:

1) Case I: Underdamped System, where $|\zeta|<1$. In this case, the poles will contain an imaginary component.
2) Case II: Critically Damped System, where $|\zeta|=1$. In this case, the poles will both be real and the same.
3) Case III: Overdamped System, where $|\zeta|>1$. In this case, the poles will both be real.

The behaviour and critical information depends on the particular case being considered.
For the underdamped case, where $|\zeta|<1$, the poles of Equation (131) will contain an imaginary component. The presence of an imaginary component will imply that there will be oscillations present in the time response. If $\zeta=0$, then the system will continuously oscillate about a mean value. In all other cases, stability will depend on the sign of $\zeta \tau_{p}$ : if positive, the system will be stable (decaying oscillations); if negative, the system will be unstable (increasing oscillations). The step response of this system can be written as

$$
\begin{equation*}
y(t)=K M\left(1-e^{-\zeta t / \tau_{p}}\left[\cos \left(\frac{\sqrt{1-\zeta^{2}}}{\tau_{p}} t\right)+\frac{\zeta^{2}}{\sqrt{1-\zeta^{2}}} \sin \left(\frac{\sqrt{1-\zeta^{2}}}{\tau_{p}} t\right)\right]\right) \tag{135}
\end{equation*}
$$

A typical underdamped step response is shown in Figure 34. Some of the key parameters that can be extracted from a step response are:

1) Time to First Peak: $t_{p}=\pi \tau / \sqrt{1-\zeta^{2}}$
2) Overshoot: $O S=\exp \left(-\pi \zeta / \sqrt{1-\zeta^{2}}\right)=\frac{a}{b}$.
3) Decay Ratio: The decay ratio is the square of the overshoot ratio, i.e.
$D R=\exp \left(-2 \pi \zeta / \sqrt{1-\zeta^{2}}\right)=\frac{c}{a}$.
4) Period of Oscillation: This is the time between 2 oscillations. It is given by

$$
P=\frac{2 \pi \tau}{\sqrt{1-\zeta^{2}}} .
$$

5) Settling Time, $\boldsymbol{t}_{s}$ : The time required for the process to remain within $5 \%$ of the steadystate value. The first time this occurs is called the settling time.
6) Rise Time, $t_{r}$ : Time required to first reach the steady state value.


Figure 34: Second-Order Underdamped Process

For the critically damped system, where $|\zeta|=1$, the poles of Equation (131) will both be the same and contain no imaginary component. The system will be stable if the sign of $\zeta \tau_{p}$ is positive and unstable if $\zeta \tau_{p}$ is negative. The step response of this system can be written as

$$
\begin{equation*}
y(t)=K M\left(1-\left(1+\frac{t}{\tau_{p}}\right) e^{-\zeta t / \tau_{p}}\right) \tag{136}
\end{equation*}
$$

A typical stable, critically damped step response is shown in Figure 35. In general, such a system looks very similar to a first-order plus deadtime system and is often analysed as such. The biggest difference is the slightly slower initial response, which is often treated as a time delay.


Figure 35: Step Response of a Critically Damped System ( $\zeta=1, \tau_{p}=10$, and $K=1$ )
For the overdamped case, where $|\zeta|>1$, the poles of Equation (131) will only be real numbers. This implies that there will be no oscillations in the step response. The system will be stable if the sign of $\zeta \tau_{p}$ is positive and unstable if $\zeta \tau_{p}$ is negative. The step response of this system can be written as ${ }^{12}$

$$
\begin{equation*}
y(t)=K M\left(1-e^{-\zeta t / \tau_{p}}\left[\cosh \left(\frac{\sqrt{\zeta^{2}-1}}{\tau_{p}} t\right)+\frac{\zeta^{2}}{\sqrt{\zeta^{2}-1}} \sinh \left(\frac{\sqrt{\zeta^{2}-1}}{\tau_{p}} t\right)\right]\right) \tag{137}
\end{equation*}
$$

A typical overdamped step response is shown in Figure 36. In most cases, this system can be analysed as a first-order plus deadtime process. It can be noted that the overdamped case has the

[^9]slowest initial response of all second-order and first-order systems. This slow initial response is often treated as a deadtime when fitting a first-order system.


Figure 36: Step Response of an Overdamped System $\left(\zeta=2, \tau_{p}=5\right.$, and $\left.K=1\right)$
The Bode plot of a second-order system can be determined as

$$
\begin{align*}
& G_{I I}(j \omega)=\frac{K}{\tau_{p}^{2}(j \omega)^{2}+2 \zeta \tau_{p} j \omega+1}=\frac{K}{1-\tau_{p}^{2} \omega^{2}+2 \zeta \tau_{p} j \omega}=\frac{K\left(1-\tau_{p}^{2} \omega^{2}-2 \zeta \tau_{p} j \omega\right)}{\left(1-\tau_{p}^{2} \omega^{2}\right)^{2}+4 \zeta^{2} \tau_{p}^{2} \omega^{2}} \\
& \Rightarrow \operatorname{AR}=\sqrt{\frac{K^{2}\left(1-\tau_{p}^{2} \omega^{2}\right)^{2}+K^{2}\left(-2 \zeta \tau_{p} \omega\right)^{2}}{\left(\left(1-\tau_{p}^{2} \omega^{2}\right)^{2}+4 \zeta^{2} \tau_{p}^{2} \omega^{2}\right)^{2}}}=\frac{|K|}{\sqrt{\left(1-\tau_{p}^{2} \omega^{2}\right)^{2}+4 \zeta^{2} \tau_{p}^{2} \omega^{2}}} \\
& \phi=\arctan \left(\begin{array}{cc}
-\arctan \frac{2 \zeta \tau_{p} \omega}{1-\tau_{p}^{2} \omega^{2}} & K>0,\left|\tau_{p} \omega\right|<1 \\
K\left(1-\tau_{p}^{2} \omega^{2}\right)
\end{array}\right)=\left\{\begin{aligned}
-\arctan \frac{2 \zeta \tau_{p} \omega}{1-\tau_{p}^{2} \omega^{2}}-180^{\circ} & K>0,\left|\tau_{p} \omega\right|>1 \\
180^{\circ}-\arctan \frac{2 \zeta \tau_{p} \omega}{1-\tau_{p}^{2} \omega^{2}} & K<0,\left|\tau_{p} \omega\right|<1 \\
-\arctan \frac{2 \zeta \tau_{p} \omega}{1-\tau_{p}^{2} \omega^{2}} & K<0,\left|\tau_{p} \omega\right|>1
\end{aligned}\right.
\end{align*}
$$

When dealing with second-order systems, there is a potential that the amplitude ratio can be greater than the original starting value of $|K|$. Specifically, this occurs whenever the denominator is less than 1. Taking the denominator of the amplitude ratio for a second-order system and solving it gives

$$
\begin{align*}
\left(1-\tau_{p}^{2} \omega^{2}\right)^{2}+4 \zeta^{2} \tau_{p}^{2} \omega^{2} & <1 \\
1-2 \tau_{p}^{2} \omega^{2}+\tau_{p}^{4} \omega^{4}+4 \zeta^{2} \tau_{p}^{2} \omega^{2} & <1 \\
-2+\tau_{p}^{2} \omega^{2}+4 \zeta^{2} & <0  \tag{139}\\
\omega^{2} & <\frac{2-4 \zeta^{2}}{\tau_{p}^{2}} \\
|\omega| & <\sqrt{\frac{2-4 \zeta^{2}}{\tau_{p}^{2}}}=\frac{\sqrt{2-4 \zeta^{2}}}{\left|\tau_{p}\right|}
\end{align*}
$$

Since the frequency is only a positive real number, the term under the square root must also only be positive (since a negative would give a complex number), that is,

$$
\begin{align*}
2-4 \zeta^{2} & \geq 0 \\
|\zeta| & \geq \sqrt{0.5} \approx 0.707 \tag{140}
\end{align*}
$$

If the damping coefficient lies within the region given by Equation (140), then there will be a hump in the Bode plot, as shown in Figure 37 (top). Since these frequencies magnify the system response, it is important that special care be taken when designing controllers for such systems.

If a lead term has been added to a second-order system, as given by Equation (132), then the system will display an inverse response if the zeros of the transfer function are positive, for example, as shown in Figure 38. The Bode plot for this composite system can be obtained by invoking the rules for combinations of transfer functions, that is,

$$
\begin{align*}
G_{I I} & =\frac{K\left(\tau_{L} s+1\right)}{\left(\tau_{p}^{2} s^{2}+2 \zeta \tau_{p} s+1\right)}=G_{I I} G_{L} \\
\mathrm{AR} & =\mathrm{AR}_{I I} \mathrm{AR}_{L}=\frac{|K| \sqrt{1+\tau_{L}^{2} \omega^{2}}}{\sqrt{\left(1-\tau_{p}^{2} \omega^{2}\right)^{2}+4 \zeta^{2} \tau_{p}^{2} \omega^{2}}}  \tag{141}\\
\phi & =\phi_{I I}+\phi_{L}=\arctan \tau_{L} \omega-\arctan \frac{2 \zeta \tau_{p} \omega}{1-\tau_{p}^{2} \omega^{2}}
\end{align*}
$$

(ignoring all angle complications)
An example is shown in Figure 39.
Representative Bode plots are shown in Figure 37 for different combinations of the parameters. The key properties of the second-order system are summarised in Table 17.

Table 17: Basic Information About a Secord-Order System

| Property |  | Value |
| :---: | :---: | :---: |
| Laplace Transform |  | $G_{I I}=\frac{K}{\left(\tau_{p}^{2} s^{2}+2 \zeta \tau_{p} s+1\right)}$ |
| Step Response (Time Domain) |  | $\begin{aligned} & \|\zeta\|<1: \text { Equation }(135) \\ & \|\zeta\|=1: \text { Equation }(136) \\ & \|\zeta\|>1: \text { Equation }(137) \end{aligned}$ |
| $\begin{aligned} & \frac{0}{0} \\ & \frac{0}{1} \\ & \frac{0}{0} \\ & 0 \end{aligned}$ | AR | $\frac{\|K\|}{\sqrt{\left(1-\tau_{p}^{2} \omega^{2}\right)^{2}+4 \zeta^{2} \tau_{p}^{2} \omega^{2}}}$ |
|  | $\phi$ | $-\arctan \frac{2 \zeta \tau_{p} \omega}{1-\tau_{p}^{2} \omega^{2}}$ <br> see Equation (138) for greater detail |
| Stable |  | Yes, if $\zeta \tau_{p}>0$ |
| Comment |  | This transfer function is commonly used to describe oscillations and inverse responses in a system. |



Figure 37: Bode Plots for $\tau_{p}=5$, (top) $\zeta=0.5$, (middle) $\zeta=1$, (bottom) $\zeta=2$; (left) $K=1$ and (right) $K=-1$


Figure 38: Step Response of a Second-Order System with Inverse Response ( $\tau_{L}=-25, \zeta=5 / \sqrt{ } 7, \tau_{p}=10 \sqrt{ } 7$ and $K=1$ )


Figure 39: Bode Plots for (left) $\tau_{L}=-5$, (right) $\tau_{L}=5$, and (top) $\zeta=0.5$ and (bottom) $\zeta=2$

## Section 3.3.4.5: Higher-Order Systems

Analysing higher-order systems is based on the results obtained from the analysis of the simpler systems. The following are some key points to consider for the analysis in the time domain:

1) Poles: The poles of the process transfer function determine the stability of the system. If the poles are less than zero, then the system is stable; otherwise, it is unstable. If the poles contain an imaginary component, then there will be oscillations in the system.
2) Zeros: If the zeros of the transfer function are positive, then there will be an inverse response in the system.
3) Time Constant: The largest stable pole in absolute magnitude can be treated as the dominating time constant of the process.
In the frequency domain, the Bode plots are obtained by combining the appropriate simple transfer functions, using the rules for composition of the amplitude ratio (multiplication) and phase angle (addition).

## Example 14: Sketching the Expected Time-Domain Response

For the following three transfer functions, sketch the expected time-domain response of the system when a positive step change is made:

1) $G_{1}=\frac{1.54(-5 s+1)}{(5 s+1)(4 s+2)(2 s+1)} e^{-10 s}$
2) $G_{2}=\frac{1.54(5 s+1)}{100 s^{2}-200 s+1}$
3) $G_{3}=\frac{1.54}{\left(-100 s^{2}+10 s-1\right)(4 s+1)}$

## Solution:

## First Transfer Function

For the first transfer function, the gain is 1.54 (by inspection), the zero is 0.2 , and the poles are $-0.2(=-1 / 5),-0.5(=-2 / 4)$, and $-0.5(=-1 / 2)$, and the time delay is 10 . Since the zero is positive, we expect an inverse response, while all poles are negative, which implies that the system is stable. As well, none of the poles have an imaginary component and so there are no oscillations. The plot would therefore be as given in Figure 40 (left).

## Second Transfer Function

For the second transfer function, the zero is -0.2 . Instead of computing the poles directly, we can note that this is a second-order system. Therefore, writing it into the standard form for a second-order system, it can be seen that $\tau_{p}=10$ and $\zeta=-10$. Since $\tau_{p} \zeta<0$, the system is unstable and without any oscillations $(|\zeta|>1)$. The "gain" is positive, which implies that the function will go towards $+\infty$. Thus, the plot would be as given in Figure 40 (middle).

## Third Transfer Function

For the third transfer function, there are no zeros or time delay. Similar to the second transfer function, let us rewrite the denominator into the required form to give

$$
G_{3}=\frac{1.54}{\left(-100 s^{2}+10 s-1\right)(4 s+1)} \frac{-1}{-1}=\frac{-1.54}{\left(100 s^{2}-10 s+1\right)(4 s+1)} .
$$

By comparison, we can see that the second-order term has the following characteristics, $\tau_{p}=10$ and $\zeta=-0.5$. This implies that the poles are unstable and oscillatory. The second term will give a stable pole $(-0.25=-1 / 4)$. Therefore, the plot of this transfer function will be given as Figure 40 (right).


Figure 40: Sketch of the Transfer Function Step Responses: (left) first transfer function, (middle) second transfer function, and (right) third transfer function.

It can be noted that when sketching the transfer function, it is important to only give the general characteristics, while the exact values can be ignored.

## Section 3.3.4.6: Summary of Functional Behaviour in Continuous- and Discrete-Time Domains

Table 18 summarises the relationship between the location of the poles of the transfer function and the generalised behaviour in both the continuous- and discrete-time domains. The
ringing cases can only occur in the discrete-time domain as a result of how the discretisation is taken.

Table 18: Graphical Representation of the Different Types of Functions (The ringing cases can only occur in the discrete domain.)


| Case | Location of the Poles |  | Unit-Step Response |  |
| :---: | :---: | :---: | :---: | :---: |
|  | Continuous | Discrete | Continuous | Discrete |
| Ringing, integrator |  |  |  |  |
| Ringing, unstable |  |  |  | - |
| Ringing, unstable, oscillatory |  |  |  | Emmumummumbundy |

## Example 15: Origin of Ringing in Discrete-Time Systems

Consider the continuous exponential and cosine system

$$
\begin{equation*}
y(t)=e^{a t} \cos (\omega t) \tag{142}
\end{equation*}
$$

that is sampled with a sampling time of $T_{s}$ to give a discrete-time system of the form

$$
\begin{equation*}
y_{k}=e^{a k T_{s}} \cos \left(\omega k T_{s}\right) \tag{143}
\end{equation*}
$$

which, from Table 6, has the $z$-transform of

$$
\begin{equation*}
\frac{1-e^{a T_{s}} \cos \left(\omega T_{s}\right) z^{-1}}{1-2 e^{a T_{s}} \cos \left(\omega T_{s}\right) z^{-1}+e^{2 a T_{s}} z^{-2}} \tag{144}
\end{equation*}
$$

We will assume that $a \in \mathbb{R}$ so that values of $\alpha \geq 0$ will lead to an unstable system. It can be noted that the results for sine will be the same mutatis mutanda. Examine the impact of the sampling time on the resulting discrete-time function and its $z$-transform.

## Solution

The behaviour of the transfer function is determined by the values of its poles, that is, the roots of the denominator. Using the quadratic formula gives the general form of the poles as

$$
\begin{equation*}
z=\frac{2 \psi \cos \left(\omega T_{s}\right) \pm \sqrt{4 \psi^{2} \cos ^{2}\left(\omega T_{s}\right)-4 \psi^{2}}}{2}=\psi \cos \left(\omega T_{s}\right) \pm \psi \sqrt{\cos ^{2}\left(\omega T_{S}\right)-1} \tag{145}
\end{equation*}
$$

where to simplify notation $\psi=e^{a T_{s}}$. We can note that $\psi$ will always be positive since the exponential of any real number is positive.

In order to understand the behaviour of the discrete function, we will need to examine the determinant of the equation (the part inside the square root) and consider three cases: when the determinant is greater than zero, exactly zero, and less than zero.

Case 1: Determinant greater than zero
For the determinant to be greater than zero, it follows that

$$
\begin{equation*}
\cos ^{2}\left(\omega T_{S}\right)-1>0 \Rightarrow \cos ^{2}\left(\omega T_{s}\right)>1 \tag{146}
\end{equation*}
$$

However, the cosine function is never greater than 1 . Thus, this situation cannot occur and there will not be two distinct real roots.

Case 2: Determinant equal to zero
For the determinant to be exactly zero, it follows that

$$
\begin{equation*}
\cos ^{2}\left(\omega T_{S}\right)-1=0 \Rightarrow \cos ^{2}\left(\omega T_{s}\right)=1 \tag{147}
\end{equation*}
$$

Taking the square root of the right-hand side and noting that there are two solutions gives

$$
\begin{equation*}
\cos \left(\omega T_{s}\right)= \pm 1 \tag{148}
\end{equation*}
$$

which occurs when

$$
\omega T_{s}=\pi n
$$

for $n \in \mathbb{Z}$. If we restrict ourselves to the domain $[0,2 \pi[$, we see that we have two solutions: at $n=$ 0 with a value of 1 and $n=1$ with a value of -1 . In such a case, the double root will be located on the $x$-axis at the location of $\psi$ for $n$ even and at the location of $-\psi$ for $n$ odd. When $n$ is odd, the poles will lie in the left-hand plane of the discrete domain, which implies that we will have ringing behaviour. Stability will be determined by the value of $\psi$. However, there will not be any oscillatory behaviour. The general form of the discrete function is then

$$
\begin{equation*}
y_{k}=\psi^{k} \cos (k \pi n) \tag{149}
\end{equation*}
$$

When $n$ is odd, the value of $\cos (k \pi n)$ will oscillate between 1 and -1 which will give the characteristic ringing behaviour.
Case 3: Determinant less than zero
For the determinant to be less than zero, it follows that

$$
\begin{equation*}
\cos ^{2}\left(\omega T_{S}\right)-1<0 \Rightarrow \cos ^{2}\left(\omega T_{s}\right)<1 \tag{150}
\end{equation*}
$$

In such cases, there will be two imaginary roots that are complements of each other, that is, $\psi+\gamma i$ and $\psi-\gamma i$, where $\gamma$ is equal to $\psi \sqrt{1-\cos ^{2}\left(\omega T_{S}\right)}$. This implies that there will be a clear oscillatory
behaviour. If we restrict ourselves to the domain $\left[0,2 \pi\left[\right.\right.$, we can note that $\cos \left(\omega T_{s}\right)$ will be positive in the region $[0,0.5 \pi[\mathrm{U}] 1.5 \pi, 2 \pi[$ and negative in the region $] 0.5 \pi, 1.5 \pi[$. The function will be exactly zero at $0.5 \pi$ and $1.5 \pi$. In the positive region, this will place the poles in the right-hand side of the discrete system, while when negative it will place them in the left-hand side leading to ringing. When $\cos \left(\omega T_{s}\right)$ is exactly zero, we will have two roots at the origin of the system and we will have no information about the system. This brings us directly to the constraint given by the Shannon sampling theorem that the sampling time must lie be greater than $2 / \omega$.

## Section 3.4: Event-Based Representations

In event-based systems, where the changes in events drive the process, a different type of model is required, where the impact of the events can be more clearly seen on the system. An automaton or, in the older literature machine, is a way of representing how a process changes from state to state based on discrete events. At each state, a series of inputs are recognised and cause the automaton to move to another (or perhaps even, the same) state.

Before considering a formal definition of an automaton, let us examine a simple case and how it could be modelled as an automaton. Consider a system that consists of only two inputs $a$ and $b$. The objective of this system is to determine when the sequence of inputs is $b a b a$. In automaton theory, the acceptable inputs are called the alphabet, while the actual sequence of inputs is called a word. Thus, in this example, the alphabet is the set $\{a, b\}$ and the word of interest is baba. A state is defined as an internal representation of the system and its expected behaviour. In this example, we can define 5 states: Z 0 that represents the initial state, Z 1 when we have received the first $b, \mathrm{Z} 2$, when we have received $b a, \mathrm{Z} 3$, when we have received $b a b$, and Z 4 , when we have received baba. Based on the inputs and current state, a transition function shows how the system will move to the next state. For example, if we are in state Z 0 and receive an $a$, then the system will transition to Z 0 (remain in the same place), while if a $b$ is received, the system will transition to Z1. States can be classified as either accepting (also called marked states) or rejecting. An accepting state is one that has the desired outcome. In this particular example, it would be state Z 4 . A rejecting state is one that does not have the desired outcome. In this particular example, it would be all the other states. When an automaton has finished reading a word, we can note the final state of the automaton. If it is an accepting state, then we can say that the automaton
accepts the word; otherwise it rejects the word. The set of all words accepted by the automaton is called the language recognised (or marked) by the automaton.

Mathematically, for a finite-state automaton $A$, we can write this as the quintuple

$$
\begin{equation*}
A \equiv\left(\mathbb{X}, \Sigma, f, x_{0}, \mathbb{X}_{m}\right) \tag{151}
\end{equation*}
$$

where $\mathbb{X}$ is the finite set of states, $\Sigma$ the finite set of symbols, called the alphabet of the automaton, $x_{0} \in \mathbb{X}$ the initial state, $\mathbb{X}_{m} \subseteq \mathbb{X}$ the set of all accepting states, and $f$ the transition function defined as

$$
\begin{equation*}
f: \mathbb{X} \times \Sigma \rightarrow \mathbb{X} \tag{152}
\end{equation*}
$$

The alphabet represents all possible values that the automaton will recognise. It can consist of letters, values, or any other acceptable symbol. An automaton generates a language, $L$, which simply consists of all possible strings (irrespective of the final state) generated by the automaton. An automaton is said to recognise (or mark) a specific language $L_{m}$.

As well, it is possible to define an output function, $g$, that determines what the output from the automaton will be. In general, it is defined as

$$
\begin{equation*}
g: \mathbb{X} \times \Sigma \rightarrow Y, \tag{153}
\end{equation*}
$$

that is, the output function depends on the state and inputs. In such cases, this automaton is called a Mealy Automaton. On the other hand, when the output function, $g$, is defined as

$$
\begin{equation*}
g: \mathbb{X} \rightarrow Y \tag{154}
\end{equation*}
$$

that is, the output function only depends on the state, then we have a Moore Automaton.
It is also possible to provide a graphical representation of an automaton. Figure 41 shows a typical representation of an automaton. The most important components are:

- The states are shown with " $Z$ " followed by a number.
- Normally, zo is the initial state.
- Each state is enclosed by a circle.
- The initial state is shown by an arrow pointing to the state and without any additional markings.
- The transitions between the states are shown with arrows.
- On the arrows, the combination of inputs and outputs that lead to the next state are shown.
- Accepting states are enclosed by a double circle.


Figure 41: Graphical Representation of an Automaton
The sending and receiving of symbols for the inputs and outputs is assumed to be instantaneous. A finite-state automaton must allocate for each input symbol at each time point a valid transition. Often such a transition can be a self-loop, that is, a transition where the initial and final states are the same.

## Example 16: Automaton for a Process

Consider the previously mentioned process where it was desired to find the word baba. Draw the automaton for this process and define the all the components in the mathematical description of the automaton.

## Solution

The automaton is shown in Figure 42.


Figure 42: Automaton for the example
The mathematical description can be stated as:

$$
\begin{gathered}
\mathbb{X}=\{\mathrm{Z} 0, \mathrm{Z} 1, \mathrm{Z} 2, \mathrm{Z} 3, \mathrm{Z} 4\} \\
\Sigma=\{\mathrm{a}, \mathrm{~b}\} \\
x_{0}=\mathrm{Z} 0 \\
\mathbb{X}_{m}=\{\mathrm{Z} 4\}
\end{gathered}
$$

The transition function, $f$, is defined as follows:

$$
\begin{gathered}
f(\mathrm{Z} 0, a)=\mathrm{Z} 0(\text { which is a self-loop }) \\
f(\mathrm{Z} 0, b)=\mathrm{Z} 1 \\
f(\mathrm{Z} 1, a)=\mathrm{Z} 2
\end{gathered}
$$

$$
\begin{aligned}
& f(\mathrm{Z} 1, b)=\mathrm{Z} 1 \\
& f(\mathrm{Z} 2, a)=\mathrm{Z} 0 \\
& f(\mathrm{Z} 2, b)=\mathrm{Z} 3 \\
& f(\mathrm{Z} 3, a)=\mathrm{Z} 4 \\
& f(\mathrm{Z} 3, b)=\mathrm{Z} 1 \\
& f(\mathrm{Z} 4, a)=\mathrm{Z} 0 \\
& f(\mathrm{Z} 4, b)=\mathrm{Z} 1
\end{aligned}
$$

Note that the transition function should match the arrows drawn in the schematic for the automaton!

Transitions in an automaton can be spontaneous, that is, we do not know exactly when a given transition occurs, for example, the transition from filling a tank to full can be spontaneous (especially, if we do not know the height of the tank). Such a transition is denoted using the symbol " $\varepsilon$ " on the arrow. Spontaneous transitions often arise when there is incomplete knowledge of the system. Automata containing spontaneous transitions are said to be nondeterministic, since it is not possible to know in which states the automaton currently is. Another form of nondeterminism is the presence of multiple transitions with the same label, for example, two transitions labelled a leading to two different states. In such cases, it is likewise impossible to know in which state the automaton is in. In all other cases, the automaton is said to be deterministic since it is possible to precisely determine the next state given all past information.

The states in an automaton can be classified as follows:

1) Periodic States: Periodic states are a set of states between which the automaton can oscillate. Normally, one takes the largest set of states between which the automaton can oscillate.
2) Ergodic States: Ergodic states are those states once reached the automaton cannot leave.
3) Transient States: All states that are not ergodic are called transient states.

## Section 3.4.1.1: Analysis of Automata

With the above mathematical model of an automaton, it is possible to analyse and manipulate the automata. This section will briefly look at some of the ways in which we can manipulate and analyse automata.

Deadlock is said to occur if an automaton ends up in a rejecting state from which it cannot leave. Such an automaton has come to a standstill and cannot take further action. Since automata represent real processes, this is an undesirable state of events and should be avoided. A similar concept is livelock which is said to occur if an automaton ends up in a periodic set that consists solely of rejecting states. This means that although the automaton can make a next move it can never reach an accepting state and terminate. Both such situations are highly undesired. Together deadlock and livelock are referred to as blocking, since the automaton is blocked from completing its task.

## Example 17: Blocking in an Automaton

Determine if the automaton shown in Figure 43, is blocking. If it is blocking, determine what types of blocking occur.


Figure 43: Automaton

## Solution

Looking at Figure 43, state Z3 is an ergodic state which is not accepting. Therefore, deadlock will occur in this state. Thus, the system is blocking. State Z 3 is a deadlock. States Z 5 and Z6 form a periodic set from which there is no escape. However, neither state is accepting. Therefore, livelock will occur for these two states.

It is possible to manipulate automata. The following are some common manipulations and their definitions:

1) Accessible Operator (Acc): This removes all unreachable states and their associated transitions. This operation is necessary when performing more advanced manipulations on automata to clean up the resulting automaton. This operation will not change the generated
or recognised languages. A state is defined as being unreachable if there exists no path from the initial state to the given state.
2) Co-accessible Operator (CoAc): This removes any states and their associated transitions from which one cannot end up in an accepting state. An automaton where $A=\operatorname{CoAc}(A)$ is called a co-accessible automaton and is never blocking. This operation can change the generated language but not the recognised language.
3) Trim Operator (Trim): This operation creates an automaton that is both co-accessible and accessible. The order of operation is immaterial, that is, $\operatorname{Trim}(A)=\operatorname{CoAc}(\operatorname{Acc}(A))=$ $\operatorname{Acc}(\operatorname{CoAc}(A))$.

## Example 18: Trimming an Automaton

Apply the trim operation to the automaton shown in Figure 43.

## Solution

Quickly looking at the automaton in Figure 43, we that there are no states that cannot be accessed from the initial state. Thus, we need to consider the co-accessible operator. Here, we see that we need to remove states $\mathrm{Z} 5, \mathrm{Z} 6$, and Z 3 since we know that they are blocking. However, by removing Z 3 , we now make Z 2 blocking. Therefore, we must also remove it. In general, this process is iterative and continues until we have removed all states or there are no more states to remove. The final automation in shown in Figure 44.


Figure 44: The Trimmed Automaton

## Section 3.4.1.2: Combining Automata

Having looked at various operations on single automata, it is worthwhile to consider manipulating two or more different automata. When we wish to combine two automata, we can perform two different operations: product, denoted by $\times$, and parallel composition, denoted by $\|$. Parallel composition is often also called synchronous composition.

Product composition is defined as combination of two automaton considering only the letters of the alphabet that the two automata have in common, that is, $\Sigma_{1} \cup \Sigma_{2}$. Formally, for two automata $G_{1}$ and $G_{2}$, their product can be written as

$$
\begin{equation*}
G_{1} \times G_{2} \equiv \operatorname{Acc}\left(\mathbb{X}_{1} \times \mathbb{X}_{2}, \Sigma_{1} \cup \Sigma_{2}, f,\left(x_{0_{1}}, x_{0_{2}}\right), \mathbb{X}_{m_{1}} \times \mathbb{X}_{m_{2}}\right) \tag{155}
\end{equation*}
$$

where

$$
f\left(\left(x_{1}, x_{2}\right), e\right) \equiv\left\{\begin{array}{cc}
\left(f_{1}\left(x_{1}, e\right), f_{2}\left(x_{2}, e\right)\right) & \text { if } e \text { is a valid input at the given point }  \tag{156}\\
\text { undefined } & \text { otherwise }
\end{array}\right.
$$

In product composition, the transitions of the two automata are always synchronised on a common event. This implies that a transition occurs only if the input is a valid input for both automata. The states of $G_{1} \times G_{2}$ are given as the pair $\left(x_{1}, x_{2}\right)$, where $x_{1}$ is the current state of $G_{1}$ and $x_{2}$ is the current state of $G_{2}$. It follows from the definition of the product composition that

$$
\begin{align*}
& L\left(G_{1} \times G_{2}\right)=L\left(G_{1}\right) \cap L\left(G_{2}\right) \\
& L_{m}\left(G_{1} \times G_{2}\right)=L_{m}\left(G_{1}\right) \cap L_{m}\left(G_{2}\right) \tag{157}
\end{align*}
$$

Product composition has the following properties:

1) It is commutative up to a re-ordering of the state components in the composed states.
2) It is associative. This implies that $G_{1} \times G_{2} \times G_{3} \equiv\left(G_{1} \times G_{2}\right) \times G_{3}=G_{1} \times\left(G_{2} \times G_{3}\right)$.

## Example 19: Product of Two Automata

Consider the automata shown in Figure 45 and Figure 46. Determine the product of these two automata.


Figure 45: $G_{1}$


Figure 46: $G_{2}$

## Solution

Before we can start with drawing the final automaton, it makes sense to first consider which inputs are valid for the final automaton. From Figure 45, it can be seen that, for $G_{1}$, the inputs are $\{a, b, g\}$. Similarly, from Figure 46, it can be seen that, for $G_{2}$, the inputs are $\{a, b\}$. Therefore, the common set is $\{a, b\}$.

When drawing the final automaton, it helps to start from the initial states of both and work through all the possible transitions and draw the next state. Thus, starting from ( $x, 0$ ), with an input of $a, G_{1}$ remains in state $x$, but $G_{2}$ goes to state 1 . Therefore, the new state in the product automaton will be $(x, 1)$. This is the only valid transition since from $(x, 0)$, an input of $b$ is not valid for $G_{1}$. An input must be valid for both automata for it to occur in the product. Note that $g$ will not occur since it is not a common input. We will then continue in the same fashion by looking at the possible states and how the inputs would affect them. A state will be marked if all the states in the original automaton are marked. The final automaton is shown in Figure 47.


Figure 47: The product of $G_{1}$ and $G_{2}$

As can be seen, product composition is relatively restrictive in that a given input must be valid for both automata. One way to relax this constraint is to consider parallel composition. In parallel composition, the inputs are split into two parts: common inputs and private inputs. A private input only pertains to a given automaton, while common inputs are shared with other automata. Formally, the parallel composition of two automata $G_{1}$ and $G_{2}$, denoted as $G_{1} \| G_{2}$, is defined as

$$
\begin{equation*}
G_{1} \| G_{2} \equiv \operatorname{Acc}\left(\mathbb{X}_{1} \times \mathbb{X}_{2}, \Sigma_{1} \cup \Sigma_{2}, f,\left(x_{0_{1}}, x_{0_{2}}\right), \mathbb{X}_{m_{1}} \times \mathbb{X}_{m_{2}}\right) \tag{158}
\end{equation*}
$$

where
$f\left(\left(x_{1}, x_{2}\right), e\right) \equiv\left\{\begin{array}{cc}\left(f_{1}\left(x_{1}, e\right), f_{2}\left(x_{2}, e\right)\right) & \text { if } e \text { is a valid common input for both automata } \\ \left(f_{1}\left(x_{1}, e\right), x_{2}\right) & \text { if } e \text { is a private event to } G_{1} \\ \left(x_{1}, f_{2}\left(x_{2}, e\right)\right) & \text { if } e \text { is a private event to } G_{2} \\ \text { undefined } & \text { otherwise }\end{array}\right.$
Note that the only difference between parallel and product composition is how the transition function is defined.

Product composition has the following properties:

1) It is commutative up to a re-ordering of the state components in the composed states.
2) It is associative. This implies that $G_{1}\left\|G_{2}\right\| G_{3} \equiv\left(G_{1} \| G_{2}\right)\left\|G_{3}=G_{1}\right\|\left(G_{2} \| G_{3}\right)$.

## Example 20: Parallel Composition of Two Automata

Consider the same two automata as for Example 19. Determine the parallel composition of these two automata.

## Solution

The general procedure for solving such a problem is similar to that of the product composition. First, we need to determine which inputs belong to which categories. Since input $g$ only affects $G_{1}$, it is a private input for $G_{1}$. The other two inputs $\{a, b\}$ are the common inputs to both automata.

Again, we start with the initial states of both automata and work our way through. From the initial state of $(x, 0)$, we have two valid inputs (the common input $a$ and the private input $g$ ). The common input $a$ will bring us as before to the state $(x, 1)$, while the private input $g$ will bring us to the state $(z, 0)$. In the state $(z, 0)$, there are three valid inputs (the common events $a$ and $b$ as well as the private input $g$ ). Input $a$ will bring us to state $(x, 1)$, while input $b$ will be a self-loop. Private input $g$, which only affects $G_{1}$, will bring us to state $(y, 0)$. The final automaton is shown in Figure 48.


Figure 48: The parallel composition of $G_{1}$ and $G_{2}$

## Section 3.4.1.3: Timed Automata

A timed automaton is an automaton that is linked with a clock. A timed automaton has four components: the finite-state automaton, the clock, the invariants, and the guards. Such automata can model minimal and maximal times associated with a given action. Figure 49 shows part of a timed automaton. The clock variable $c$ represents the elapsed time since the last reset, which is denoted as $c \equiv 0$. A guard shows when a transition is activated and can be implemented, for example, in Figure 49, we see that the transition $\mathrm{Z} 0 \rightarrow \mathrm{Z1}$ can only be considered once $c \geq$ 180 s . An invariant shows how long the automaton can remain in a given state, for example, in Figure 49 , we see that the process can remain in state Z 0 until $c=240 \mathrm{~s}$, at which point it must leave the state. The implies that a time automaton has an infinite number of possible realisations, since the system can switch states at any point between 180 and 240 s . Thus, a timed automaton has an infinite state space.

> Condition for the transition $\mathrm{Z} 0 \rightarrow \mathrm{Z} 1$


Figure 49: Timed Automaton

Mathematically, a timed automaton can be represented as

$$
\begin{equation*}
A \equiv\left(\mathbb{X}, \Sigma, f, \mathbb{C}, I, x_{0}, \mathbb{X}_{m}\right) \tag{160}
\end{equation*}
$$

where $\mathbb{C}$ is a finite set of clock variables and $I$ is the invariant function that links the states with the transition constraints, that is,

$$
\begin{equation*}
I: \mathbb{X} \rightarrow \Phi(\mathbb{C}) \tag{161}
\end{equation*}
$$

where $\Phi(\mathbb{C})$ is the set of transition constraints $\delta$. The transition constraints $\delta$ are always either true or false based on the current clock variables. The allowed transition constraints are (where $a \in \mathbb{R}$ ):

- $\delta \equiv(c \leq a)$
- $\delta \equiv(c=a)$
- $\delta \equiv(c \geq a)$
- $\delta \equiv(\delta 1$ OR $\delta 2$ )
- $\delta \equiv \neg(\delta)$
- $\delta \equiv \emptyset$ or $\}$

The transition function is defined as

$$
\begin{equation*}
f: \mathbb{X} \times U \times \Phi(\mathbb{C}) \rightarrow \mathbb{X} \times 2^{|\mathbb{C}|} \tag{162}
\end{equation*}
$$

where $2^{|C|}$ represents the power set of $\mathbb{C}$.
In certain formalisms, transitions can be specified as either urgent or nonurgent. Urgent transitions will always be taken as soon as possible, while nonurgent transitions can wait. Normally, spontaneous transitions are assumed to be nonurgent.

## Section 3.5: Chapter Problems

Problems at the end of the chapter consist of three different types: (a) Basic Concepts (True/False), which seek to test the reader's comprehension of the key concepts in the chapter; (b) Short Exercises, which seek to test the reader's ability to compute the required parameters for a simple data set using simple or no technological aids. This section also includes proofs of theorems; and (c) Computational Exercises, which require not only a solid comprehension of the basic material, but also the use of appropriate software to easily manipulate the given data sets.

## Section 3.5.1: Basic Concepts

Determine if the following statements are true or false and state why this is the case.

1) A process satisfying the principles of superposition and homogeneity is said to be linear.
2) In a time-variant model, the parameters themselves vary with respect to time.
3) In a lumped-parameter model, there are space derivatives.
4) A noncausal system depends on future values.
5) A system with memory only cares about the current value of the process.
6) A state-space representation provides a model linking states, inputs, and outputs.
7) A transfer function can only be written for linear processes.
8) Every transfer function has a unique state-space representation.
9) Prediction error models are discrete-time models of the process.
10) A white noise signal depends on past values of the noise.
11) In a Box-Jenkins model, the order of the $A$-polynomial is fixed to zero.
12) In an autoregressive exogenous model, the only orders of the $C$ - and $D$-polynomials are zero.
13) It is not possible to convert a continuous model into a discrete model.
14) A process at steady state will experience wild, unpredictable swings in values.
15) The gain of a process represents the transient behaviour of the process.
16) The process time constant represents the delay before a process responds.
17) A continuous transfer function with poles of $-2,-1$, and 0 is stable.
18) A continuous transfer function with poles of 1,2 , and 5 is stable.
19) A discrete transfer function with poles of $0.5,-0.5$, and 1 is unstable.
20) A discrete transfer function with poles of $0.25,0.36,0.25 \pm 0.5 i$ is stable.
21) A continuous state-space model with eigenvalues of $0.25 \pm 2 i$ is stable.
22) A discrete state-space model with eigenvalues of $\pm 0.25 i$ is stable.
23) The alphabet of an automaton represents the allowed inputs into the process.
24) An accepting state is a state that has the desired outcome.
25) The language recognised by an automaton is the set of all words accepted by the automaton.
26) In a Mealy automaton, the output function depends only on the states.
27)Blocking occurs when an automaton cannot reach an accepting state.
27) We denote spontaneous transitions by $\varepsilon$.
28) An automaton with spontaneous transitions is called a deterministic automaton.
29) The co-accessible operator removes all states and their associated transitions that are unreachable from the initial state.
30) In a timed automaton, a guard determines the maximal time in which we can remain in a given state.
32)For a timed automaton, $\delta \equiv(c \leq 1.000)$ is an acceptable time constraint.
33)For a timed automation, $\delta \equiv$ ( $c=a b c d)$ is an acceptable time constraint.
31) There is no escape from an ergodic set.

## Section 3.5.2: Short Questions

These questions should be solved using pen and paper. Appropriate software for drawing the required diagrams can also be used to assist with the design.
35)For the following models, classify them based on the information in this chapter. Are the models linear, time-invariant, lumped parameter, memoryless, or causal?
a. $\quad y_{k+1}=4 y_{k}+7 u_{k+5}-5 e_{k}$.
b. $\frac{\partial^{2} T}{\partial x^{2}}=-\alpha(t) \frac{\partial T}{\partial t}$, where $\alpha$ is a parameter that depends on time.
c. $y_{k+1}=-4 y_{k}-3 e_{k}$
d. $\frac{\partial T}{\partial t}=-\alpha(t) u(t+1)$, where $\alpha$ is a parameter that depends on time.
36)For the following continuous-time transfer function, determine their stability. For the stable transfer functions, determine the gain, time constant, and time delay.
a. $\quad G(s)=\frac{5(s+1)}{6 s^{3}+11 s^{2}+6 s+1} e^{-3 s}$
b. $\quad G(s)=\frac{-5}{150 s^{3}+65 s^{2}+2 s-1} e^{-10 s}$
c. $\quad G(s)=\frac{4.5}{15 s^{2}+8 s+1} e^{-10 s}$
d. $\quad G(s)=\frac{4.5}{15 s^{2}+26 s+7} e^{-7 s}$
37)For the following continuous-time state-space models, determine their stability. Convert the models to a transfer function. For the stable models, determine their gain and time constant.
a. $\frac{d \vec{x}}{d t}=\left[\begin{array}{cc}5 & 0 \\ 0 & -2\end{array}\right] \vec{x}+\left[\begin{array}{l}2 \\ 1\end{array}\right] \vec{u}$
$\begin{aligned} y & =\left[\begin{array}{ll}1 & 0 \\ 0 & 1\end{array}\right] \vec{x} \\ \frac{d \vec{x}}{d t} & =\left[\begin{array}{cc}-5 & 0 \\ 0 & -2\end{array}\right] \vec{x}+\left[\begin{array}{c}-2 \\ 1\end{array}\right] \vec{u}\end{aligned}$
b.

$$
\begin{aligned}
y & =\left[\begin{array}{ll}
2 & 0 \\
0 & 1
\end{array}\right] \vec{x} \\
\frac{d \vec{x}}{d t} & =\left[\begin{array}{ccc}
-2 & 1 & 2 \\
0 & -3 & 2 \\
0 & 0 & -1
\end{array}\right] \vec{x}+\left[\begin{array}{c}
1 \\
-0.5 \\
2
\end{array}\right] \vec{u}
\end{aligned}
$$

$$
y=\left[\begin{array}{ccc}
2 & 0 & 0 \\
0 & -1 & 0 \\
0 & 0 & 1
\end{array}\right] \vec{x}
$$

38)For the following discrete-time models, determine their stability:
a. $y_{k+1}=4 y_{k}+7 u_{k+5}-5 e_{k}$
b. $y_{k+1}=\frac{z^{-5}}{1-4 z^{-1}} u_{k}$
c. $y_{k+1}=\frac{z^{5}+z^{4}}{z^{6}+z^{5}+z^{4}+z^{3}+z^{2}+z^{1}+1} u_{k}$

$$
\vec{x}_{k+1}=\left[\begin{array}{ccc}
-0.5 & 1 & 2 \\
0 & 0.5 & 2 \\
0 & 0 & 0.25
\end{array}\right] \vec{x}_{k}+\left[\begin{array}{c}
1 \\
-0.5 \\
2
\end{array}\right] \vec{u}_{k}
$$

d.

$$
\vec{y}_{k}=\left[\begin{array}{ccc}
2 & 0 & 0 \\
0 & -1 & 0 \\
0 & 0 & 1
\end{array}\right] \vec{x}_{k}
$$

$$
\begin{aligned}
& \vec{x}_{k+1}=\left[\begin{array}{ccc}
-1.5 & 0 & 0 \\
3 & 1.5 & 0 \\
1 & 2 & 2.25
\end{array}\right] \vec{x}_{k}+\left[\begin{array}{c}
-1 \\
0.5 \\
3
\end{array}\right] \vec{u}_{k} \\
& \text { e. } \vec{y}_{k} \\
&=\left[\begin{array}{lll}
1 & 0 & 0 \\
0 & 1 & 0 \\
0 & 0 & 1
\end{array}\right] \vec{x}_{k}
\end{aligned}
$$

39) What is the relationship between the eigenvalues of the state-space model and the time constant (as determined from the transfer function)?
40)For the automata shown in Figure 50, classify the states into marked, ergodic, periodic, and transient. Determine if there is blocking. If present, give the type of blocking encountered.



Figure 50: Automata for Questions 40 and 41
41)Perform the trim operation on the automata in Figure 50
42) Using the automata given in Figure 51, perform the following operations: $K_{1} \times K_{2}, K_{1} \| K_{2}$, $\mathrm{K}_{1} \times \mathrm{K}_{3}, \mathrm{~K}_{1}\left\|\mathrm{~K}_{2}\right\| \mathrm{K}_{3}$, and $\mathrm{K}_{1} \times \mathrm{K}_{2} \times \mathrm{K}_{3}$.



Figure 51: Automata for Question 42
43)Draw the automata for the following processes:
a. Given the letters $a$ and $b$, find the string $a b a b$.
b. Given the letters $c, d$, and $e$, find the string $d e c d$.
c. Given the letters $g, h$, and $i$, find the strings hig and high.

## Section 3.5.3: Computational Exercises

The following problems should be solved with the help of a computer and appropriate software packages, such as MATLAB ${ }^{\circledR}$ or Excel ${ }^{\circledR}$.
44) Model a system that you are familiar with. Be sure to include all the differential equations required for the process.

## Chapter 4: Schematic Representation of a

## Process

This section will describe the most common schematic methods for representing a process, including block diagrams, piping and instrumentation diagrams (P\&IDs), process flow diagrams (PFDs), and electrical and logic circuit diagrams.

## Section 4.1: Block Diagrams

A block diagram is an abstract way of representing a system in the frequency domain that allows for all the messy details to be hidden and only the essential elements shown. ${ }^{13}$ The basic block diagram consists of three parts, as shown in Figure 52. On the left, entering the block diagram, is the input, denoted by $U$, while on the right, leaving the block diagram, is the output, denoted by $Y$. Inside the block, the process model, denoted by $G$, is shown. In most cases, the exact form of the process model will not be specified, but it can be any relationship between the input and output.


Figure 52: The basic block diagram
Another common block is the summation block, which shows how two or more signals are to be combined. Figure 53 shows a typical summation block. The signs inside the circle show whether the signal coming in should be added or subtracted. Given the common nature of a summation block, two simplifications can be made. Rather than using a full circle, the signals are simply shown to connect and the signs are shown beside each signal. This is shown in the bottom part of Figure 53. A further simplification is to completely ignore any positive signs and only give the negative signs beside the appropriate signals. A commonly used standard is to place the summation signs on the left-hand side of the arrow.

[^10]

Summation block that subtracts stream $U_{2}(\mathrm{~s})$ from stream $U_{1}(\mathrm{~s})$ to give $Y(\mathrm{~s})$.

Summation block that subtracts stream $U_{2}(\mathrm{~s})$ from stream $U_{1}(\mathrm{~s})$ to give $Y(\mathrm{~s})$.

Figure 53: Summation block: (top) full form and (bottom) short-hand equivalent
One of the most useful features of block diagrams is the ability to easily compute the relationship between the different signals. For the basic block shown in Figure 52, the relationship between the input and output can be written as:

$$
\begin{equation*}
Y=G U \tag{163}
\end{equation*}
$$

while for the summation block, it can be written as

$$
\begin{equation*}
Y=U_{1}-U_{2} \tag{164}
\end{equation*}
$$

Now, consider the process shown in Figure 54, where it is desired to determine the relationship between $Y$ and $U$, where there are three blocks in series. The naïve approach would be to denoted each of the outputs from the two intermediate steps to be denoted as $Y_{1}$ and $Y_{2}$ and then write relationships between each of these variables to obtain the final result, that is,

$$
\begin{gather*}
Y_{1}=G_{1} U  \tag{165}\\
Y_{2}=G_{2} Y_{1}=G_{2} G_{1} U  \tag{166}\\
Y=G_{3} Y_{2}=G_{3} G_{2} G_{1} U \tag{167}
\end{gather*}
$$

This shows that the final relationship can be written as $Y=G_{3} G_{2} G_{1} U$. The easier approach is to note that since these process blocks are in series, the process models can be multiplied together to give the desired result (as the final result shows). However, the naïve approach is very useful if the system is very complex with many different summation blocks and functions.


Figure 54: Block diagram algebra: In order to relate $U$ and $Y$, the transfer functions between the two points need to be multiplied, thus, $Y=G_{3} G_{2} G_{1} U$.

## Example 21: Complex Block Diagrams

Consider the closed-loop system shown in Figure 55 and derive the expression for the relationship between $R$ and $Y$. Assume that all signals are in the frequency domain.


Figure 55: Generic closed-loop, feedback control system

## Solution:

Starting from the desired signal $R$ and moving towards $Y$, the following relationships can be written

$$
\begin{gathered}
\varepsilon=R-G_{m} Y \\
U=G_{c} \varepsilon \\
Y=G_{p} G_{a} U+G_{d} D
\end{gathered}
$$

Substituting the first relationships into the second and then the combined into the third gives

$$
Y=G_{p} G_{a} G_{c}\left(R-G_{m} Y\right)+G_{d} D
$$

Re-arranging this equation gives

$$
Y=G_{p} G_{a} G_{c} R-G_{p} G_{a} G_{c} G_{m} Y+G_{d} D
$$

Solving this equation for $Y$ gives

$$
Y=\frac{G_{p} G_{a} G_{c}}{1+G_{p} G_{a} G_{c} G_{m}} R+\frac{G_{d}}{1+G_{p} G_{a} G_{c} G_{m}} D
$$

Since we are only interested in the relationship between $Y$ and $R$, we can set $D=0$ to get

$$
Y=\frac{G_{p} G_{a} G_{c}}{1+G_{p} G_{a} G_{c} G_{m}} R
$$

These equations are commonly encountered in closed-loop control.

The creation of block diagram for a given process can be quite a complex task, but it is worthwhile as it allows the essential features of the system to be abstracted out and understood.

## Section 4.2: Process Flow Diagrams

The process flow diagram is a simplified diagram of the process, where only the key components and connections are shown. For complex processes, the process flow diagram is often created using blocks, where the blocks show a subprocess. In such cases, they can approximate the block diagrams described in Section 4.1.

A process flow diagram for a single process normally contains the following elements: process piping, key components, key valves and control valves, connections to other systems, key bypass and recycle streams, and process flow names.

Figure 56 shows a typical process flow diagram. The rules for creating a process flow diagram are similar to the rules for creating a piping and instrumentation diagram, given in Section 4.3. The only difference is the level of detail.


Figure 56: Process Flow Diagram for Alkylate Splitter

## Section 4.3: Piping and Instrumentation Diagrams (P\&ID)

The piping and instrumentation diagram (P\&ID) is a detailed description of the process, where all connections and components are shown. A P\&ID contains, in addition to the information found in a process flow diagram, the following information:

1) Type and Identification number for all components
2) Piping, fittings with nominal diameters, pressure stages, and materials
3) Drives
4) All measurement and control devices

Figure 57 shows a typical P\&ID for part of a chemical plant.


Figure 57: P\&ID for a Gas Chilling and Separation Plant According to Canadian Design Standards (Note the engineering stamp in the bottom middle box.)

## Section 4.3.1: P\&ID Component Symbols According to the DIN EN

 62424Table 19 shows the typical component symbols according to the DIN EN 62424 Standard. ${ }^{14}$ Additional symbols can be found in the standard itself.

Table 19: Component Symbols for P\&IDs According to the DIN EN 62424

| Symbol | Name |
| :---: | :---: | :---: | :---: |
| Pipe ${ }^{15}$ |  |
| Insulated pipe |  |
| Jacketed pipe |  |
| Int |  |

[^11]Name
Axial fan
Radial fan
Furnace


## Section 4.3.2: Connections and Piping in P\&IDs

The type of connection must be shown in a P\&ID. Table 20 shows the most common possibilities for such connections.

Table 20: Connections Types for P\&IDs

| Symbol | Name |
| :---: | :---: |
| $\begin{aligned} & H \quad H \quad H \\ & --------- \\ & L \quad L \quad L \end{aligned}$ | Pipe (process flow) <br> Pneumatic signal <br> Electrical signal <br> Hydraulic signal |


| Symbol | Name |
| :---: | :---: |
| $\forall 8$ | Electromagnetic Signal |

## Section 4.3.3: Labels in P\&IDs

Another important feature of $\mathrm{P} \& \mathrm{IDs}$ is that the various components must be clearly identified. Sensors, valves, and other actuators should be named according to different guidelines. As well, the type and location of the components must be clearly shown. Table 21 and Table 22 shows the symbols that are used to show the location of actuators and sensors.

Table 21: Location Symbols

| Location | Local/in the field | In a central location | In a local central point |
| :---: | :---: | :---: | :---: |
| Symbol | $\begin{gathered} \hline \text { PI } \\ 123.4 \\ \hline \end{gathered}$ | $\begin{gathered} \text { PI } \\ 123.4 \\ \hline \end{gathered}$ | $\begin{gathered} \mathrm{PI} \\ 123.4 \\ \hline \end{gathered}$ |
| Comments | Component is found in the neighbourhood of the process itself. | Component is found in some central location. Often, this central location is a computer. | Component is found in some local central point, for example, a process control cabinet. |

Table 22: Type Symbols

|  | General | Process Control Function |
| :--- | :---: | :---: |
| Symbol | PI |  |
| 123.4 |  |  |

There are many different fields in the equipment label. These are shown in Figure 58. The left fields (\#1 to 3 ) are only used if needed and there are no restrictions on what can be placed there. Field \#1 often gives the supplier, while Field \#2 gives the standard value of the component. The two central fields (\#4 and 5) show the important information about the component. Field \#4 shows the PCE category ${ }^{16}$ und the PCE processing category (see Table 23 for additional information). Field \#5 shows the PCE tag that is arbitrary. The right fields (\#6 to 12) show

[^12]additional information about the component. For example, Fields \#6 to \#8 show alarms and notifications related to upper limits, while Fields \#10 to 12 contain alarms and notifications related to lower limits. The fields located furthest from the center contain the most important information. Field \#9 shows the importance of the component. A triangle (A) shows a safety-relevant component, while a circle $(\bullet)$ shows a component required for the good manufacturing process (GMP). A square (■) shows a quality relevant component.


Figure 58: Fields in a P\&ID Label
Table 23 shows the PCE categories, while Table 24 shows the PCE processing functions. For motorised drives (PCE category N ), there are only two possibilities for the processing function: S, which implies an on/off motor, and C, which implies motor control. For valves (PCE category Y), there are also limitations: S implies an on/off valve, C implies a control valve, Z implies an on/off valve with safety relevance, and IC implies a control valve with a continuous position indicator. For process control functions, there are some very specific rules:

1) The PCE category is always the same $U$.
2) The following letters are often $\mathrm{A}, \mathrm{C}, \mathrm{D}, \mathrm{F}, \mathrm{Q}, \mathrm{S}, \mathrm{Y}$, or Z . Combinations of these letters are possible.

| Letter | Meaning |
| :---: | :---: |
| A | Analysis |
| B | Burner Combustion |
| C $^{17}$ | Conductivity |
| D | Density |
| E | Voltage |
| F | Flow |
| G | Gap |
| H | Hand |
| I | Current |
| J | Power |
| K | Time Schedule |
| L | Level |
| M | Moisture |
| N | Motor |
| O | Free |
| P | Pressure |
| Q | Quantity/Event |
| R | Radiation |
| S | Speed, Frequency |
| T | Temperature |
| U | anticipated for PCE control functions |
| V | Vibration |
| W | Weight |
| X | Free |
| Y | Valve |
| Z | Free |

Table 24: PCE Processing Categories

| Letter | Meaning | Comment |
| :---: | :---: | :---: |
| A | Alarming | only in Fields $\# 6,7,8,10,11$, and 12 ; in Field \#4 for process control functions |
| B | Condition, Limitation | Field \#4 for process control functions only in Field 4 |
| C | Control | only in Field 4 |
| D | Difference | only in Field 4 |
| F | Fraction | only in Field 4 |
| H | High, on, open | only in Fields \#6,7, 8, 10, 11, and 12 |
| I | Indicator | only in Field 4 |
| L | Low, off, closed | only in Fields \#6,7, 8, 10, 11, and 12 |
| O | Local or PCS status indictor from a binary signal | only in Fields \#6,7, 8, 10, 11, and 12 |

[^13]| Letter | Meaning | Comment |
| :---: | :--- | :--- |
| Q | Quantity | only Field 4 |
| R | Recording | only Field 4 |
| S | Switching | only in Fields \#6,7, 8, 10, 11, and 12; in <br>  <br> Y |
| Z | Computation | Fmergency | | Field \#4 for process control functions |
| :--- |
| only in Field 4 |
| only in Fields \#6,7, 8, 10, 11, and 12; in |

## Example 22: P\&ID Tags

What is the meaning of the following P\&ID tags: PI-512, UZ-512 und MDI-512?

## Solution

For PI-512, the first letter is "P". From Table 23, we can see that "P" represents pressure. The following letter is "I". From Table 24, we can see that "I" represents indicator. Thus, "PI-512" represents a pressure indicator.

For UZ-512, the first letter is "U". From Table 23, we can see that "U" represents a process control function. The following letter is " $Z$ ". From Table 24 , we can see that " $Z$ " represents a safety-critical control function. Thus, "UZ-512" represents a safety-critical control function that is computed using a computer/PLC.

For MDI-512, the first letter is "M". From Table 23, we can see that "M" represents moisture. The following letters are "DI". From Table 24, we can see that "D" represents difference and "I" represents indicator. Thus, "MDI-512" represents a moisture difference indicator.

The labels for the other fields are not standardised. However, in a P\&ID, the same label should be used for the same component/idea. All P\&IDs require a legend, where all the components are briefly described, including such information as their name, engineering data (size, material), and operating conditions. Appropriate stamps and signatures may need to be affixed depending on local laws, for example, in Canada, all engineering documents require a stamp and signature from a practising engineer.

## Section 4.4: Electric and Logic Circuit Diagrams

Since much automation is implemented using electrical signals and logic statements, it is useful to understand the basics of how such circuit diagrams are constructed. As well, some of these symbols will re-appear when considering graphical programming languages.

There exist two main standards for drawing the shapes in electric and logic circuit diagrams: the official DIN EN 60617 (a.k.a. the IEC 617) standard and the nonpreferred, but commonly encountered, ANSI IEEE 91-1991 standard. The official standard, although not explicitly showing the nonpreferred symbols, states that it is permissive to use locally accepted national versions of the symbols. Common symbols for both standards are shown in Table 25.

Furthermore, when it comes to showing connections, for example, the splitting of a wire, different formats are possible. The recommended approach is to use T-junctions to show a split (as shown by Figure 59a). However, often, the split will be shown as in Figure 59b. A simple crossing of wires is shown by Figure 59c.

Table 25: Common Symbols in Circuit Diagrams

| Item | DIN-60617 Symbol | ANSI Symbol | Comments |
| :---: | :---: | :---: | :---: |
| Resistor | $-\square-$ | $-M-$ |  |
| Inductor | -mm | 3002 |  |
| Direct-Current <br> Voltage Source | $-1+$ | $-1+$ | Essentially a battery |
| Alternating- |  |  |  |
| Current Voltage |  | $\sim$ |  |
| Source |  |  |  |
| Diode |  |  |  |
| Capacitor |  | $-\mathrm{H}$ |  |
| AND Gate |  |  |  |


| Item | DIN-60617 Symbol | ANSI Symbol | Comments |
| :---: | :---: | :---: | :---: |
| OR Gate |  |  |  |
| Inverter |  |  |  |
| NAND Gate |  |  |  |
| NOR Gate |  |  |  |
| XOR Gate |  |  |  |
| Ground | $\stackrel{\perp}{=}$ | $\stackrel{\perp}{=}$ |  |
| Variable | $y$ | $!$ | The arrow is placed over the element that is variable, for example, a variable |
| General Negation | $\bigcirc$ | $\bigcirc$ | resistor: <br> Negation is placed where the signal enters or leaves a block (see, for example, the NAND or NOR gates above). |



Figure 59: Connections: a) recommend form for contact; b) commonly encountered form for contact; and c) no contact

## Section 4.5: Chapter Problems

Problems at the end of the chapter consist of three different types: (a) Basic Concepts (True/False), which seek to test the reader's comprehension of the key concepts in the chapter; (b) Short Exercises, which seek to test the reader's ability to compute the required parameters for a simple data set using simple or no technological aids. This section also includes proofs of theorems; and (c) Computational Exercises, which require not only a solid comprehension of the basic material, but also the use of appropriate software to easily manipulate the given data sets.

## Section 4.5.1: Basic Concepts

Determine if the following statements are true or false and state why this is the case.

1) A block diagram is a frequency-domain representation of the process.
2) A pneumatic signal is denoted using a line with sinusoids.
3) An electric signal is denoted using a dashed line.
4) A tag placed in a round symbol with a single line through the middle represents a component located in some central location.
5) An instrument with the tag TIC is a density-indicator controller.
6) An instrument with the tag PIC is a pressure-indicator controller.
7) An instrument with the tag RI is a radiation indicator.
8) An instrument with the tag TDI is a temperature-difference indicator.
9) An instrument with the tag JI is a current indicator.
10) A P\&ID should contain a legend explaining the symbols and notation used.

## Section 4.5.2: Short Questions

These questions should be solved using pen and paper. Appropriate software for drawing the required diagrams can also be used to assist with the design.
11) Write the process model between $U$ and $Y$ for the block diagrams shown in Figure 60.


Figure 60: Block Diagrams for Question 11
12) Draw the process flow diagrams for the following process descriptions. Hint: Knowledge of chemical engineering and physical behaviour of systems required.
a. The DEA solution enters the treatment unit via a throttle valve and is heated in three consecutive heat exchangers. The heated solution enters the first evaporation column which is operated under vacuum at 80 kPa and a temperature of $150^{\circ} \mathrm{C}$. Most of the water is evaporated here. The water vapour is used in the first heat exchanger to heat the incoming solution. The concentrated DEA solution is transferred from the first evaporator to a second evaporation column via a second throttle valve, so that the pressure can be lowered to 10 kPa . The second evaporation column is a thin film evaporator (similarly built to a scraped wall heat exchanger), which is heated by high pressure steam. Here most of the DEA is evaporated. The remaining HSS slurry is removed from the column by a positive displacement pump. The DEA vapour is used to heat the feed in the second heat exchanger. The third heat exchanger in the feed line is heated by medium pressure steam to achieve sufficient evaporation in the first evaporation column. After exchanging heat with the feed stream, the water vapour and DEA vapour streams
are completely liquefied in two condensers operated with cooling water, before being pumped to a T -junction and returned to the gas scrubbing process.
b. A concentrated solution of $99 \%$ triethylene glycol $\left(\mathrm{C}_{6} \mathrm{H}_{14} \mathrm{O}_{4}\right)$ and $1 \%$ water is commonly used to absorb moisture from natural gas at 4.1 MPa . Natural gas is mainly methane, with smaller amounts of ethane, propane, carbon dioxide and nitrogen. The water-saturated natural gas is contacted with the triethylene glycol at $40^{\circ} \mathrm{C}$ in a trayed absorber column to remove the water. The dried gas in then ready for piping to market. The triethylene glycol solution leaving the absorber column is diluted by water, and must be regenerated for re-use. It first passes through a flash drum at a pressure of 110 kPa to release dissolved gases. Then, it is heated in a heat exchanger with the hot regenerated glycol. The dilute solution then enters the top of a packed regeneration column. As it flows downward over the packing, it passes steam flowing upwards which evaporates water from the glycol solution. The reboiler at the bottom of the regeneration column uses medium-pressure steam to heat the triethylene glycol solution to $200^{\circ} \mathrm{C}$ to generate the vapour that passes up the column. At the top of the column, the vapour (i.e. the water evaporated from the triethylene glycol solution) is vented to the atmosphere. The hot regenerated solution from the bottom of the regenerator column is cooled in a heat exchanger with the dilute solution from the absorber. It is then pumped through an air cooler and then back to the absorber column at approximately $45^{\circ} \mathrm{C}$.
13) Using the sketch of the process flow diagram in Figure 61 and the process description below, provide a properly formatted process flow diagram that includes all the required components and information. The process description is as follows. Ethyl chloride $\left(\mathrm{C}_{2} \mathrm{H}_{5} \mathrm{Cl}\right)$ is produced in a continuous, stirred tank reactor (CSTR) filled with a slurry of catalyst suspended in liquid ethyl chloride. Most of the heat of reaction is absorbed by vaporising $25 \mathrm{kmol} / \mathrm{hr}$ of the liquid ethyl chloride. This vapour leaves the reactor with the product stream. The reactor is jacketed to allow additional temperature control if required. All of the ethyl chloride in the product stream is condensed and enough ethyl chloride is returned to the reactor to maintain steady state. The waste gas is sent to the flare system.


Figure 61: Sketch of Process Flow Diagram for Question 13
14) Evaluate the P\&IDs shown in Figure 62. Have they been corrected drawn? Is there anything that is missing? Are all the streams correctly shown?


Figure 62: P\&IDs for Question 14
15) Given the P\&ID in Figure 63, determine where the controller is located and what signals it requires.


Figure 63: P\&ID for Question 15
16)For the process shown in Figure 64, answer the following questions:
a. What kind of alarm does the controller provide?
b. How is the level measured? What kind of signal is produced?
c. What kind of valve is used to control the level?
d. What do the double lines for item LI-135 mean?


Figure 64: P\&ID for Question 16
17) Imagine that you producing maple syrup. The $P \& I D$ along with the current values is shown in Figure 65. Determine if there are any measurement errors.


Figure 65: Maple Syrup P\&ID for Question 17

## Section 4.5.3: Computational Exercises

The following problems should be solved with the help of a computer and appropriate software packages, such as MATLAB ${ }^{\circledR}$ or Excel ${ }^{\circledR}$.
18) Take a complex process and draw the P\&ID for it. Make sure to include all the relevant sensors and controllers.
19) Using the $P \&$ ID from Question 18, create a simplified process flow diagram for the process.

# Chapter 5: Control and Automation 

## Strategies

Having looked at the different instruments that can be used for automation and how to describe the process and its interconnections, it is now necessary to see how the desired automation can actually be implemented. The control or automation strategy is the method used to control/automate the process to achieve the desired objectives. Although two different words can be used here, the overall concepts are very similar: produce a system that can operate on its own with minimal human interference. Depending on the field and even industry, one of these two words may be more common.

There exist many different ways to classify the different types of strategies that can be implemented. Often, in practice, these strategies can be combined to produce a final overall control strategy. The following are the most common types of control strategies:

1) Open-Loop Control (or steering, servo response): In open-loop control, the desired trajectory is set and implemented. The object follows this trajectory without taking into consideration any variations in the environment or surroundings. Clearly, if there are any changes in the environment, the object may not achieve its trajectory.
2) Closed-Loop Control (or regulation, regulatory response): In closed-loop control, the actual value of the system is always compared against the setpoint value. Any deviations are then corrected using a controller. This allows for the system to make corrections based on imperfections in the original specifications, changes in the environment, or unexpected events. In most control strategies, some aspect of closed-loop control will be implemented.
3) Feedforward Control: In feedforward control, known disturbances are measured and corrective action is taken so that the disturbance does not impact the plant. This is a way of taking into consideration future information about the process and taking action now.
4) Discrete-Event Control: In discrete-event control, control only occurs, when some logic condition is triggered. This is often used in safety-relevant systems to trigger immediate action or raise an alarm, for example, if the pressure is above a certain threshold, then the pressure release valve should be opened. Often, such systems are modelled using automata.
5) Supervisory Control: In supervisory control, the control loop does not control a process, but it controls another control loop. Supervisory control loops are common in complex industrial systems, where there may be multiple objectives and variables to control.

## Section 5.1: Open- and Closed-Loop Control

Since the two control strategies are often combined together, it is useful to understand the difference between them.

## Section 5.1.1: Open-Loop Control

Consider a system as shown in Figure 66, where the manipulated variable $u(t)$ influences the output variable, $y(t)$. The desired behaviour is specified by the setpoint $r(t)$. The path from $r(t)$ to $u(t)$ to $y(t)$ is called the open-loop control path. The objective here is to design a controller, $G_{c}$, such that given a setpoint trajectory, it can produce a sequence of value for the manipulated variable $u(t)$, so that the output $y(t)$ attains the values specified by $r(t)$. The disturbance affecting the system is given as $e(t)$. Often, it will be assumed that $e(t)$ is a Gaussian, white noise signal.


Figure 66: Open-Loop Control
In principle, if we are given a mathematical function that converts $u(t)$ into $y(t)$, more commonly called a plant model and denoted by $G_{p}$, then by taking the inverse of this model as the controller will allow us to attain the goal of making $y(t)=r(t)$. However, in practice, the following issues arise:

1. Unrealisability of the Model Inverse: Due to time delays and unstable zeros of the process, it may not be possible to obtain a realisable inverse model.
2. Plant-Model Mismatch: It is very likely that the mode for $G_{p}$ is not exact, which implies that the incorrect parameters are given to the controller leading to incorrect results.
3. Disturbances: In reality, there will be disturbances that prevent $y(t)$ from reaching the value specified by $r(t)$. One solution to this problem is to implement feedforward control
that measures the disturbance and takes corrective action. However, this only solves the problem if the disturbance variable can be measured.

Example 23: Heating the House: Part I: Open-Loop Control


Figure 67: Open-Loop Control for the Temperature in the House
Figure 67 shows a schematic diagram for the open-loop control of the temperature in the house, where the radiator setting is used to determine the temperature in the room. However, once the radiator setting is set, no further actions are taken if a disturbance should occur, for example, the temperature outside changes, more people enter the room, the window is opened, or the sun starts shining. All of these disturbances will change the temperature in the room, but no automatic change will occur in radiator setting. Of the disturbances, the largest is probably the outside temperature. If, as shown in Figure 67, we can measure the outside temperature, then we can design a feedforward controller to correct for the disturbance caused by this variable.

## Section 5.1.2: Closed-Loop Control

In closed-loop control, information about the process is used to update the control action. Figure 68 shows the overall closed-loop control system, where the difference between the reference signal and the actual output is used as the input into the controller. The controller error, $\varepsilon(t)$, is defined as the difference between the setpoint and the measured output. This error is then
the input into the controller. In order to emphasise that the output must be measurable in order for closed-loop control to occur, the measuring transfer function, $G_{m}$, is explicitly shown. In most applications, it will be assumed that $G_{m}$ has a minimal impact on the overall system, and hence, it will be ignored.


Figure 68: Closed-Loop Control
Closed-loop control provides the following advantages over open-loop control:

1) Better accuracy, especially since it can compensate for unknown and unmeasurable disturbances or imprecise models of the system.
2) Stability, since closed-loop control can stabilise unstable processes.

On the other hand, closed-loop control requires the ability to measure the output variables, which can lead to a greater cost. As well, a poorly designed closed-loop controller can destabilise the system.

## Example 24: Temperature Control: Closed-Loop Case

Consider the same set-up as in Example 23, but modified as shown in Figure 69, so that the temperature in the room can now be measured and an appropriate controller designed. Based on the temperature in the room, the controller can set the valve opening which will determine the flow of heat into the room. Opening the valve more will cause more heat to enter the room; thus, increasing its temperature. Closing the valve will cause less heat to enter the room; thus, decreasing its temperature.


Figure 69: Closed-Loop Control of the House Temperature
Furthermore, other disturbances may occur that are not easily measured, but whose impact can be controlled by the closed-loop controller. Figure 69 shows people in the room (S1), an open window (S2) and direct sunlight (S3). Since these will influence the room temperature, their impact will be noticed by the room temperature sensor that will cause the controller to make appropriate changes to maintain the temperature at its desired value.

In closed-loop control, the following are the key ideas to consider (Seborg, Edgar, Mellichamp, \& Doyle, 2011):

1) The closed-loop system should be stable; otherwise, the system is worse than before.
2) The impact of disturbances should be minimised.
3) The controller should provide quick and effective change between different setpoints, that is, setpoint tracking should be good.
4) There should not be any bias, that is, the output should reach the desired value.
5) Large actuator values should be minimised.
6) The controller should be robust, that is, small changes in the process characteristics should not make the overall system unstable.

When designing a closed-loop control strategy, there are two common approaches to take:

1) State Feedback, where it is assumed that the controller is a constant $K$ and all the states are measurable and available, that is, $y_{t}=x_{t}$. The control law is given as $u_{t}=-K x_{t}$.
2) Proportional, Integral, and Derivative (PID) Control, where the control law can be written as

$$
\begin{equation*}
U(s)=K_{c}\left(1+\frac{1}{\tau_{I} s}+\tau_{D} s\right) \varepsilon(s) \tag{168}
\end{equation*}
$$

where $K_{c}, \tau_{I}$, and $\tau_{D}$ are parameters to be determined. There exist different forms of the PID control law depending on the industry, era, and implementation. It is common to drop the $D$ term, that is, set $\tau_{D}=0$, and have a PI controller.

In addition to these common methods, there also exist various nonlinear approaches that can be applied to improve overall control, such as deadbanding or squared-error control.

## Section 5.1.2.1: State Feedback

A state feedback controller is an effective manner for designing a controller when it is possible to determine the behaviour of the states. A state feedback controller consists of two parts: an observer and a controller. The observer uses the output from the process to estimate the states, while the controller uses the (estimated) states to determine the input. Figure 70 shows a diagram of such a state-feedback controller and observer.


Figure 70: State-feedback control

It is possible to design different types of observers, depending on the intended application. The most common observer is the Luenberger observer which seeks to minimise the error in the predicted and actual states. Assume that the Luenberger observer can be written as

$$
\begin{align*}
& \dot{\hat{\vec{x}}}=\mathcal{A} \hat{\vec{x}}+\mathcal{L}(\vec{y}-\hat{\vec{y}})+\mathcal{B} \vec{u}  \tag{169}\\
& \hat{\vec{y}}=\mathcal{C} \hat{\vec{x}}+\mathcal{D} \vec{u}
\end{align*}
$$

where $\mathcal{L}$, the observer gain, is an appropriately sized matrix that needs to be designed and $\hat{\mathrm{o}}$ represents an estimated value. Rewriting Equation (169) gives

$$
\begin{align*}
& \dot{\overrightarrow{\vec{x}}}=(\mathcal{A}-\mathcal{L C}) \hat{\vec{x}}+\mathcal{L} \vec{y}+(\mathcal{B}-\mathcal{L D}) \vec{u}  \tag{170}\\
& \hat{\vec{y}}=\hat{\mathcal{C}} \hat{\vec{x}}+\mathcal{D} \vec{u}
\end{align*}
$$

Let the error be defined as

$$
\begin{equation*}
\vec{e}=\vec{x}-\hat{\vec{x}} \tag{171}
\end{equation*}
$$

Taking the derivative (or difference) of Equation (171) gives

$$
\begin{equation*}
\dot{\vec{e}}=\dot{\vec{x}}-\dot{\overrightarrow{\vec{x}}} \tag{172}
\end{equation*}
$$

Substituting Equations (30) and (170) into Equation (172) gives

$$
\begin{align*}
\dot{\vec{e}} & =\mathcal{A} \vec{x}+\mathcal{B} \vec{u}-(\mathcal{A}-\mathcal{L C}) \hat{\vec{x}}-\mathcal{L}(\mathcal{C} \vec{x}+\mathcal{D} \vec{u})-(\mathcal{B}-\mathcal{L D}) \vec{u}  \tag{173}\\
& =(\mathcal{A}-\mathcal{L C}) \vec{e}
\end{align*}
$$

From Equation (173), it can be seen that for the error to reach zero $\mathcal{A}-\mathcal{L C}$ must be stable, that is, the eigenvalues of $\mathcal{A}-\mathcal{L C}$ must lie to make the process stable. ${ }^{18}$

Now that we can estimate the state values, it makes sense to design a state controller. Assume that the state controller law can be written as

$$
\begin{equation*}
\vec{u}=-\mathcal{K} \hat{\vec{x}} \tag{174}
\end{equation*}
$$

Before examining the general case, let us consider the situation, where perfect information can be obtained about the states and there is no need to design an observer, that is, $\mathcal{C}=\mathcal{J}$, which implies that $\vec{y}=\vec{x}$. In this situation, we can use the basic state-space equations, given by Equation (30), to describe the system. Substituting Equation (174) into Equation (30) and re-arranging gives

[^14]\[

$$
\begin{align*}
\dot{\vec{x}} & =\mathcal{A} \vec{x}-\mathcal{B}(\mathcal{K} \vec{x})=(\mathcal{A}-\mathcal{B K}) \vec{x}  \tag{175}\\
\vec{y} & =\mathcal{C} \vec{x}+\mathcal{D} \vec{u}
\end{align*}
$$
\]

Thus, it can be seen that the stability of the controlled (overall, closed-loop) system is determined by the location of the eigenvalues of the $\mathcal{A}-\mathcal{K C}$ matrix.

Now consider the case, where we do not have full information about the process. In this case, we will need to consider two different equations: the observer equation and the true statespace model. Substituting the control law given by Equation (174) into the observer equation, Equation (170), gives

$$
\begin{align*}
& \dot{\overrightarrow{\hat{x}}}=(\mathcal{A}-\mathcal{L C}-\mathcal{B K}+\mathcal{L D K}) \hat{\vec{x}}+\mathcal{L} \vec{y}  \tag{176}\\
& \hat{\vec{y}}=\mathcal{C} \hat{\vec{x}}-\mathcal{D K} \hat{\vec{x}}
\end{align*}
$$

The fact that both terms are present in this equation raises the question if the controller and observer must be designed simultaneously or could they be designed separately.

In fact, the linear, state-space law presented here coupled with the linear observer are fully separable and stability will be assured if both components are stable. In order to show this, rewrite the controller law as

$$
\begin{equation*}
\vec{u}=-\mathcal{K}(\vec{x}-\vec{e}) \tag{177}
\end{equation*}
$$

From Equation (175), we can write the true state-space system as

$$
\begin{equation*}
\dot{\vec{x}}=\mathcal{A} \vec{x}-\mathcal{B K}(\vec{x}-\vec{e})=(\mathcal{A}-\mathcal{B K}) \vec{x}+\mathcal{B K} \vec{e} \tag{178}
\end{equation*}
$$

Now, combining this equation with the error equation, given as Equation (173), into a single equation, we can see that the combined system consisting of the states and errors can be written as

$$
\left[\begin{array}{l}
\dot{\vec{x}}  \tag{179}\\
\dot{\vec{e}}
\end{array}\right]=\left[\begin{array}{cc}
\mathcal{A}-\mathcal{B K} & \mathcal{B K} \\
0 & \mathcal{A}-\mathcal{L K}
\end{array}\right]\left[\begin{array}{l}
\vec{x} \\
\vec{e}
\end{array}\right]
$$

Since the combined system matrix is triangular, the eigenvalues can be obtained by considering the eigenvalues of the diagonal elements, that is, the eigenvalues of the $\mathcal{A}-\mathcal{B K}$ and $\mathcal{A}-\mathcal{L K}$ matrices. Furthermore, since neither of the two matrices interact with each other, it can be concluded that both can be designed separately and then combined. As long as each is stable, then the overall system will also be. This result is called the separation principle and is a very powerful result in designing state-feedback systems.

## Section 5.1.2.2: Proportional, Integral, and Derivative (PID) Control

Another class of common controllers is the proportional, integral, and derivative (PID) controllers. This class of controllers provides effective control for many industrial processes and can be configured to deal with a wide range of different situations. The general form of the PID controller can be written as

$$
\begin{equation*}
G_{c}=K_{c}\left(1+\frac{1}{\tau_{I} S}+\tau_{D} S\right) \tag{180}
\end{equation*}
$$

where $K_{c}$ is the proportional term or controller (proportional) gain, $\tau_{I}$ is the integral term or the integral time constant in units of time, and $\tau_{D}$ is the derivative term or the derivative time constant in units of time. Another common representation is given as

$$
\begin{equation*}
G_{c}=K_{c}+K_{I} \frac{1}{s}+K_{D} s \tag{181}
\end{equation*}
$$

where $K_{I}$ is the integral gain and $K_{D}$ is the derivative gain. In many industrial plants, it is possible to find the terms reset for integral and rate for derivative, for example, $\tau_{I}$ would be called the reset time constant.

In industry, different combinations of this controller can be found including the rather common proportional and integral controller (PI), where $\tau_{D}=K_{D}=0$, that is, the $D$-term is ignored. In order to understand the behaviour of this controller, each of the three terms $(P, I$, and $D)$ will be investigated separately. As well, the two common configurations (PI and PID) will be considered in greater detail.

## Section 5.1.2.2.a: Proportional Term

The proportional term, $K_{c}$, has the greatest impact on the overall control performance. It represents the contribution of the current error to the overall controller action, and thus, controls two key aspects: stability and speed of response. The larger the absolute value of $K_{c}$ is, the faster the system will response to changes, but the less stable the system will be. Furthermore, a controller containing only the proportional term will have bias or offset, that is,

$$
\begin{equation*}
\lim _{s \rightarrow 0} \frac{s G_{c} G_{p}}{1+G_{c} G_{p}} R_{s}=\lim _{s \rightarrow 0} \frac{s K_{c} G_{p}}{1+K_{c} G_{p}} \frac{M_{r}}{s}=\frac{K_{c} K_{p} M_{r}}{1+K_{c} K_{p}} \tag{182}
\end{equation*}
$$

where $R_{s}$ is the setpoint which gives a step change of $M_{r}$ and $K_{p}$ is the process gain. It is assumed that the overall closed-loop transfer function is stable. From Equation (182), it can be seen that the
steady-state value for the closed-loop transfer function is not 1 . This implies that the final value will not be equal to the setpoint. For this reason, pure P-controllers are rarely used.

## Example 25: Investigation of the Proportional Term on Stability and Performance

Consider the following first-order system

$$
\begin{equation*}
G_{p}=\frac{1.5}{20 s+1} e^{-10 s} \tag{183}
\end{equation*}
$$

controlled with a proportional controller $G_{c}=K_{c}$. Setting $K_{c}$ equal to $-0.25,0,0.5,1.0$, and 2.0, show how the closed-loop response to a step change of 2 in the setpoint changes.

## Solution:

Figure 71 shows the results for changing the $K_{c}$. It can be seen that for none of the cases is the setpoint attained. There is offset for all of the cases. Furthermore, for $K_{c}=0$, the process remains at zero, since no control is being implemented. As $K_{c}$ increases, the amount of oscillation and length of time required to attain a steady-state value increase. In fact, if $K_{c}$ is increased further, the process will become unstable. Similarly, for $K_{c}<0$, the process is initially stable. However, increasing the value will rapidly make the system unstable.


Figure 71: Effect of Changing $K_{c}$ for a P-Controller

## Section 5.1.2.2.b: Integral Term

The integral term, $\tau_{l}$, provides a measure of additional stability to the system as well as removing any bias present. It represents the contribution of past errors on the current controller action. In general, the integral term is always combined with a proportional term to give the proportional and integral controller. Furthermore, a controller containing only the integral term will have no bias or offset, that is,

$$
\begin{equation*}
\lim _{s \rightarrow 0} \frac{s G_{c} G_{p}}{1+G_{c} G_{p}} R_{s}=\lim _{s \rightarrow 0} \frac{s \frac{1}{\tau_{I} s} G_{p}}{1+\frac{1}{\tau_{I} s} G_{p}} \frac{M_{r}}{s}=\lim _{s \rightarrow 0} \frac{M_{r} G_{p}}{\tau_{I} s+G_{p}}=M_{r} \tag{184}
\end{equation*}
$$

where $R_{s}$ is the setpoint which makes a step change of $M_{r}$. It is assumed that the overall closedloop transfer function is stable. From Equation (184), it can be seen that the closed-loop gain is exactly equal to the step change in the setpoint, which implies that there will be no bias or offset.

Since the integral term represents the effect of past errors on the current controller action, it is possible that in a real implementation, integral (reset) wind-up can occur. This occurs when the controller output is above the physical limit of the actuator. Since the controller does not have a way of knowing that the physical limit has been reached, it continues to increase the desired controller output as the error keeps on increasing. This makes the integral term very large. When there is a need to now decrease this value, the controller decreases the output value from its large, unphysical value until it goes below the upper limit. The time delay between when the actuator should have started to respond and when it is observed to be responding is called integral wind-up. This is shown in Figure 72, where it can be seen that at 100 min , when the setpoint is changed, it takes the process another 20 minutes before it even starts changing, since the true actuator value remains saturated at its upper limit. Integral wind-up can be avoided by including the physical limits of the actuator in the controller, that is, by making values beyond the physical limits of the actuator equal to the physical limit.


Figure 72: Integral Wind-Up
Example 26: Investigation of the Integral Term on Stability and Performance
Consider the following first-order system

$$
\begin{equation*}
G_{p}=\frac{1.5}{20 s+1} e^{-10 s} \tag{185}
\end{equation*}
$$

controlled with a proportional controller $G_{c}=1 / \tau_{I S}$. Varying the value of $\tau_{I}$ from 20 to 100 in increments of 10 show how the closed-loop response to a step change of 2 in the setpoint changes.

## Solution:

Figure 73 show the results. It can be seen that as $\tau_{I}$ increases the impact on the system decreases. Note that the system is originally stable so that without any control it will reach the desired setpoint. Furthermore, and unlike with only a $P$-term, there is no bias in the closed-loop system.


Figure 73: Effect of Changing $\tau_{I}$ for a I-Controller. The solid, black line is the setpoint.

## Section 5.1.2.2.c: Derivative Term

The derivative term, $\tau_{D}$, provides a measure of additional stability to the system. The derivative term represents the contribution of future (estimated) errors on the controller action. In general, the derivative term is almost always combined with a proportional and integral terms. Furthermore, a controller containing only the derivative term will be unable to track the setpoint, that is,

$$
\begin{equation*}
\lim _{s \rightarrow 0} \frac{s G_{c} G_{p}}{1+G_{c} G_{p}} R_{s}=\lim _{s \rightarrow 0} \frac{s \tau_{D} s G_{p}}{1+\tau_{D} s G_{p}} \frac{M_{r}}{s}=\lim _{s \rightarrow 0} \frac{s M_{r} G_{p}}{1+\tau_{D} s G_{p}}=0 \tag{186}
\end{equation*}
$$

where $R_{s}$ is the setpoint which makes a step change of $M_{r}$. It is assumed that the overall closedloop transfer function is stable. From Equation (186), it can be seen that irrespective of the change in the setpoint, the system will return to zero, which implies that the derivative term cannot track the setpoint. Therefore, the main role of the derivative term is for disturbance rejection, since in that case, it is desired that the system return to the same initial value.

The derivative term as written cannot be physically realised, since the numerator is greater than the denominator. In order to avoid this problem, it is common to rewrite the derivative term into the following form:

$$
\begin{equation*}
G_{c, \text { real }}=\tau_{D} \frac{N}{\left(1+\frac{N}{s}\right)} \tag{187}
\end{equation*}
$$

where $N$ is a filter parameter to be selected. It can be noted that as $N \rightarrow \infty$, this form will become equal to the original version.

Using a derivative term introduces two potential problems into the control strategy: jitter and derivative kick. Jitter occurs when error signal fluctuates greatly about some mean value causing the estimated value of the future error to also fluctuate greatly. This in turn causes unnecessary fluctuations in the controller output, which can lead to issues with the actuator. Figure 74 shows the typical situation where jitter is present. Jitter can be mitigated by filtering the error signal used in computing the derivative term with a low-pass filter, which removes the high frequency component of the noise. Derivative kick occurs when the setpoint has a sudden change in value (for example, a step change) causing the instantaneous error to be practically speaking infinite. This causes the controller error to spike for a very short period of time after the change occurs. Figure 75 shows the typical situation when derivative kick occurs. When the setpoint changes at 10 s , there is a corresponding spike in the actuator value. Note that in practice, due to saturation in the actuator, the impact of the derivative kick will be less than shown. Nevertheless, it will be present. Derivative kick can be lessened by solely using the change in the output rather than the error when computing the derivative term, that is,

$$
\begin{equation*}
u_{t}=\tau_{D} s y_{t} \tag{188}
\end{equation*}
$$

Since it has been shown that the overall impact of the derivative term is primarily for disturbance rejection, that is, making $y_{t} \approx 0$, this approach is reasonable.


Figure 74: Jitter with a Derivative Term


Figure 75: Derivative Kick

## Example 27: Investigation of the Derivative Term on Stability and Performance

Consider the following first-order system

$$
\begin{equation*}
G_{p}=\frac{1.5}{20 s+1} e^{-10 s} \tag{189}
\end{equation*}
$$

controlled with a derivative controller $G_{c}=\tau_{D} S$. Varying the value of $\tau_{D}$ from 1 to 11 in increments of 2 show the closed-loop response of the controller to a white-noise disturbance of magnitude 0.05 .

## Solution:

Figure 76 show the results. It can be seen that as the magnitude of the $D$-Term increases, so does the magnitude of the response. This implies that the $D$-Term only influences the magnitude of the response.


Figure 76: Effect of Changing $\tau_{D}$ for a D-Controller

## Section 5.1.2.2.d: Proportional and Integral (PI) Controller

A proportional and integral (PI) controller is one of the most common industrial controllers found. It combines the proportional and integral terms to give

$$
\begin{equation*}
G_{c}=K_{c}\left(1+\frac{1}{\tau_{I} s}\right) \tag{190}
\end{equation*}
$$

The behaviour of a PI controller will be similar to that of each of the two separate terms, that is, the proportional term will influence the stability and overall performance, while the integral term
will provide a bias-free controller. This controller can be interpreted as only considering the current and past information about the process and how that impacts the overall control performance.

PI control can be used for both disturbance rejection and setpoint tracking in many different applications including flows, levels, and temperatures. Large time delays and highly nonlinear systems can limit its effectiveness.

## Section 5.1.2.2.e: Proportional, Integral, and Derivative (PID) Controller

A proportional, integral, and derivative (PID) controller can be tuned on the basis of the three previously considered terms. Such a controller takes all available information about the process and predicts the future trajectory of the system. It is commonly used in large complex systems where there is a need for an additional degree-of-freedom to stabilise the system. Systems which have a very jittery output or setpoint would need to be filtered before being used as the unfiltered signal can cause undesired behaviour.

## Section 5.1.2.2.f: Discretisation of the PID Controller

Since many control algorithms are implemented on a computer, it can often be helpful to implement the PID controller in the discrete domain. To discretise the PID controller, let $s=1-z^{-1}$ in Equation (180) to give

$$
\begin{equation*}
u_{k}=K_{c}\left(1+\frac{1}{\tau_{I}\left(1-z^{-1}\right)}+\tau_{D}\left(1-z^{-1}\right)\right) \varepsilon_{k} \tag{191}
\end{equation*}
$$

Re-arranging Equation (191) gives

$$
\begin{equation*}
u_{k}=\frac{K_{c}}{\left(1-z^{-1}\right)}\left(\left(1-z^{-1}\right)+\frac{1}{\tau_{I}}+\tau_{D}\left(1-z^{-1}\right)^{2}\right) \varepsilon_{k} \tag{192}
\end{equation*}
$$

Equation (192) is often called the positional form of the discretised PID controller. Re-arranging Equation (192) to give

$$
\begin{equation*}
\left(1-z^{-1}\right) u_{k}=\Delta u_{k}=K_{c}\left(\left(1-z^{-1}\right)+\frac{1}{\tau_{I}}+\tau_{D}\left(1-z^{-1}\right)^{2}\right) \varepsilon_{k} \tag{193}
\end{equation*}
$$

produces another common PID equation called the velocity form of the discretised PID controller. In industrial implementations, it is possible to encounter both forms. Although these forms do not have an impact on controller design, they will have an impact on understanding the values obtained.

## Section 5.1.2.3: Controller Tuning

Having established the different types of controllers that can be used and understood some of behaviour of the parameters on the overall control system, the remaining question is how do we correlate a given set of performance criteria (for example, the objectives of control previously mentioned) and the different controller parameters. One easy approach would be to use some simulation software to obtain the desired parameter values using a trial-and-error method. As one can easily see, this approach could take very long and not necessarily provide the optimal parameter values. Instead, various correlations or approaches have been developed that will allow for the performance of a controller to be specified. In this section, we will consider the methods for tuning state-space controllers and PID controllers. Irrespective of the approach used, there will always be a need to simulate the resulting control system to make sure that there are no issues with the performance. Furthermore, it should be noted that for certain applications, there exist, specific controller tuning rules that provide optimal performance for the problem at hand, for example, level controllers are often tuned using special rules that seek to minimise large deviations from the setpoint.

The general approach to controller tuning can be summarised by Figure 77. Before starting the tuning procedure, it is required that the process be understood or modelled, that the performance criteria be clearly specified, and that the control strategies of interest be known. The first step will be to determine the initial controller parameters, followed by a simulation (or actual implementation, if the system can handle repeated perturbations). Based on the test and a comparison with the desired performance, new controller parameters can be obtained and new tests performed. This would be repeated until either the desired criteria are satisfied or the time available has run out. At this point, it may happen that the process understanding is found to be deficient and additional information would need to be sought. If the controller has not been implemented on the real system during the initial controller testing, then it will be necessary to repeat this procedure once more on the actual system. Very often, it will be found that the performance may not be as expected given the differences between the simulated and actual processes. However, following this procedure will minimise the risk of the occurrence of such a situation.


Figure 77: Controller Tuning Workflow

## Section 5.1.2.3.a: Tuning a State-Space Controller

A state-space controller is often tuned using pole placement, which involves selecting the desired poles of the closed-loop transfer function and designing a controller that can achieve this.

When placing the poles or eigenvalues of the closed-loop transfer function using a statespace approach, both the controller and observer will need to have their poles placed. The general rule of thumb is that the poles of the observer should be 10 to 20 times faster than the poles of the controller. Pole placement can be achieved using two different methods: characteristic polynomial and Ackermann's Formula. The approaches work for both continuous and discrete controllers. It should be noted that for multi-input systems, the computed value of $\mathcal{K}$ is not necessarily unique.

The general problem statement is: given the desired eigenvalues $\left\{\lambda_{1}, \lambda_{2}, \ldots, \lambda_{n}\right\}$ and the system $\{\mathcal{A}, \mathcal{B}, \mathcal{C}\}$, determine the state feedback controller gain $\mathcal{K}$. For the characteristic polynomial approach, perform the following steps:

1) Compute $\bar{\Delta}(s)=\prod_{i=1}^{n}\left(s-\lambda_{i}\right)=s^{n}+\alpha_{1}^{\prime} s^{n-1}+\cdots+\alpha_{n-1}^{\prime} s+\alpha_{n}^{\prime}$.
2) Compute $\Delta(s)=s^{n}+\alpha_{1} s^{n-1}+\cdots+\alpha_{n-1} s+\alpha_{n}$, which is the characteristic polynomial of $\mathcal{A}$.
3) Compute $\overline{\mathcal{A}}=\left[\begin{array}{cccc}-\alpha_{1}^{\prime} & \cdots & -\alpha_{n-1}^{\prime} & -\alpha_{n}^{\prime} \\ 1 & & 0 & 0 \\ 0 & \ddots & 0 & 0 \\ 0 & 0 & 1 & 0\end{array}\right]$, $\overline{\mathcal{B}}=\left[\begin{array}{l}1 \\ 0 \\ 0 \\ 0\end{array}\right]$.
4) Compute $\overline{\mathcal{C}}=\left[\begin{array}{llll}\mathcal{B} & \mathcal{A B} & \cdots & \mathcal{A}^{n-1} \mathcal{B}\end{array}\right]$ and $\overline{\mathcal{C}^{\prime}}=\left[\begin{array}{llll}\overline{\mathcal{B}} & \overline{\mathcal{A}} \overline{\mathcal{B}} & \cdots & \overline{\mathcal{A}}^{n-1} \overline{\mathcal{B}}\end{array}\right]$.
5) $\mathcal{K}=\left[\begin{array}{lll}\alpha_{1}^{\prime}-\alpha_{1} & \cdots & \alpha_{n}^{\prime}-\alpha_{n}\end{array}\right] \overline{\mathcal{C}}^{\prime} \overline{\mathcal{C}}^{-1}$.

Similarly, Ackermann's formula, uses the following steps to place the poles:

1) Compute $\bar{\Delta}(s)=\prod_{i=1}^{n}\left(s-\lambda_{i}\right)=s^{n}+\alpha_{1}^{\prime} s^{n-1}+\cdots+\alpha_{n-1}^{\prime} s+\alpha_{n}^{\prime}$.
2) Compute $\Delta(s)=s^{n}+\alpha_{1} s^{n-1}+\cdots+\alpha_{n-1} s+\alpha_{n}$, which is the characteristic polynomial of $\mathcal{A}$.
3) Compute $\overline{\mathcal{A}}=\left[\begin{array}{cccc}-\alpha_{1}^{\prime} & \cdots & -\alpha_{n-1}^{\prime} & -\alpha_{n}^{\prime} \\ 1 & & 0 & 0 \\ 0 & \ddots & 0 & 0 \\ 0 & 0 & 1 & 0\end{array}\right]$, $\overline{\mathcal{B}}=\left[\begin{array}{l}1 \\ 0 \\ 0 \\ 0\end{array}\right]$.
4) Compute $\overline{\mathcal{C}}=\left[\begin{array}{llll}\mathcal{B} & \mathcal{A B} & \cdots & \mathcal{A}^{n-1} \mathcal{B}\end{array}\right]$ and $\overline{\mathcal{C}^{\prime}}=\left[\begin{array}{llll}\overline{\mathcal{B}} & \overline{\mathcal{A}} \overline{\mathcal{B}} & \cdots & \overline{\mathcal{A}}^{n-1} \overline{\mathcal{B}}\end{array}\right]$.
5) $\mathcal{K}=\left[\begin{array}{llll}0 & \cdots & 0 & 1\end{array}\right]_{1 \times n} \overline{\mathcal{C}}^{-1} \bar{\Delta}(\mathcal{A})$.

It is also possible to design the observer using Ackermann's formula which gives

$$
\mathcal{L}=\bar{\Delta}(\mathcal{A}) \mathcal{O}^{-1}\left[\begin{array}{llll}
0 & \cdots & 0 & 1 \tag{194}
\end{array}\right]_{1 \times n}^{T}
$$

When using pole placement, the following points should be borne in mind:

1) The magnitude of $\mathcal{K}$ gives the amount of effort required to control the process. The further the desired poles are from the actual poles of the system, the larger the controller gain $\mathcal{K}$.
2) For multi-input systems, $\mathcal{K}$ is not unique.
3) $(\mathcal{A}, \mathcal{B})$ and $(\mathcal{A}-\mathcal{B} \mathcal{K}, \mathcal{B})$ are controllable. However, due to pole-zero cancellations, the resulting system may not be observable.
4) Discrete systems can be controlled in the same manner, mutatis mutandi.

## Section 5.1.2.3.b: Tuning a PID Controller

When designing a PID controller, there exist various different approaches that can be taken to specifying the initial controller parameters. In practice, once these initial parameters have been
obtained, they will be fine-tuned to obtain the desired performance. There are two main approaches to controller tuning: model-based and structure-based.

In model-based controller tuning, a model of the process is required and the closed-loop behaviour of the resulting system is specified. Using this information, it is possible to obtain the resulting form of the controller transfer function. Comparing the resulting controller transfer function with the standard PID controller allows equations for the constants to be determined. The most common model-based approach is the internal model control (IMC) framework. The main advantage of this approach is that it allows the engineer to specify the desired closed-loop behaviour of the system, while the main disadvantage of this approach is that a relatively accurate model of the system is required. Note that the specification of the closed-loop transfer function can be used to add robustness to the system by changing the speed (time constant) of the response. The faster the system responds to changes the less robust is the resulting closed-loop system. In many cases, the simplified IMC (SIMC) rules provide better control performance for industrial systems. The SIMC rules take into consideration additional factors, such as disturbances in the input and faster response for systems with large time constants. SIMC rules recommend PI controllers for first-order systems and PID controllers for second-order systems.

In structure-based controller tuning, the structure (or type) of a controller, the objective (disturbance rejection or setpoint tracking), and metric for measuring good control are specified. Minimising this metric allows the values of the parameters to be determined. The most common structure-based approach is using the integrated time-averaged error (ITAE) as the metric. Using the ITAE metric penalises persistent errors, which can be small errors that last for a long period of time, and provides a conservative approach to control. The main advantage of this method is that a model of the system need not be provided, while the main disadvantages are that the engineer has no control over the closed-loop response of the system and that the resulting system is not very robust to changes in the process.

Table 26 and Table 27 present some common tuning methods for first-order transfer functions. Table 28 presents the SIMC tuning method for a second-order transfer function. It should be noted that the PID controller has the following series form:

$$
\begin{equation*}
G_{c}=\tilde{K}_{c}\left(1+\frac{1}{\tilde{\tau}_{I} s}\right)\left(1+\tilde{\tau}_{D} s\right) \tag{195}
\end{equation*}
$$

where the controller parameters with the tilde (\%) have the same meaning as the corresponding controller parameters without a tilde.

Table 26: PI controller constants for first-order-plus-deadtime models

| Controller Method |  | $K_{c}$ | $\tau_{I}$ |
| :---: | :---: | :---: | :---: |
| 己0000000000 | ITAE <br> (setpoint tracking) | $\frac{0.586}{K}\left(\frac{\theta}{\tau}\right)^{-0.916}$ | $\frac{\tau}{1.03-0.165\left(\frac{\theta}{\tau}\right)}$ |
|  | ITAE <br> (disturbance rejection) | $\frac{0.859}{K}\left(\frac{\theta}{\tau}\right)^{-0.977}$ | $\frac{\tau}{0.674}\left(\frac{\theta}{\tau}\right)^{0.680}$ |
|  | SIMC <br> ( $\tau_{c}$ chosen so that $\tau_{c} / \theta>0.8$ and $\tau_{c}>\tau / 10$ ) | $\frac{1}{K} \frac{\tau}{\theta+\tau_{c}}$ | $\min \left(\tau, 4\left(\tau_{c}+\theta\right)\right)$ |

Table 27: PID controller constants for first-order-plus-deadtime models

| Controller Method |  | $K_{c}$ | $\tau_{I}$ | $\tau_{D}$ |
| :---: | :---: | :---: | :---: | :---: |
|  | $\begin{gathered} \text { ITAE } \\ \text { (setpoint tracking) } \end{gathered}$ | $\frac{0.965}{K}\left(\frac{\theta}{\tau}\right)^{-0.85}$ | $0.796-0.147\left(\frac{\theta}{\tau}\right)$ | $0.308 \tau\left(\frac{\theta}{\tau}\right)^{0.929}$ |
|  | ITAE <br> (disturbance rejection) | $\frac{1.357}{K}\left(\frac{\theta}{\tau}\right)^{-0.947}$ | $\frac{\tau}{0.842}\left(\frac{\theta}{\tau}\right)^{0.738}$ | $0.381 \tau\left(\frac{\theta}{\tau}\right)^{0.995}$ |
| $\begin{aligned} & \ddot{0} \\ & 0 \\ & 0 \\ & 0 \\ & \frac{1}{0} \\ & 0 \\ & \vdots \\ & \hline \end{aligned}$ | $\begin{gathered} \hline \hline \text { IMC } \\ \left(\tau_{c}\right. \text { chosen so that } \\ \tau_{c} / \theta>0.8 \text { and } \\ \left.\tau_{c}>\tau / 10\right) \end{gathered}$ | $\frac{1}{K}\left(\frac{2 \frac{\tau}{\theta}+1}{2 \frac{\tau_{C}}{\theta}+1}\right)$ | $\frac{\theta}{2}+\tau$ | $\frac{\tau}{2\left(\frac{\tau}{\theta}\right)+1}$ |

Table 28: PID controller constants for a second-order-plus-deadtime models

|  | Controller Method | $\tilde{K}_{c}$ | $\tilde{\tau}_{I}$ | $\tau_{D}$ |
| :---: | :---: | :---: | :---: | :---: |
|  | $\begin{aligned} & \text { SIMC } \\ & G_{p}=\frac{K}{\left(\tau_{1} s+1\right)\left(\tau_{2} s+1\right)} e^{-\theta s} \\ & \text { with } \tau_{1}>\tau_{2} \end{aligned}$ | $\frac{1}{K}\left(\frac{\tau_{1}}{\theta+\tau_{c}}\right)$ | $\min \left(\tau_{1}, 4\left(\theta+\tau_{c}\right)\right)$ | $\tau_{2}$ |

## Example 28: Designing a PI Controller

Consider the following first-order system

$$
\begin{equation*}
G_{p}=\frac{1.5}{20 s+1} e^{-10 s} \tag{196}
\end{equation*}
$$

Determine the controller parameters for a PI controller using the IMC method.

## Solution

Before solving the problem, it is helpful to determine the values of the various constants

$$
\text { Gain: } K=1.5
$$

Time Constant: $\tau=20$
Time Delay: $\theta=10$
For the IMC method, there is a need to also specify the value of $\tau_{c}$, which is essentially the closedloop time constant. The smaller the value, the faster the response is, but less robust is the overall system. The constraints given with this method imply that $\tau_{c}>0.8 \theta=8$ and $\tau_{c}>0.1 \tau=2$. For the purposes of this example, set the value of $\tau_{c}$ to be 10 , which satisfies both constraints.

Therefore, the controller parameters can be computed from Table 26 as follows:

$$
\begin{align*}
K_{c}=\frac{1}{K} \frac{\tau}{\theta+\tau_{c}} & =\frac{1}{1.5}\left(\frac{20}{10+10}\right)=\frac{2}{3}  \tag{197}\\
\tau_{I} & =\tau=20 \tag{198}
\end{align*}
$$

The resulting behaviour of the closed-loop system is shown in Figure 78. It can be seen that the process responds almost exactly as expected.


Figure 78: Closed-Loop Performance of the PI Controller

## Section 5.1.2.4: Controller Performance

Once a controller has been designed, it is often necessary to measure its performance and determine if the controller is attaining its objectives. As with any such measurements, performance is measured against some (engineering-related) specification. Unlike most other equipment, the specification for a controller involves time, either explicitly when speed of response is a criterion, or implicitly, for statistics such as standard deviation. The performance of a controller can be measured based on two different sets of criteria: how the controller responds to a change in setpoint, that is, the servo response, and how the controller responds to disturbances, that is, the regulatory response.

Since the objective in servo response performance is how well the controller responds to a setpoint change, it is easy to determine an appropriate measure for performance assessment. For this reason, servo response is often used to design a controller. As a first approximation, a closedloop system can be considered to be a second-order system with time delay. Therefore, the response of the closed loop to a step change in setpoint would be the step response of a secondorder system. The most common servo control performance measures are just characteristics of a second-order system responding to a step input, that is, rise time, overshoot, and settling time of the controlled variable, as shown in Figure 79. The rise time, $\tau_{r}$, is defined as the first time the process reaches the setpoint value. The overshoot is defined as the ratio of how much the process goes over (or under) the setpoint divided by the magnitude of the step change. Note that if the step change is negative, then the overshoot will also be negative (that is, it is really an undershoot) so that the ratio remains positive. The settling time, $\tau_{s}$, is defined as the last time for which the process lies outside the $5 \%$ bounds. The $5 \%$ bounds are defined as a boundary on either side of the new setpoint that is defined as $2.5 \%$ of the difference $\left(y_{s s, 2}-y_{s s, 1}\right)$, where $y_{s s}$ is the steady-state value and the numeric subscripts represent the initial and final values. The settling time is roughly three times the closed-loop time constant.


Figure 79: Measures of Servo Control Performance
On the other hand, for performance assessment of regulatory response, the objective is less certain. Instead, various quantitative measures are used to quantify performance assessment, including:

1) Standard deviation of manipulated variable: This is the simplest and most important measure of control performance. However, it is difficult to specify the desired value $a$ priori since it depends on the nature and magnitude of disturbances affecting the process. What can be specified, or at least described qualitatively, is the trade-off between control error and control action. One way of thinking of a controller is that it moves variation from one variable (usually the controlled variable) to another (usually the manipulated variable). As disturbances affect the process, the controller will respond by changing the
manipulated variable. The response will then mitigate the effect of the disturbances. Larger disturbances will require greater control action than small disturbances.
2) Mean of the output variable: This can be used to assess how well the controller maintains the given setpoint in the presence of disturbances.
3) Advanced performance assessment: More advanced performance assessment methods are available, such as the minimum variance controller and the Harris Index, which can provide appropriate benchmarks for determining good performance.

## Section 5.2: Feedforward Control

In feedforward control, a measurable disturbance is used as the input to a controller, so as to take corrective action before the disturbance impacts the process. The goal when using feedforward control is to design a controller that can correct the disturbance at the same time as the disturbance affects the process. In practice, such a controller can rarely be designed, since there will be time delay in the system, that is, there is some time period between when the variable is measured and it impacts the system. If the time delay between the disturbance and the process is smaller than the time delay between the input and the process, then it is not possible to compensate the disturbance before it has impacted the system. Figure 80 shows a typical feedforward control loop.


Figure 80: Block Diagram for Feedforward Control
A special type of feedforward controller is a decoupler, which seeks to separate two or more interacting processes so that they can be treated as independent systems controlled by their
own SISO control loops. A decoupler basically takes the controller output from the other controllers and treats them as a measurable disturbance that needs to be countracted. Decouples can be very effective in controlling a multivariate system when the interactions are relatively straight forward and not too nonlinear.

In feedforward control design, the objective to design a controller that can compensate for a measured disturbance before it can affect the process. Assume that the process model is $G_{p}$ and the disturbance model is $G_{d}$, then for the feedforward controller shown in Figure 80, the basic form of the feedforward controller can be written as

$$
\begin{equation*}
G_{f f}=-\frac{G_{d}}{G_{p}} \tag{199}
\end{equation*}
$$

In the ideal situation, this will exactly cancel out the effect of the measured disturbance. However, in practice, such a controller cannot be implemented for the following two reasons:

1) Time Delay: If the time delay in the process response is larger than the time delay in the disturbance response, then the overall time delay for the process will be negative requiring knowledge of future information about the process. Since such a situation is impossible, the general solution is to drop the unrealisable time delay from the final controller.
2) Unstable Zeroes in $\boldsymbol{G}_{\boldsymbol{p}}$ : If the process contains unstable zeroes, then the resulting transfer function will also be unstable. In such a case, it will not be possible to realise the controller. A common solution in this case is to drop the offending zeroes from the final representation.

There are two common feedforward controllers that can be designed:

1) Static controllers where only the gains of the process are considered, that is,

$$
\begin{equation*}
G_{f f}=-\frac{K_{d}}{K_{p}} \tag{200}
\end{equation*}
$$

where $K_{d}$ is the disturbance gain and $K_{p}$ is the process gain.
2) Dynamic controllers where the process and disturbance dynamics are included in the final control law. Very often, the controller takes the form of a lead-lag controller, that is,

$$
\begin{equation*}
G_{f f}=-\frac{K_{d}\left(\tau_{p} s+1\right)}{K_{p}\left(\tau_{d} s+1\right)} e^{-\theta_{f f} s} \tag{201}
\end{equation*}
$$

where $\tau_{p}$ is the process time constant, $\tau_{d}$ is the disturbance time constant, and $\theta_{f f}=\theta_{d}-\theta_{p}$. If $\theta_{f f}$ is negative, then the term is often ignored. Ideally, $\left|\tau_{p} / \tau_{d}\right|<1$, which avoids large spikes when this controller is used.

## Example 29: Designing a Feedforward Controller

Given the following system, design both a dynamic and a static feedforward controller. Compare the performance of both. The system parameters are:

$$
\begin{align*}
& G_{p}=\frac{-1.5(s-1)}{(s+1)(10 s+1)} e^{-10 s}  \tag{202}\\
& G_{d}=\frac{-0.5}{(5 s+1)(7 s+1)} e^{-5 s}  \tag{203}\\
& \quad G_{c}=\frac{2}{3}\left(1+\frac{1}{20 s}\right) \tag{204}
\end{align*}
$$

For the simulation, assume that the measured disturbance is driven by Gaussian white noise with a magnitude of 0.5 and there is a step change in the setpoint of 2 units after 100 s .

## Solution

Using Equation (200), the static feedforward controller can be computed as

$$
\begin{equation*}
G_{f f, s}=-\frac{K_{d}}{K_{p}}=-\frac{-0.5}{1.5}=\frac{1}{3} \tag{205}
\end{equation*}
$$

Note that the gain can be obtained by setting $s=0$ and evaluating the resulting value.
When creating the dynamic feedforward controller, it is important that the transfer functions be written in their standard form, that is, all roots are of the form $\tau s+1$. In this particular example, the zero of the process transfer function needs to be written into this form to give

$$
\begin{equation*}
G_{p}=\frac{1.5(-s+1)}{(s+1)(10 s+1)} e^{-10 s} \tag{206}
\end{equation*}
$$

The dynamic feedforward controller can be designed using Equation (199). This gives

$$
\begin{equation*}
G_{f f, d}=-\frac{G_{d}}{G_{p}}=-\frac{\frac{-0.5}{(5 s+1)(7 s+1)} e^{-5 s}}{\frac{1.5(-s+1)}{(s+1)(10 s+1)} e^{-10 s}}=\frac{1}{3} \frac{(s+1)(10 s+1)}{(5 s+1)(7 s+1)(-s+1)} e^{5 s} \tag{207}
\end{equation*}
$$

However, this controller as written cannot be realised since there is an unstable pole and the time delay requires future values. In order to obtain a realisable dynamic feedforward controller, both
the unstable pole and time delay terms will be dropped. This gives the following form for the controller

$$
\begin{equation*}
G_{f f, d}=\frac{1}{3} \frac{(s+1)(10 s+1)}{(5 s+1)(7 s+1)} \tag{208}
\end{equation*}
$$

Figure 81 shows the effects of both feedforward controllers on the process. It can be seen that without the feedforward controller the process has greater oscillations. Adding the static feedforward controller tends to decrease the impact of the oscillations. However, the behaviour is still rather jittery. The dynamic feedforward controller has smoothed out the jitter and the behaviour is cleaner.


Figure 81: Effect of Feedforward Control on a Process

## Section 5.3: Discrete-Event Control

In discrete-event control, a specific event or series of events triggers a given control sequence. This can often be seen when dealing with safety-oriented control situations, where, for
example, a pressure relief valve will only open if the pressure is above a certain value and will close once the pressure falls below the threshold.

Another type of discrete event control is interlocking, where a specified set of circumstances must hold before a given action can occur, for example, a microwave cannot be started until the microwave door has been closed, which minimises the impact of the radiation on the user. Interlocking is an important component of safety considerations in order to prevent operators from taking inappropriate action. Designing the correct interlocks can be a challenging proposition.

A special type of discrete-event control occurs when binary signals are used, for example, a sensor will report when the elevator has reached the correct floor, and then the motor of the drive will be switched off. The exact height of the elevator above the ground is not continuously measured and monitored. When dealing with binary signals, there are two possible types of controllers: logic control and sequential control.

In logic controllers, also called combinatorial controllers, one obtains the control variable as a combination, that is, a logical operation, of signals, for example, system output signals (sensor values such as door open or closed) or user inputs (e.g emergency button pressed or not pressed). Consider a simple example of a lathe. A lathe is turned on when the user presses the power button and the chuck is closed. The lathe is switched off when an end position is reached or the emergency stop button is pressed.

In sequential control, individual control operations are performed in certain steps. The advancement to the next step occurs either after a certain time has passed (time-dependent sequence control), for example, a simple traffic light, or in the presence of a specific event (process-dependent sequence control), for example, traffic lights that change their signal when a pedestrian presses the button. In contrast to the logic controllers, it is therefore not possible to clearly determine the manipulated variable even if all signal values are known, since one also has to know at which point in the process one is currently located.

## Section 5.4: Supervisory Control

Supervisory control focuses on designing controllers that can deal with complex, interacting systems incorporating physical, process, and economic constraints, as well as dealing with changing economic circumstances. Supervisory control systems will often set the setpoint for
subsidiary (or slave) controllers. They can be used to control large systems containing many different variables and even locations. A supervisory control system can range from a simple nested control loop system to complex model predictive control systems. There are two primary control systems cascade control and model predictive control (MPC), which is often called advanced process control (APC) in the petrochemical industry.

## Section 5.4.1: Cascade Control

In cascade control, at least two control loops are nested within each other. A typical cascade control strategy is shown in Figure 82, where it is desired to ensure that the flow rate controlled by the valve is as accurate as possible, so that the process experiences the tightest control possible. The primary or master control loop is the outer control loop, which sets the setpoint for the secondary or slave control loop. The general procedure for tuning a cascade control loop is to first tune the innermost loop and then work up towards the uppermost loop, treating all tuned controllers as part of the process. The closed-loop time constant of the inner loop should be smaller (faster) than the closed-loop time constant of the outer loop. A common rule of thumb is that the ratio of closed-loop time constant of the primary loop and that of the secondary loop ( $\tau_{c, p} / \tau_{c, s}$ ) should be between 4 and 10. Too small a ratio implies that the secondary loop will not have reached steady state and there will be interactions between the two loops leading to a loss of overall performance. Too large a ratio implies that the advantages of cascade control will be lost as it will slow down the overall response of the system. However, it will make it more robust to changes in the underlying process models.


Figure 82: Block Diagram for Cascade Control

## Section 5.4.2: Model Predictive Control

Model predictive control is an advanced control strategy that can take into consideration not only the deviations from the setpoint, but also various economic and physical constraints, for example, it can explicitly handle the situation where the water in the tank should not go above a given value. In order to accurately handle these constraints, this approach requires a good model and proper engineering design. Model predictive control works by optimising the control actions over a period of time, often called the control horizon, using the predicted values from the model. Any constraints are then implemented as necessary on the predicted results. The controller then implements the next control action and repeats the optimisation procedure again. This ensures that the impact of disturbances, plant-model mismatches, and other imperfections in the system do not have a too large impact on the overall performance.

Model predictive control is effective in handling complex process with multiple inputs and outputs that interact strongly. Although the original version was designed for linear systems, variants exist for nonlinear systems.

One of the key disadvantages of model predictive control is that it requires a good model that must be updated as necessary. A poor model can cause the control system to degrade and forgo the benefits of the approach.

There are many different implementations of model predictive control. The most popular implementation in the industry is the dynamic-matrix controller (DMC) approach, which is detailed here. The objective function for model predictive control is

$$
\begin{align*}
& \min _{\Delta \vec{u}}\left[(\vec{r}-\hat{\vec{y}})^{T} \mathcal{Q}(\vec{r}-\hat{\vec{y}})+(\Delta \vec{u})^{T} \mathcal{R}(\Delta \vec{u})\right] \\
& \text { subject to }  \tag{209}\\
& \hat{\vec{y}}=\vec{y}^{*}+\mathcal{A} \Delta \vec{u}
\end{align*}
$$

where $\vec{r}$ is the reference vector, $\hat{\vec{y}}$ the vector of the predicted process values with control, $Q$ the process scaling matrix, $\Delta \vec{u}$ the vector of the changes in the controller action, $\mathscr{R}$ the input scaling matrix, $\vec{y}^{*}$ the vector of the process values without control, and $\mathcal{A}$ the dynamic matrix. Furthermore, let $m$ be the control horizon, $p$ the prediction horizon, $d$ the process deadtime, and $n$ the settling time. It should be noted that $1 \leq m \leq p-d$ and $p>d$.

The solution requires the step response of the process, that is,

$$
\begin{equation*}
y_{t}=\sum_{i=1}^{\infty} a_{i} z^{-i} \Delta u_{t} \tag{210}
\end{equation*}
$$

where $a_{i}$ is the step-response coefficient. The step-response coefficient can be calculated either by polynomial division of the transfer function model or using the coefficients of the impulse response. The coefficients $a_{i}$ of the step response have the following relationship to the coefficients of the impulse response $h_{j}$

$$
\begin{equation*}
a_{i}=\sum_{j=1}^{i} h_{j} \tag{211}
\end{equation*}
$$

Furthermore, the settling time is defined as the time at which the first coefficient $a_{n}$ is within a range of between 0.975 and 1.025 times the value of $a_{\infty}$, where $a_{\infty}$ is the steady-state value. All step response coefficients after the settling time $n$ can be assumed to be equal to $a_{n+1}$. The $p \times m$ dynamic matrix $\mathcal{A}$ can written as

$$
\mathcal{A}=\left[\begin{array}{cccc}
a_{1} & 0 & \cdots & 0  \tag{212}\\
a_{2} & a_{1} & \ddots & \vdots \\
\vdots & \vdots & \ddots & 0 \\
\vdots & \vdots & & a_{1} \\
a_{p} & a_{p-1} & \cdots & a_{p-m+1}
\end{array}\right]
$$

For a univariate system with only one input variable and one output variable, the solution for Equation (209) is

$$
\begin{equation*}
\Delta \vec{u}=\vec{K}_{C} \vec{e}=\left(\mathcal{A}^{T} \mathcal{Q} \mathcal{A}+\mathcal{R}\right)^{-1} \mathcal{A}^{T} \mathcal{Q}^{T}\left(\vec{r}-\vec{y}^{*}\right) \tag{213}
\end{equation*}
$$

where

$$
\begin{equation*}
y_{t+l}^{*}=y_{t}-\sum_{i=1}^{n}\left(a_{i}-a_{l+i}\right) \Delta u_{t-i} \tag{214}
\end{equation*}
$$

Once the solution has been implemented, then the first control action $\Delta u_{1}$ will be implemented by the controller. At the next sampling time, the above optimisation procedure is repeated, a new optimal value is calculated, and the first control action is implemented. This allows the system to take unexpected process changes into consideration.

## Example 30: Design of a Model Predictive Controller

Design a model predictive controller for the following SISO system:

$$
\begin{equation*}
G_{p}=\frac{2 z^{-1}}{1-0,75 z^{-1}} \tag{215}
\end{equation*}
$$

Set up the required matrices and perform the first step of the iteration. For this, assume that $m=p$ $=3$. Let $\mathcal{Q}=\mathcal{J}_{p}$ and $\mathcal{R}=\mathcal{J}_{3}$, where $\mathcal{J}_{n}$ is the $n \times n$ identity matrix. Furthermore, a step change occurs at $t=0$, with the process having previously been in steady state.

## Solution

To determine the required model predictive controller, we must first determine the stepresponse model. This can be determined using long division, which gives the impulse response coefficients from which the step response coefficients can be easily calculated. Thus,

$$
\begin{array}{r}
\frac{2 z^{-1}+1.5 z^{-2}+1.125 z^{-3}}{\left.1-0.75 z^{-1}\right)} \begin{array}{l}
\frac{2 z^{-1}}{} \frac{-2 z^{-1}+1.5 z^{-2}}{1.5 z^{-2}} \\
\frac{-1.5 z^{-2}+1.125 z^{-3}}{1.125 z^{-3}} \ldots
\end{array} \tag{216}
\end{array}
$$

It can be seen that $h_{i}=2(0.75)^{i-1}$ for $i \geq 1$. In general, for a transfer function of the form

$$
\begin{equation*}
G_{p}=\frac{\beta z^{-d}}{1-\alpha z^{-1}} \tag{217}
\end{equation*}
$$

the coefficients of the impulse response can be given as

$$
h_{i}=\left\{\begin{array}{cc}
\beta \alpha^{i-d} & i \geq d  \tag{218}\\
0 & i<d
\end{array}\right.
$$

The required step response coefficients can be found using Equation (211)

$$
\begin{equation*}
a_{i}=\sum_{j=0}^{i} h_{j}=\sum_{j=1}^{i} 2\left(0.75^{j-1}\right) \tag{219}
\end{equation*}
$$

Equation (219) gives a geometric series that can be written as

$$
\begin{equation*}
a_{i}=\beta\left(\frac{1-\alpha^{i-d+1}}{1-\alpha}\right)=2 \frac{\left(1-0.75^{i}\right)}{1-0.75} \tag{220}
\end{equation*}
$$

It follows that

$$
\begin{align*}
& a_{1}=2 \frac{(1-0.75)}{1-0.75}=2 \\
& a_{2}=2 \frac{\left(1-0.75^{2}\right)}{1-0.75}=3.5  \tag{221}\\
& a_{3}=2 \frac{\left(1-0.75^{3}\right)}{1-0.75}=4.625
\end{align*}
$$

Incidentally, since $d=1$ for this example, the step response coefficients with a smaller $d$ have a value of zero and can thus be ignored in the summation.

In the next step, the settling time is calculated. For a converging geometric series (implies that the process of interest is stable), the value to which the series converges is given by

$$
\begin{equation*}
K_{p}=a_{\infty}=\frac{\beta}{1-\alpha}=\frac{2}{1-0.75}=8 \tag{222}
\end{equation*}
$$

This gives the settling time for the first value for which the process lies in the interval $0.975 \times 8$ and $1.025 \times 8$. Since the process under consideration has no oscillations, only the lower limit is of interest to us. Given that the settling time is equal to the value of $i$ in Equation (220), we get

$$
\begin{equation*}
\beta\left(\frac{1-\alpha^{n-d+1}}{1-\alpha}\right)=\frac{0.975 \beta}{1-\alpha}, \tag{223}
\end{equation*}
$$

which can be solved for $n$ to give

$$
\begin{equation*}
n=\frac{\ln 0,025}{\ln \alpha}+d-1 \tag{224}
\end{equation*}
$$

In our case, we obtain

$$
\begin{equation*}
n=\frac{\ln 0,025}{\ln 0,75}+1-1=12,8=13 \tag{225}
\end{equation*}
$$

Note that $n$ is always an integer, so we need to round the resulting value up to the nearest integer value.

The $3 \times 3$ dynamic matrix $\mathcal{A}$ is then

$$
\mathcal{A}=\left[\begin{array}{cccc}
a_{1} & 0 & \cdots & 0  \tag{226}\\
a_{2} & a_{1} & \ddots & \vdots \\
\vdots & \vdots & \ddots & 0 \\
\vdots & \vdots & & a_{1} \\
a_{p} & a_{p-1} & \cdots & a_{p-m+1}
\end{array}\right]=\left[\begin{array}{ccc}
2 & 0 & 0 \\
3.5 & 2 & 0 \\
4.625 & 3.5 & 2
\end{array}\right]
$$

This gives

$$
\left.\begin{array}{rl}
\mathcal{A}^{T} \mathcal{Q} \mathcal{A}+\mathcal{R} & =\left[\begin{array}{ccc}
2 & 0 & 0 \\
3.5 & 2 & 0 \\
4.625 & 3.5 & 2
\end{array}\right]^{T}\left[\begin{array}{lll}
1 & 0 & 0 \\
0 & 1 & 0 \\
0 & 0 & 1
\end{array}\right]\left[\begin{array}{ccc}
2 & 0 & 0 \\
3.5 & 2 & 0 \\
4.625 & 3.5 & 2
\end{array}\right]+\left[\begin{array}{lll}
1 & 0 & 0 \\
0 & 1 & 0 \\
0 & 0 & 1
\end{array}\right]  \tag{227}\\
& =\left[\begin{array}{ccc}
38.640 & 625 & 23.1875
\end{array} 9.25\right. \\
23.1875 & 17.25
\end{array}\right]
$$

The required inverse is then

$$
\begin{align*}
\left(\mathcal{A}^{T} \mathcal{Q} \mathcal{A}+\mathcal{R}\right)^{-1} & =\left[\begin{array}{ccc}
38.640625 & 23.1875 & 9.25 \\
23.1875 & 17.25 & 7 \\
9.25 & 7 & 5
\end{array}\right]^{-1}  \tag{228}\\
& =\left[\begin{array}{ccc}
0.134046 & -0.184200 & 0.009896 \\
-0.184200 & 0.387349 & -0.201518 \\
0.009896 & -0.201518 & 0.463818
\end{array}\right]
\end{align*}
$$

For the controller gain, we obtain the result

$$
\begin{align*}
\vec{K}_{C} & =\left(\mathcal{A}^{T} \mathcal{Q} \mathcal{A}+\mathcal{R}\right)^{-1} \mathcal{A}^{T} \mathcal{Q}^{T} \\
& =\left[\begin{array}{ccc}
0.134046 & -0.184200 & 0.009896 \\
-0.184200 & 0.387349 & -0.201518 \\
0.009896 & -0.201518 & 0.463818
\end{array}\right]\left[\begin{array}{ccc}
2 & 0 & 0 \\
3.5 & 2 & 0 \\
4.625 & 3.5 & 2
\end{array}\right]^{T}\left[\begin{array}{lll}
1 & 0 & 0 \\
0 & 1 & 0 \\
0 & 0 & 1
\end{array}\right]  \tag{229}\\
& =\left[\begin{array}{ccc}
0.268091 & 0.100759 & -0.004948 \\
-0.368400 & 0.129997 & 0.100759 \\
0.019792 & -0.368400 & 0.268091
\end{array}\right]
\end{align*}
$$

The reference signal for the next three sample periods is

$$
\vec{r}=\left[\begin{array}{l}
1  \tag{230}\\
1 \\
1
\end{array}\right]
$$

The predicted uncontrolled position will be equal to the steady state value, i.e. zero since no control has been made. This implies that

$$
\vec{y}^{*}=\left[\begin{array}{l}
0  \tag{231}\\
0 \\
0
\end{array}\right]
$$

The controller action is then

$$
\begin{align*}
\Delta \vec{u} & =\vec{K}_{C} \vec{e}=\vec{K}_{C}\left(\vec{r}-\vec{y}^{*}\right) \\
& =\left[\begin{array}{ccc}
0.268091 & 0.100759 & -0.004948 \\
-0.368400 & 0.129997 & 0.00759 \\
0.019792 & -0.368400 & 0.268091
\end{array}\right]\left[\begin{array}{l}
1-0 \\
1-0 \\
1-0
\end{array}\right]  \tag{232}\\
& =\left[\begin{array}{c}
0.363902 \\
-0.137 \\
-0.080 \\
-0
\end{array}\right]
\end{align*}
$$

Only the first controller action $\Delta u_{1}=0.363902$ will be implemented. At the next time point, Equation (232) will be recomputed using the new values for $r$ and $y^{*}$.

For a multivariate system, the vectors and matrices become "supervectors" and "supermatrices", meaning that a vector consists of many vectors, for example

$$
\Delta \vec{u}=\left[\begin{array}{c}
\Delta \vec{u}_{1}  \tag{233}\\
\vdots \\
\Delta \vec{u}_{h}
\end{array}\right]
$$

where $\Delta \vec{u}_{i}$ is the input vector for the $i^{\text {th }}$ input. Let us consider a MIMO System with $s$ outputs and $h$ inputs. This gives a dynamic matrix $\mathcal{A}$ with the following form

$$
\mathcal{A}=\left[\begin{array}{ccc}
\mathcal{A}_{11} & \cdots & \mathcal{A}_{1 h}  \tag{234}\\
\vdots & & \vdots \\
\mathcal{A}_{s 1} & \cdots & \mathcal{A}_{s h}
\end{array}\right]
$$

where $\mathcal{A}_{i j}$ is the dynamic matrix between the $j^{\text {th }}$ input and $i^{\text {th }}$ output. From Equation (213), it follows that

$$
\begin{equation*}
y_{i, t+l}^{*}=y_{i, t}-\sum_{j=1}^{h} \sum_{k=1}^{n}\left(a_{i j, k}-a_{i j, l+k}\right) \Delta u_{j, t-k} \tag{235}
\end{equation*}
$$

and

$$
\begin{equation*}
\Delta u_{j}=\sum_{i=1}^{p s} k_{j i}\left(r_{i}-y_{i}^{*}\right), \quad j=k m+1, k=0,1, \ldots, h-1 \tag{236}
\end{equation*}
$$

where $a_{i j, k}$ is the $k^{\text {th }}$ coefficient of the step response for the process between the $j^{\text {th }}$ input and $i^{\text {th }}$ output and $\Delta u_{j, t-k}$ is the change in the control action for the $j^{\text {th }}$ input at time point $t-k$.

## Section 5.5: Advanced Control Strategies

In addition to the control strategies mentioned so far, there exist various useful strategies that can be combined with the aforementioned approach to obtain a better result. Such methods include the Smith predictor, deadbanding, squared control, ratio control, input position control, and characterisation of nonlinearities.

## Section 5.5.1: Smith Predictor

The Smith predictor is an approach that seeks to minimise the effect of time delay on the control strategy. It is useful when there is a larger time delay in the system that needs to be dealt with. The control strategy for this approach is shown in Figure 83.


Process-Model Deviation

Figure 83: Block Diagram for Smith Predictor Control

## Section 5.5.2: Deadbanding and Gain Scheduling

Deadbanding refers to the idea that when the control error is within some band of the reference point, then there is no need to perform additional control. Deadbanding means that as long the value is within the band then no control action will be made. In turn, once the error strays outside the deadband, then control will be performed. Mathematically, this can be written as

$$
u_{t}=\left\{\begin{array}{cc}
0 & \left|\varepsilon_{t}\right|<K_{d b}  \tag{237}\\
G_{c} & \text { otherwise }
\end{array}\right.
$$

where $K_{d b}$ is the deadband constant. The deadband can be specified as either some fixed value of the setpoint (for example, within $5^{\circ} \mathrm{C}$ of the setpoint) or as a percentage (within $5 \%$ of the setpoint). Deadbanding is useful when tight control is not desired, that is, small deviations are permissible. It is often used with level control for surge tanks as these will often need to maintain a general level rather than a specific value. Furthermore, the controller can be made more aggressive outside the band to quickly drive the system to the desired value.

A more generalised approach to deadbanding is called gain scheduling, where the controller gain changes depending on the value of the scheduling variable. Most often, the domain of the scheduling variable is partitioned into different regions. For each region, a different controller gain may be assigned. Mathematically, this can be written as

$$
u_{t}=G_{c}\left\{\begin{array}{cc}
K_{c, 1} & s_{t}<K_{g s, 1}  \tag{238}\\
K_{c, 2} & K_{g s, 1} \leq s_{t}<K_{g s, 2} \\
\vdots & \vdots \\
K_{c, n} & K_{g s, \mathrm{n}}<s_{t}
\end{array}\right.
$$

where $K_{c, i}$ is the controller gain for the $i^{\text {th }}$ region, $K_{g s, i}$ is the $i^{\text {th }}$ scheduling limit, and $s_{t}$ is the scheduling variable. This approach allows for a nonlinear process to be controller with a series of linear controllers, which would not be possible with a single controller.

## Section 5.5.3: Squared Control

Squared control is useful if it is desired to penalise large deviation from the setpoint more strongly than those close to the setpoint. This control law can be written as

$$
\begin{equation*}
u_{t}=K_{c} \operatorname{sign}\left(e_{t}\right) e_{t}^{2} \tag{239}
\end{equation*}
$$

where sign is the sign function that returns -1 if $e_{t}$ is negative, 0 if the value is zero, and 1 otherwise.

## Section 5.5.4: Ratio Control

In certain systems, it may be desirable to keep two variables in a constant ratio, for example, in a mixing process, to keep the composition of the mixture constant, the ratio between the flowrates of the two inlets should be maintained constant. In such cases, we can implement ratio control to make sure that the ratio, $R$, is maintained at the desired value. A schematic for ratio control is shown in Figure 84. The ratio control is implemented by the multiplication block that takes the measured value of the solids and multiplies it by the desired ratio set by the VC controller.

When implementing ratio control, the ratio itself should not be controlled, but rather the ratio controller should set the setpoint for one of the variables using the ratio, for example, if $R=$ $u_{1} / u_{2}$, then we could set $u_{2}=R u_{1}$ and use this $u_{2}$ as the setpoint to the $u_{2}$ controller. In all cases, division should be avoided since if one of the variables becomes zero, this will then cause an error. As well, it should be noted that ratio control is implemented using the absolute values of the variables and not their deviational values as is common with other forms of control.


Figure 84: Ratio Control with Trim Feedback Control
Ratio control is useful when it is desired to control some intensive (amount-invariant) variable, such as composition, density, viscosity, or temperature, using extensive (amountvarying) variables, such as flowrates. Note that when implementing ratio control, we need to be able to scale all extensive variables by the same amount and keep all independent intensive variables constant. Thus, for example, a heat exchanger cannot be effectively controlled by ratio control, since it is not feasible to change the surface area during operation. One solution to the problem of keeping all independent intensive variables constant is to implement a form of feedback control on the ratio itself. Such control is often called feedback trim control. It shown using dashed lines in Figure 84, where the ratio is set based on the measured and desired values for $y$. Another advantage of ratio control is that it does not require a model to implement meaning that we can obtain feedforward-like control without requiring any models of the system.

## Section 5.5.5: Input-Position Control

Input-position control, also called valve-position control, allows controlling multi-input, single-output systems using more advanced methods. In this control strategy, it is normally assumed that there are two inputs: a fast-acting but expensive (or otherwise restricted) input, $u_{1}$, and a slow-acting but cheap (or otherwise abundant) input, $u_{2}$. An example would be cooling an exothermic reactor using both coolant (expensive, but fast) and cooling water (abundant, but slow). In general, the cheap input is used for controlling the process, but the expensive input can be used to improve the speed of response. Basically, at steady state, it is assumed that the expensive input will attain some steady-state value $u_{1 \mathrm{~s}}$, so that the overall process will be controlled by $u_{2}$. A schematic of this control strategy is shown in Figure 85.


Figure 85: Input-Position Control
Since input-position control can be treated as a modification of cascade control, the fast loop given by $u_{1}$ is tuned first and then the slower loop given by $u_{2}$. This also implies that the rule regarding time constants between the loops for cascade control should hold, that is, $\tau_{c 2} / \tau_{c 1}$ should be between 4 and 10. In general, both controllers are designed as simple PI controllers without any anti-windup features.

## Section 5.5.6: Nonlinear Characterisation

Nonlinear characterisation allows nonlinearities present in actuators to be dealt with separately from the controller itself. The nonlinear characterisation block converts the linear controller output into the nonlinear actuator output, for example, if the controller gives a flow rate, then this block can convert the flow rate into the corresponding value. Very often, this block is a look-up table with interpolation between the given data points. Such an approach can allow for the system to be focused on those areas that are highly nonlinear.

## Section 5.5.7: Bumpless Transfer

Bumpless transfer is the smooth transfer from one control mode or strategy to another without causing any visible, undesirable changes (bumps) in the process variables. This often arises from a mismatch between the requested input values in the two modes. In general, bumpless transfer is ensured using an appropriate anti-wind-up method. However, there is one case, where it may be necessary to take special action and that is when transferring from manual to automatic mode. In these cases, the automatic input value will not be exactly equal to the manual value meaning that it can lead to a bump in the output. The solution is to match the current manual value with the expected automatic value. As well, it may be helpful to reset the integration so that it equals zero at the moment of the switch.

## Section 5.6: Chapter Problems

Problems at the end of the chapter consist of three different types: (a) Basic Concepts (True/False), which seek to test the reader's comprehension of the key concepts in the chapter; (b) Short Exercises, which seek to test the reader's ability to compute the required parameters for a simple data set using simple or no technological aids. This section also includes proofs of theorems; and (c) Computational Exercises, which require not only a solid comprehension of the basic material, but also the use of appropriate software to easily manipulate the given data sets.

## Section 5.6.1: Basic Concepts

Determine if the following statements are true or false and state why this is the case.

1) In open-loop control, the controller uses the measured output to determine how to control the process.
2) In closed-loop control, it is important to be able to measure the output.
3) Plant-model mismatch can derail open-loop control.
4) Closed-loop control should provide stable, biasfree, and robust control.
5) A state-space controller requires measurement of the states.
6) In state-space control, the separation principle implies that we can separate the design of the observer and controller.
7) A proportional-only controller will always provide biasfree control.
8) The integral term considers the effect of future values on the system.
9) Integral wind-up occurs when the disturbance changes a lot causing the manipulated variable to fluctuate.
10) Derivative kick occurs in PID controllers when there is a step change in the setpoint.
11) Increasing the absolute value of $K_{c}$ in a PI controller will cause the system to become more stable.
12) The proportional term in a PID controller considers the current values of the error.
13) A PI controller can be tuned using the IMC method.
14) The rise time of a closed-loop system can be used to determine how well a controller regulates the disturbance.
15) The settling time is defined as three times the closed-loop time constant.
16)For regulatory controller performance, it is easy to specify the desired values.
17)Feedforward control seeks to minimise the effects of unmeasurable disturbances on the process.
16) A decoupler is a type of feedforward control.
17) In feedforward control, it may be necessary to remove terms such as time delay or unstable zeros from the final controller.
18) Interlocking occurs when a series of discrete events must be fulfilled before some action can occur.
19) Batch control is an example of sequential control.
20) Supervisory control allows us to create a network of controllers each of which can be controlling another controller.
21) In cascade control, the master controller should always be faster than the slave controller.
22) Model predictive control allows constraints and economic conditions to be considered in controlling the process.
23) Model predictive control requires good models.

## Section 5.6.2: Short Questions

26)Describe in words how the following control works. What kind of methods are being used? What are the objectives and what are some potential disturbances?
a. An elevator reaching a given floor.
b. Temperature control in a home oven.
c. Temperature control in a fridge/freezer.
d. Driving a car on the highway.
e. Driving a car on the highway using cruise control.
27) Design PI controllers using the formulae in Table 26 for the following processes:
a. $\quad G_{p}=\frac{2}{30 s+1} e^{-15 s}$
b. $\quad G_{p}=\frac{-2}{30 s+1} e^{-15 s}$
c. $\quad G_{p}=\frac{2}{3 s+1} e^{-150 s}$
d. $\quad G_{p}=\frac{-5}{s-0.5} e^{-15 s}$

You should use the smallest $\tau_{c}$ possible.
28) Design PID controllers using the formulae in Table 27 for the same processes as in Question 27). You should use the smallest $\tau_{c}$ possible.
29) Simulate the controllers from Questions 27) and 28) for a setpoint change of +2 and -2 . You can assume that there is no disturbance affecting the process. Explain what you see. Which controller would you recommend for each process? Why?
30)Design dynamic feedforward controllers for the following processes.
a. $\quad G_{p}=\frac{2}{30 s+1} e^{-15 s}, G_{d}=\frac{-2}{15 s+1} e^{-5 s}$
b. $\quad G_{p}=\frac{-2(5 s-1)}{(30 s+1)(25 s+1)} e^{-15 s}, G_{d}=\frac{2}{5 s+1} e^{-20 s}$
c. $\quad G_{p}=\frac{2}{(3 s+1)(10 s+1)} e^{-s}, G_{d}=\frac{-2}{15 s+1} e^{-5 s}$
31)Design static feedforward controllers for the processes in Question 31). Simulate and compare the performance of the two types of feedforward controllers for a setpoint change of +2 . You can assume that the disturbance is driven by Gaussian, white noise.
32) Consider a cascade loop with the following two transfer functions:

$$
\begin{equation*}
G_{p, \text { slave }}=\frac{2}{3 s+1} e^{-s}, G_{p, \text { master }}=\frac{-2}{30 s+1} e^{-20 s} \tag{240}
\end{equation*}
$$

Design appropriate PI controllers for this cascade loop. What aspects should you take into consideration when designing this controller? Simulate the system.
33) Consider the same situation as in Question 32), but now you also wish to ensure that the slave loop does not have any oscillations and reaches the setpoint as fast as possible. Using simulations, design appropriate controllers for this system.
34) You need to design a controller for a process whose model has been determined experimental to be

$$
\begin{equation*}
\hat{G}_{p}=\frac{1}{15 s+1} e^{-10 s} \tag{241}
\end{equation*}
$$

However, you do not know how well the experimental model reflects the true process. Using the IMC tuning method, design a PID controller assuming that the experimental model is correct. Knowing that the model may be incorrect, do not pick a too small $\tau_{c}$. Simulate your model assuming Equation (241) is correct. If your controller is satisfactory, then try simulating the closed-loop assuming that the true process model is given by
a. Small Mismatch: $G_{p}=\frac{1.1}{16 s+1} e^{-11 s}$
b. Large Mismatch in Gain: $G_{p}=\frac{2}{14 s+1} e^{-9 s}$
c. Large Mismatch in the Time Constant: $G_{p}=\frac{1.05}{4 s+1} e^{-10 s}$
d. Large Mismatch in Time Delay: $G_{p}=\frac{1}{14 s+1} e^{-25 s}$

If your system is unstable, design a new controller that makes your system stable using both the experimental model and the true model. What conclusions can you draw about the controller design?
35)For the following two-input, two-output system, design a decoupler between $u_{2}$ and $y_{1}$. The PID couplings are $y_{1}$ with $u_{1}$ and $y_{2}$ with $u_{2}$.

$$
\left[\begin{array}{l}
y_{1}  \tag{242}\\
y_{2}
\end{array}\right]=\left[\begin{array}{cc}
\frac{10}{(10 s+1)} e^{-15 s} & \frac{-2}{(15 s+1)(10 s+1)} e^{-10 s} \\
\frac{5(10 s+1)}{(15 s+1)(20 s+1)} e^{-25 s} & \frac{5}{(10 s-1)} e^{-30 s}
\end{array}\right]\left[\begin{array}{l}
u_{1} \\
u_{2}
\end{array}\right]
$$

36) Simulate the system given by Equation (242) with and without the decoupler. What is the effect of the decoupler? The PI controllers are given as

$$
\begin{align*}
& \varepsilon_{1}=\frac{1}{10}\left(1+\frac{1}{25 s}\right)\left(r_{1}-y_{1}\right)  \tag{243}\\
& \varepsilon_{2}=\frac{1}{5}\left(1+\frac{1}{40 s}\right)\left(r_{2}-y_{2}\right)
\end{align*}
$$

37) Using the DMC method, design a model predictive controller for the following system:

$$
\begin{equation*}
y_{t}=\frac{z^{-10}}{1-0.25 z^{-1}} u_{t} \tag{244}
\end{equation*}
$$

38) Using the DMC method, design a model predictive controller for the following system:

$$
\left[\begin{array}{l}
y_{1}  \tag{245}\\
y_{2}
\end{array}\right]=\left[\begin{array}{ll}
\frac{2}{10 s+1} e^{-5 s} & \frac{-2}{5 s+1} e^{-15 s} \\
\frac{5}{15 s+1} e^{-10 s} & \frac{4}{10 s+1} e^{-10 s}
\end{array}\right]\left[\begin{array}{l}
u_{1} \\
u_{2}
\end{array}\right]
$$

Assume that the sampling time is 1 s .

## Chapter 6: Boolean Algebra

Boolean algebra is the algebra of binary variables that can only take two values, for example, true and false or 1 and 0 . It is very useful for solving problems in logic and is a requirement for good programming. George Boole (1815-1864) is the discoverer of Boolean algebra.

A Boolean expression is a group of elementary terms that are linked with connectors (operators). The mathematical space in which a Boolean algebra is defined will be denoted using a double-struck $\mathrm{B}(\mathbb{B}, \mathrm{U}+1 \mathrm{D} 539)$.

## Section 6.1: Boolean Operators

In Boolean algebra, there are 5 operators: conjunction, disjunction, negation, implication, and equivalency. These 5 operators are described in Table 29. It should be noted that negation is a unary operator, that is, it only requires a single variable. All other operators are binary operators, that is, they require two variables. When we write a Boolean expression, it is common to write conjunction as multiplication and disjunction as addition, that is, $\mathrm{a} \wedge \mathrm{b}$ is written as ab and a V b as a +b .

Table 29: Boolean operators, where $a, b \in \mathbb{B}$

| Operator | Symbol | Statement | Other <br> Representations | Word Representation |
| :---: | :---: | :---: | :---: | :---: |
| Conjunction | $\wedge(\mathrm{U}+2227)$ | a ^ b | $\begin{aligned} & \mathrm{a} \cdot \mathrm{~b}, \mathrm{ab}, \\ & \mathrm{a} \text { AND } \mathrm{b}, \\ & \mathrm{a} \& \mathrm{~b} \end{aligned}$ | AND |
| Disjunction | $\mathrm{V}(\mathrm{U}+2228)$ | $\mathrm{a} \vee \mathrm{b}$ | $\begin{aligned} & a+b, a \text { OR } \\ & b, a \\| b \end{aligned}$ | OR |
| Negation | $\neg(\mathrm{U}+00 \mathrm{AC})$ | $\neg \mathrm{b}, \mathrm{b}^{\prime} \overline{\mathrm{b}}$ | NOT b, ! b | NOT |
| Implication | $\rightarrow(\mathrm{U}+2192)$ | $\mathrm{a} \rightarrow \mathrm{b}$ | - | IF-THEN |
| Equivalence | $\leftrightarrow(\mathrm{U}+2194)$ | $\mathrm{a} \leftrightarrow \mathrm{b}$ | - | IF-\&ONLY-IF |

All Boolean expressions can be written using $\wedge, \vee, \neg, 0$ and 1 . The precedence of the operators is such that $\wedge$ has a higher precedence than $\vee$, for example, $a \vee b \wedge c=a \vee(b$ $\wedge c)$.

## Section 6.2: Boolean Axioms and Theorems

For the Boolean space $\mathbb{B}=\{0,1\}$ with variables $a, b$, and $c$ and operators $\wedge$ und $\vee$, the following axioms hold:

1) Closure
a. $a \vee b \in \mathbb{B}$
b. $a \wedge b \in \mathbb{B}$

## 2) Commutativity

a. $a \vee b=b \vee a$
b. $a \wedge b=b \wedge a$
3) Distributivity
a. $a \wedge(b \vee c)=(a \wedge b) \vee(a \wedge c)$
b. $a \vee(b \wedge c)=(a \vee b) \wedge(a \vee c)$
4) Identity
a. $\mathrm{a} \vee 0=\mathrm{a}$
b. $a \wedge 1=a$

## 5) Complementation

a. $a \wedge \neg a=0$
b. $a \vee \neg a=1$
6) Annihilator
a. $a \vee 1=1$
b. $a \wedge 0=0$

## 7) Associativity

a. $(a \wedge b) \wedge c=a \wedge(b \wedge c)$
b. $(\mathrm{a} \vee \mathrm{b}) \vee \mathrm{c}=\mathrm{a} \vee(\mathrm{b} \vee \mathrm{c})$
8) Idempotence
a. $\mathrm{a} \vee \mathrm{a}=\mathrm{a}$
b. $a \wedge a=a$

## 9) Involution

a. $\neg \neg a=a$
10)Absorption
a. $a \wedge(a \vee c)=a$
b. $a \vee(a \wedge c)=a$

With these axioms, it is possible to define a complete Boolean algebra.
An important law is De Morgan's Law:

$$
\begin{align*}
& \neg(\mathrm{a} \vee \mathrm{~b})=\neg \mathrm{a} \wedge \neg \mathrm{~b}  \tag{246}\\
& \neg(\mathrm{a} \wedge \mathrm{~b})=\neg \mathrm{a} \vee \neg \mathrm{~b} \tag{247}
\end{align*}
$$

This law can be used to simplify many complex Boolean expressions or convert between two representations.

## Section 6.3: Boolean Functions

A Boolean function is a function where all variables are Boolean variables, for example,

$$
\begin{equation*}
F=f\left(X_{1}, X_{2}\right) \text {, where } X_{1}, X_{2} \in \mathbb{B} \tag{248}
\end{equation*}
$$

Typically, a Boolean variable is shown using a capital letter.
A truth table shows the values of the expression for each combination of values of the variables, that is, each possible input value. This implies that a truth table will have $2^{N}$ rows, where $N$ is the number of variables (input variables), for example, for two variables, we will have $2^{2}=4$ rows and for four variables $2^{4}=16$ rows.

The truth table for the Boolean AND operator is shown in Table 30 (left). Since the AND operator requires that both inputs be TRUE, there is only a single row that evaluates to TRUE. For the Boolean OR operator, the truth table is shown in Table 30 (right). Since the OR operator requires that at least one of the inputs be TRUE, there are three rows that evaluate to TRUE and one, where both inputs are FALSE, that evaluates to FALSE.

Table 30: Truth table for the Boolean Operators (left) AND and (right) OR

| AND |  |  |
| :---: | :---: | :---: |
| $\mathbf{A}$ | $\mathbf{B}$ | $\mathbf{A B}$ |
| 0 | 0 | 0 |
| 0 | 1 | 0 |
| 1 | 0 | 0 |
| 1 | 1 | 1 |


| OR |  |  |
| :---: | :---: | :---: |
| $\mathbf{A}$ | $\mathbf{B}$ | $\mathbf{A + B}$ |
| 0 | 0 | 0 |
| 0 | 1 | 1 |
| 1 | 0 | 1 |
| 1 | 1 | 1 |

Example 31: Truth Table
What is the truth table for the Boolean function $F=A B^{\prime}$ ?

## Solution

Since we have two variables (A and B), $N=2$. This means that there will be $2^{2}=4$ rows in the truth table.

| A | B | $\mathbf{B}^{\prime}$ | $\mathbf{F}$ |
| :---: | :---: | :---: | :---: |
| 0 | 0 | 1 | 0 |
| 0 | 1 | 0 | 0 |
| 1 | 0 | 1 | 1 |
| 1 | 1 | 0 | 0 |

There are two common representations for Boolean functions: the sum-of-products (SOP) form and the product-of-sums (POS) form.

## Section 6.3.1: Sum-of-Products Form and Minterms

The sum-of-products form is a representation of a Boolean function where all products (or terms) are products of single variable, for example, $F=A B C+B^{\prime} C D E '+A^{\prime} B^{\prime}$ or $\mathrm{F}=\mathrm{BCDE}+\mathrm{AB} \mathrm{B}^{\prime}+\mathrm{HI}^{\prime}+\mathrm{C}$. Using the axioms, it is possible to convert all functions to a sum-of-products form.

Example 32: Sum-of-Products Form
Which of the following functions are in the sum-of-product form:

1) $\mathrm{F}=\mathrm{ABC}+\mathrm{B}^{\prime} \mathrm{CDE}{ }^{\prime}$
2) $F=(A+B) C$
3) $F=A B C+B^{\prime}(D+E)$
4) $\mathrm{F}=\mathrm{BC}+\mathrm{DE}$ '?

## Solution

Only 1) and 4) are in the sum-of-product form, since these representations are a sum of the products of individual variables. In 2) and 3), we have a product of many variables (A and B) in 2) and ( $D$ and $E$ ) in 3).

Example 33: Converting into the Sum-of-Products Form
Convert this function into the sum-of-products form: $\mathrm{F}=(\mathrm{A}+\mathrm{B}) \mathrm{C}$.

## Solution

Using the distributive property, we can obtain the sum-of-products form, that is,

$$
F=(A+B) C=A C+B C
$$

A minterm is a row in the truth table where the value is " 1 ". The symbol for a minterm is $m_{i}$, where $i$ is the decimal row number. The compact sum-of-products form is $\Sigma m(i, \ldots)$, where $i$ is the row number of the minterms. When the minterms are converted into a functional representation, each variable, whose value is 1 , will be written in its plain form, while each variable, whose value is 0 will be written in its negated form, for example, for $m_{2}$, the minterm with the value $010_{2}$, the term will be given as $A^{\prime} B C^{\prime}$.

## Example 34: Compact Sum-of-Products Form

What is the compact sum-of-products form for the function $F=(A+B) C$.

## Solution

First, we require a truth table. Then, we can simply find the rows with a value of 1.

| A | B | C | F |
| :---: | :---: | :---: | :---: |
| 0 | 0 | 0 | 0 |
| 0 | 0 | 1 | 0 |
| 0 | 1 | 0 | 0 |


| $\mathbf{A}$ | $\mathbf{B}$ | $\mathbf{C}$ | $\mathbf{F}$ |
| :---: | :--- | :--- | :--- |
| 0 | 1 | 1 | 1 |
| 1 | 0 | 0 | 0 |
| 1 | 0 | 1 | 1 |
| 1 | 1 | 0 | 0 |
| 1 | 1 | 1 | 1 |

The rows in red are the rows, where the function has a value of 1 . We can find the decimal row number simply by converting the binary representation into a decimal value, that is, for the first red row (011)

$$
011_{2}=2+1=3
$$

and

$$
\begin{gathered}
101_{2}=2^{2}+1=5 \\
111_{2}=2^{2}+2+1=7
\end{gathered}
$$

Thus, the compact sum-of-products form is

$$
\Sigma m(3,5,7) .
$$

It should be noted that the ordering of the Boolean variables is important. When the ordering is changed, then the compact sum-of-products form will also change.

## Section 6.3.2: Product-of-Sums Form and Maxterms

The product-of-sums form is a representation of a Boolean function, where all the factors are sums of individual variables, for example, $F=(A+B+C)\left(B^{\prime}+C+D\right) E^{\prime}$ or $F=(A+B) C$. Using the axioms, it is possible to convert all functions into a product-of-sums form.

## Example 35: Product-of-Sums Form

Which of the following Boolean functions are in the product-of-sums form:

1) $\mathrm{F}=\mathrm{ABC}+\mathrm{B}^{\prime} \mathrm{CDE}{ }^{\prime}$
2) $F=(A+B) C$
3) $F=(A+B+C) B^{\prime}(D+E)$
4) $\mathrm{F}=\mathrm{BC}+\mathrm{DE}$ ?

## Solution

Only 2) and 3) are in the product-of-sums form, since these representations are products of sums with individual variables. In 1) and 4 ), we have a sum of many variables ( $A, B$, and $C$ ) in 1) and ( $B, C, D$, and $E$ ) in 4).

## Example 36: Converting into the Product-of-Sums Form

Convert the function $\mathrm{F}=\mathrm{AC}+\mathrm{BC}$ into its product-of-sums form.

## Solution

Using the distributive property, we can obtain the product-of-sums form, that is,

$$
F=A C+B C=(A+B) C
$$

A maxterm is a row in the truth table with the value of 0 . The symbol for a maxterm is $M_{i}$, where $i$ is the decimal value of the row. The compact product-of-sums form is $\Pi M(i, \ldots)$, where $i$ is the decimal row value of the corresponding maxterm. When maxterms are converted into a functional representation, then each variable with the value of 0 is converted as written, while each variable with a value of 1 is written in its negated form, for example, $M_{2}$, the maxterm with the value $010_{2}$, would be written a $A+B^{\prime}+C$.

## Example 37: Compact Product-of-Sums Form

What is the compact product-of-sums form for the function $F=(A+B) C$.

## Solution

First, we require the truth table for the function. Then, we can easily find the rows with a value of 0 .

| $\mathbf{A}$ | $\mathbf{B}$ | $\mathbf{C}$ | $\mathbf{F}$ |
| :---: | :---: | :--- | :--- |
| 0 | 0 | 0 | 0 |
| 0 | 0 | 1 | 0 |
| 0 | 1 | 0 | 0 |
| 0 | 1 | 1 | 1 |
| 1 | 0 | 0 | 0 |


| $\mathbf{A}$ | $\mathbf{B}$ | $\mathbf{C}$ | $\mathbf{F}$ |
| :---: | :--- | :--- | :--- |
| 1 | 0 | 1 | 1 |
| 1 | 1 | 0 | 0 |
| 1 | 1 | 1 | 1 |

The rows in red are the rows where the function has a value of 0 . We can obtain the row number by converting the binary values into a decimal representation, for example, for the first row (000)

$$
000_{2}=0+0+0=0
$$

and

$$
\begin{aligned}
& 001_{2}=0+0+1=1 \\
& 010_{2}=0+2+0=2 \\
& 100_{2}=2^{2}+0+0=4 \\
& 110_{2}=2^{2}+2+0=6
\end{aligned}
$$

Therefore, the compact product-of-sums form is

$$
\Pi M(0,1,2,4,6)
$$

It is important to note that the order of Boolean variables is important. When the order is changed, then the compact product-of-sums form will also change. Also, it is obvious that any terms in the sum-of-products form are not in the product-of-sums form!

## Section 6.3.3: Don't-Care Values

It can happen that a given logical situation cannot occur. Such cases are often denoted using * or $\times$ in the truth table. It is then up to the engineer to determine which value ( 0 or 1 ) is best. Such cases are called don't-care values.

## Example 38: Don't Cares

What is the truth table for the Boolean function $F=A B$, when the case $A=B=0$ is not possible?

## Solution

The truth table is

| $\mathbf{A}$ | $\mathbf{B}$ | $\mathbf{F}$ |
| :---: | :---: | :---: |
| 0 | 0 | $\times$ |


| A | $\mathbf{B}$ | F |
| :---: | :---: | :---: |
| 0 | 1 | 0 |
| 1 | 0 | 0 |
| 1 | 1 | 1 |

## Section 6.3.4: Duality

Duality in Boolean algebra has the following definition:

- Replace all AND by OR.
- Replace all OR by AND.
- Replace 0 by 1 .
- Replace 1 by 0.

Duality is an important property when we are simplifying or minimising an expression. Often the dual of the representation can be easier to work with.

## Example 39: Dual of a Function

What is the dual of the function $F=(A+B) C$ ?

## Solution

$$
\begin{gathered}
F=(A+B) C \\
F^{D}=[(A+B) C]^{D} \\
F^{D}=[(A+B)]^{D}+C D \\
F^{D}=A B+C
\end{gathered}
$$

The dual of a function is often represented by a superscript $D$.

## Section 6.4: Minimising a Boolean Function

When a logical expression is to be implemented using electronic gates, it is often important to minimise the expression, that means, that the number of required gates is minimised. Often, each Boolean operation requires a separate logic gate. There are three ways to minimise a Boolean function:

1) Manually using the axioms and theorems of Boolean algebra;
2) Using a Karnaugh map; and
3) Using the Quine-McCluskey algorithm, which is often used for large (greater than 10 variables) Boolean functions.

In general, the Karnaugh map is the best method for minimising a Boolean function.

## Section 6.4.1: Karnaugh Map

A Karnaugh map is a visual algorithm that minimises a Boolean function so that the largest number of groups of 0 and 1 are found. We search for groups of 0 to find the minimal product-of-sums representation and groups for 1 to find the minimal sum-of-products representation. The 0 's and 1 's can only be arranged in groups of $2^{n}$, where $n \in \mathbb{N}$. There are simple Karnaugh maps for $2,3,4$, and 5 variables.

The Karnaugh map for two variables is shown in Figure 86. The diagram is created so that on one axis lies one variable and on the other axis the other variable. Then, for the minimal sum-of-products representation, we circle the largest group with a size that is a power of 2 that covers as many of the 1 as possible.


Figure 86: Karnaugh map for the function $F=B^{\prime}$
The Karnaugh map for three variables is shown in Figure 87. The diagram is created so that on one axis lies one variable and on the other axis the remaining two variables. Each row and column has one of the two possible values ( 0 or 1 ) so that each adjacent row or column only differs by one entry, for example, when the row is 01 , then the next row must be 11 , since only one entry should be changed.

| $A$ | $B$ | $C$ | 0 |
| :---: | :---: | :---: | :---: |
| 1 |  |  |  |
| 0 | 0 | 1 | 0 |
| 0 | 1 | 0 | 1 |
| 1 | 1 | 0 | 0 |
| 1 | 0 | 0 | 1 |

Figure 87: Karnaugh map for the function $F=\Sigma m(0,3,5)$
The Karnaugh map for four variables is shown in Figure 88. Each row and column has two variables. The Karnaugh map for five variables is shown in Figure 89. When we are looking for 1, it is common to not bother with writing the 0 .

|  | $A$ | 0 | 1 | 1 | 0 |
| :---: | :---: | :---: | :---: | :---: | :---: |
| $C$ | $B$ | 0 | 0 | 1 | 1 |
| 0 | 0 | 1 | 0 | 0 | 1 |
| 0 | 1 | 0 | 1 | 0 | 0 |
| 1 | 1 | 1 | 1 | 1 | 1 |
| 1 | 0 | 1 | 1 | 1 | 1 |

Figure 88: Karnaugh map for the function $F=A^{\prime} B D+B^{\prime} C^{\prime} D^{\prime}+C$

|  | $B$ | 0 | 1 | 1 | 0 |
| :---: | :---: | :---: | :---: | :---: | :---: |
| $D$ | $E$ | $C$ | 0 | 0 | 1 |
| 0 | 0 | 1 | 1 | 1 | 1 |
| 0 | 1 |  | 1 | 1 |  |
| 1 | 1 | 1 | 1 | 1 |  |
| 1 | 0 |  |  | $1 / 1$ | 1 |

A $1 / 0$

Figure 89: Karnaugh map for the function $F=\Pi M(2,5,7,9,13,15,16,17,18,20,24,25,27)$

The general procedure for finding the minimal sum-of-products form is:

1) Construct the Karnaugh map.
2) Circle the 1 into the largest groups possible.
3) Write the corresponding sum-of-product form based on the circled groups. For each group, we only write the variables whose value is not changed. Variables with a value of 0 are negated, for example, in Figure 88, for the loop in the first row, A varies between 0 and 1, which means that it is ignored. The other variables remain constant and have a value of 0 . Thus, we write the negated forms of these variables, that is, $B^{\prime} C^{\prime} D^{\prime}$. The general procedure for finding the minimal product-of-sums form is:
4) Construct the Karnaugh map.
5) Circle the 0 into the largest groups possible.
6) Write the corresponding sum-of-product form for $\mathrm{F}^{\prime}$ based on the circled groups. For each group, we only write the variables whose value is not changed. Variables with a value of 0 are negated, for example, in Figure 88, for the loop in the first row, A varies between 0 and 1 , which means that it is ignored. The other variables remain constant and have a value of 0 . Thus, we write the negated forms of these variables, that is, $\mathrm{B}^{\prime} \mathrm{C}^{\prime} \mathrm{D}^{\prime}$.
7) Convert $\mathrm{F}^{\prime}$ to F .

Of course, we can look for the loops using trial and error, but we have no guarantee that we will find the best loops. Therefore, we wish to find a procedure that we can always find the best loops.

Before we can describe such a procedure, we need to define certain words. An implicant is a single 1 or group of 1 's. A prime implicant is an implicant that cannot be further combined, that means, that a single 1 creates a prime implicant if there are not adjacent 1 . Two adjacent 1 's create a prime implicant if they cannot be combined into a group of four 1's, while four adjacent 1 's create a prime implicant if they cannot be combined into a group of eight 1 's. Figure 90 shows the difference between an implicant and a prime implicant. We can state that a sum-of-products form with nonprime implicants is not a minimal form, but not all prime implicants will necessarily be required for the minimal sum-of-products form. An essential prime implicant is a minterm that is covered only by a single prime implicant. All essential prime implicants must be in the minimal sum-of-products form. Thus, we can say that the objective is to find all the essential prime terms. Figure 92 shows the procedure for minimising a Karnaugh map.

|  | $A$ | 0 | 1 | 1 | 0 |
| :---: | :---: | :---: | :---: | :---: | :---: |
| $C$ | $B$ | 0 | 0 | 1 | 1 |
| 0 | 0 | 1 | 1 | 0 | 0 |
| 0 | 1 | 1 | 1 | 1 | 1 |
| 1 | 1 | 0 | 0 | 0 | 0 |
| 1 | 0 | 0 | 0 | 0 | 0 |

- Implicant

Prime Implicant

Figure 90: Prime implicant and implicant

## Example 40: Karnaugh map

For the Karnaugh map shown in Figure 91, find the minimal sum-of-products form.

|  | $A$ | 0 | 1 | 1 | 0 |
| :---: | :---: | :---: | :---: | :---: | :---: |
| $C$ | $B$ | 0 | 0 | 1 | 1 |
| 0 | 0 | 1 | 1 | 0 | 0 |
| 0 | 1 | 1 | 1 | 0 | 0 |
| 1 | 1 | 0 | 1 | 1 | 1 |
| 1 | 0 | 1 | 0 | 0 | 0 |

Figure 91: Karnaugh map for Example 40

## Solution

1) We start with the first 1 (cell: 0000). The largest group that we can find is four 1 . This group is an essential prime implicant.

|  | $A$ | 0 | 1 | 1 | 0 |
| :---: | :---: | :---: | :---: | :---: | :---: |
| $C$ | $D^{B}$ | 0 | 0 | 1 | 1 |
| 0 | 0 | 1 | 1 | 0 | 0 |
| 0 | 1 | 1 | 1 | 0 | 0 |
| 1 | 1 | 0 | 1 | 1 | 1 |
| 1 | 0 | 1 | 0 | 0 | 0 |

2) The next unlooped 1 is in the cell 0010 . This 1 can only be combined with the 1 in cell 0000 . This group is also an essential prime implicant.

|  | $A$ | 0 | 1 | 1 | 0 |
| :---: | :---: | :---: | :---: | :---: | :---: |
| $C$ | $D$ | 0 | 0 | 1 | 1 |
| 0 | 0 | 1 | 1 | 0 | 0 |
| 0 | 1 | 1 | 1 | 0 | 0 |
| 1 | 1 | 0 | 1 | 1 | 1 |
| 1 | 0 | 1 | 0 | 0 | 0 |

3) The next free 1 is in cell 0111 . Here we have two possibilities. We could make the loop with either cell 0101 or cell 1111. This implies that we have a prime implicant.

|  | $A$ | 0 | 1 | 1 | 0 |
| :---: | :---: | :---: | :---: | :---: | :---: |
| $C$ | $D^{B}$ | 0 | 0 | 1 | 1 |
| 0 | 0 | 1 | 1 | 0 | 0 |
| 0 | 1 | 1 | 1 | 0 | 0 |
| 1 | 1 | 0 | 1 | -1 | 1 |
| 1 | 0 | 1 | 0 | 0 | 0 |

4) The next free 1 is in cell 1011, which we can combine with cell 1111. This is the only possibility for combining this cell. Therefore, this group is also an essential prime implicant.

5) The last step is to write the minimal sum-of-product form. First, we will write all the essential prime implicants, that is, $A^{\prime} B^{\prime} D^{\prime}, A^{\prime} C^{\prime}$ and $A C D$. Then, we must select one of the remaining prime implicants ( $A^{\prime} B D$ or $B C D$ ). There is no difference between which of these two prime implicants we select. The minimal sum-of-products form is then

$$
F=A^{\prime} B^{\prime} D^{\prime}+A^{\prime} C^{\prime}+A C D+\left\{A^{\prime} B D \text { or } B C D\right\}
$$



Figure 92: Procedure for minimising a Karnaugh map

## Section 6.5: Chapter Problems

Problems at the end of the chapter consist of three different types: (a) Basic Concepts (True/False), which seek to test the reader's comprehension of the key concepts in the chapter; (b) Short Exercises, which seek to test the reader's ability to compute the required parameters for a simple data set using simple or no technological aids. This section also includes proofs of theorems; and (c) Computational Exercises, which require not only a solid comprehension of the basic material, but also the use of appropriate software to easily manipulate the given data sets.

## Section 6.5.1: Basic Concepts

Determine if the following statements are true or false and state why this is the case.

1) In Boolean algebra, there are three possibilities: $-1,0$, and 1.
2) In Boolean algebra, conjunction is denoted using $\wedge$.
3) In Boolean algebra, negation is denoted using $V$.
4) The Boolean statements $\neg b$ and $b^{\prime}$ have the same meaning.
5) The Boolean operator $\Lambda$ is a unary operator.
6) If $\mathrm{a}=1, \mathrm{~b}=0$, and $\mathrm{c}=1$, then the Boolean statement $\mathrm{a} \vee \mathrm{b} \wedge \mathrm{c}$ evaluates to "false".
7) Boolean algebra does not obey the law of associativity.
8) Boolean algebra obeys the law of distributivity.
9) De Morgan's Law states that $(A+B)^{\prime} \equiv A^{\prime} B^{\prime}$.
10) The function $F=(A+B) C$ is in a sum-of-products form.
11) The function $F=A B C+A^{\prime} B^{\prime} C^{\prime}$ is in a sum-of-products form.
12) The function $F=(A+B) C$ is in a product-of-sums form.
13) The function $F=A B C+A^{\prime} B^{\prime} C^{\prime}$ is in a product-of-sums form.
14)For a function with the representation $F=\Sigma m(1,3,5)$, the minterm $m_{2}$ is equal to 1 .
14) For the function with the representation $F=\Pi M(1,3,5)$, the maxterm $M_{3}$ is equal to 0 .
15) The dual of $F=A+B C$ is $F^{D}=A(B+C)$.
16) We must always assign a "don't care" a value of 1 .
17) All prime implicants are in the minimal sum-of-products form.
18) All essential prime implicants are in the minimal sum-of-products form.
19) A Karnaugh map can be used to determine the minimal product-of-sums form.

## Section 6.5.2: Short Questions

These questions should be solved using only a simple, nonprogrammable, nongraphical calculator combined with pen and paper.
21) Simplify the following function $Z=(A+B)(A+B C)(B+B C)$. Using a truth table, show that the original and simplified functions are the same.
22) Convert the following functions into the sum-of-products form:
a. $Z=(A+B)(A+C)(A+D)(B D C+E)$
b. $Z=(A+B+C)(B+C+D)(A+C)$
23) Convert the following functions into a product-of-sums form:
a. $\quad Z=W+X Y Z$
b. $Z=A B C+A D E+A B F$
24)Find the compact product-of-sums form for the following functions:
a. $F(A, B, C)=A^{\prime}$.
b. $F(A, B, C, D)=A^{\prime} B^{\prime}+A^{\prime} B^{\prime}\left(C D+C D^{\prime}\right)$.
25)Find the compact sum-of-products form for the following functions:
a. $F(A, B, C)=\left(A^{\prime}+B+C\right)(A+C)$.
b. $F(A, B, C, D)=A^{\prime} B^{\prime}+A^{\prime} B\left(C D+C D^{\prime}\right)$.
26) Using a Karnaugh map, find the minimal sum-of-products form for the following functions:
a. $F(A, B, C, D)=A^{\prime} B^{\prime}+A^{\prime} B\left(C D+C D^{\prime}\right)$.
b. $F(A, B, C, D)=\Sigma m(0,1,2,3,6,7)$.
c. $F(A, B, C, D)=\Sigma m(0,4,5,6,8,9,10,11)$.
d. $F(A, B, C, D)=\Sigma m(10,12,14)$.
27)For the Karnaugh map shown in Figure 93, find all prime implicants and essential prime implicants.


Figure 93: Karnaugh map for Question 27

## Chapter 7: PLC Programming

When using a PLC, it is necessary to transfer the required information to the PLC using some methods. There are two possible approaches. We could use an $a d-h o c$ approach that depends on the PLC or situation or we could use some type of standard. It is obvious that using standards is better since we can re-use the information, for example, the same code can be re-used for different cases without having to worry about compatibility problems. The IEC/EC 61131 is the standard for PLC programming and will be described in the following sections. This standard consists of five different programming languages and a common set of rules that apply to all the programming languages.

## Section 7.1: The Common IEC-Standard Hierarchy

The foundational component of the IEC 61131-3 standard is the programme organisation unit (POU). The POU is the smallest self-standing component of a PLC programme. There are three types of POUs:

1) Function (FUN): A function is a parametrisable POU without any static variables or state information (memory) that given the same input parameters will give the same output.
2) Function Block (FB): A function block is a parametrisable POU with static variables (that is, memory). Given the same input values, a function block can give different values that depend on the internal function block values, as well as external values.
3) Programme (PROG): A programme represents the "main programme." All the variables for the programme and their physical addresses must be specified. Otherwise, a programme is like a function block.

Programmes and function blocks can have input and output parameters, while functions only have input parameters and the function value as the return value. A POU consists of three parts:

1) Declaration of the POU type with POU name (and data type for functions). The possibilities are:
a. Function: FUNCTION Name DataType ... END_FUNCTION
b. Function block: FUNCTION_BLOCK Name ... END_FUNCTION_BLOCK
c. Programme: PROGRAM Name ... END_PROGRAM
2) Declaration of variables
3) Remainder of the POU with the instructions.

Before a POU can be used, each programme must be associated with a task. Before we can load a task into the PLC, we must first define:

1) Which resources do we need? In the standard, a resource is defined as either a CPU (central processing unit) or a special processor.
2) How is the program executed and with what priority?
3) Do variables have to be assigned to physical PLC addresses?
4) Do references to other programs have to be made using global or external variables?

The priority of a task shows how the given task should be executed. There two important parts to define the priority:

1) Scheduling: Scheduling focuses on how the programme is executed. There are two possibilities: cyclically, that is, the programme will be continually executed; and on demand, that is, the programme will only be executed as need.
2) Priority Type: The priority type represents whether the programme can be interrupted. Again, there are two possibilities: nonpre-emptive and pre-emptive. A nonpre-emptive task must always be completed before another task can be started. A pre-emptive task can be stopped if a task higher priority occurs. Figure 95 shows the two possibilities. As well, the priority level needs to be set. This ranges from 0 for the highest priority to 3 for the lowest priority.
Defining the above components creates a configuration. Figure 94 shows a visual representation of the components of a configuration and how these are combined together.

The call hierarchy is defined as follows: a programme can only call function blocks and functions. A function block can only call other function blocks and functions, while a function can only call other functions. Recursion cannot be implemented in this standard. POUs cannot call themselves nor can they be called as a result of a chain of POUs.


Figure 94: Visual representation of a configuration

$T_{1}$ : High priority, every 100 ms
$T_{2}$ : Low priority, nonpre-emptive, every 200 ms
$T_{3}$ : Low priority, pre-emptive, every 200 ms

Figure 95: Pre-emptive and nonpre-emptive tasks

## Section 7.2: Types of Variables

In the IEC standard, there are many different types of variables:

1) Variables (VAR): These are general variables that can be used by all POUs.
2) Input Variables (VAR_INPUT): The actual parameter will be transferred by value to the POU, that is, the variable itself is not passed to the POU, but only a copy. This ensures that
the input variable outside of the POU cannot be changed. This concept is also often called call by value. All POUs can use this type.
3) Output Variables (VAR_OUTPUT): The output variable will be returned to the calling POU as a value. This concept is also called return by value. All POUs can use this type.
4) Input-and-Output Variables (VAR_IN_OUT): The actual parameter will be transferred to the called POU as a pointer, that is, the location of the variable will be given so that any changes made to the variable are directly stored. The concept is also called call by reference. For variables with complex data structures, this can lead to efficient programming. However, since the variable location is passed, this means that (undesired) changes will impact the variable even outside the calling function. All POUs can use this type.
5) External Variables (VAR_EXTERNAL): This variable can be changed outside of the POU using the variable. An external variable is required to have read and write access to a global variable of another POU within a given POU. It is only visible for POUs that list this global variable under VAR_EXTERNAL, all others have no access to this global variable. The identifier and type of a variable under VAR_EXTERNAL must match the corresponding VAR_GLOBAL declaration in the program. Only programs and function blocks can use this type.
6) Global Variables (VAR_GLOBAL): A globally declared variable can be read and written to by several POUs. To do this, the variable must be specified in the other POUs under the VAR_EXTERNAL with the same name and type.
7) Access Variables (VAR_ACCESS): Access variables are global variables for the configurations that act as a communication channel between components (resources) of the configurations. It can be used like a global variable within the POU.

## Section 7.3: Variables, Data Types, and Other Common Elements

The IEC standard defines common elements that apply to all programmes. These elements are not only components, but also rules that determine how the elements can be used.

## Section 7.3.1: Simple Elements

Each PLC programme consists of basic elements that are defined as the smallest unit that, when combined together, build declarations and instructions, which form a complete programme. These simple elements can be classified into delimiters, keywords, literals, and identifiers.

## Section 7.3.1.1: Delimiters

Delimiters are symbols that separate the individual components from one another. Typical delimiters include the space, + , the comma (, ), and *. Table 31 shows all the delimiters in the IEC standard.

Table 31: Delimiters in the IEC IEC 61131-3 standard

| Delimiter | Meaning, Clarification |
| :--- | :--- |
| Space | Can be inserted anywhere, except within keywords, literals, <br> identifiers, directly represented variables, or combinations <br> of delimiters, such as (* or *). IEC 61131-3 does not <br> specifically make any statements about tabs and, hence, they <br> are usually treated as spaces. |
| End-of-Line | At the end of an instruction line in instruction list, in <br> structured text also permitted within an instruction. Not <br> permitted within comments in instruction list. |
| Start of Comment (* | Starts a comment (not nestable) |
| End of Comment *) | Ends a comment |
| Plus + | 1. Leading sign of a decimal literal <br> 2. In the exponent of a floating-point literal |
| Minus - | 3. Addition operator in expressions |
| Octothorpe \# | Leading sign of a decimal literal <br> 2. In the exponent of a floating-point literal <br> 3. Negation operator in expressions |
| Period . | 4. Year-month-day separator in time literals |
| 1. Base-number separator in literals |  |
| 2. Time-literal separator |  |


| Delimiter | Meaning, Clarification |
| :---: | :---: |
|  | 4. Separator for components of a FB instance (when accessing it) |
| e, E | Leading character for exponents of floating-point literals |
| Quotation Mark ' | Start and end of strings |
| Dollar sign \$ | Start of a special symbol in a string |
| ```Prefix for Time literal t#,T#; d#, D#; d, D; h, H; m, M; s, S; ms, MS; date#, DATE#; time#, TIME#; time_of_day#; TIME_OF_DAY#; tod#, TOD#; date_and_time#; DATE_AND_TIME#; dt#, DT#``` | Introductory characters for time literals, combinations of lowercase and uppercase letters are permitted |
| Colon : | Separator for: <br> 1. Time within time literals <br> 2. Definition of the data type for variable declaration <br> 3. Definition of a data-type name <br> 4. Step names <br> 5. PROGRAM ... WITH ... <br> 6. Function name / data type <br> 7. Access path: data / type <br> 8. Jump label before the next instruction <br> 9. Network name before the next instruction |
| Assignment (Walrus) <br> Operator : = | 1. Operator for initial value assignment <br> 2. Input connection operator (assignment of actual parameters to formal parameters when calling the POU) <br> 3. Assignment operator |
| (Round) Brackets (...) | Start and end of: <br> 1. Initial value list, also: multiple initial values (with repetition number) <br> 2. Range specification <br> 3. Field index <br> 4. Sequence length |


| Delimiter | Meaning, Clarification |
| :---: | :---: |
|  | 5. Operator in instruction list (calculation level) <br> 6. Parameter list when calling the POU <br> 7. Subexpression hierarchy |
| Square Brackets [...] | Start and end of: <br> 1. Array index (access to an array) <br> 2. String length (when declaring a string) |
| Comma , | Separator for: <br> 1. Lists <br> 2. Initial-value lists <br> 3. Array indices <br> 4. Variable names (when there are multiple variables with the same data type) <br> 5. Parameter list when calling a POU <br> 6. Operator in instruction list <br> 7. CASE list |
| Semicolon ; | End of: <br> 1. Definition of a (data) type <br> 2. Declaration (of a variable) <br> 3. Structured text command |
| Period-Period | Separator for: <br> 1. Range specifications <br> 2. CASE branches |
| Percent \% | Introductory character for hierarchical addresses for directly represented and symbolic variables |
| Assignment Operator => | Output binding operator (assignment of formal parameters to actual parameter when calling a PROGRAM) |
| $\begin{aligned} & \text { Comparison }>,<; \quad>=, \quad<=; \\ & =,<> \end{aligned}$ | Comparison operators in expressions |
| Exponent ** | An operator in expressions |
| Multiplication * | Multiplication in expressions |
| Division / | Division in expressions |
| Ampersand \& | AND operator in expressions |

## Section 7.3.1.2: Keywords

In the IEC 61131-3 standard, keywords are the elementary "word". Normally, keywords are written in bold. They are standard identifiers that are clearly specified in the IEC 61131-3 standard in terms of spelling and purpose. Therefore, they cannot be used in a user-defined manner for variable or other names. Capitalisation is not significant for keywords, i.e. they can be written
using all lowercase, all uppercase, or a mixture of the two. In this book, all keywords are shown in capitals. The keywords include the following possibilities:

1) Names of elementary data types
2) Names of standard functions
3) Names of standard function blocks
4) Names of the input variables for the standard functions
5) Names of the input and output variables of the standard function blocks
6) The variables EN und ENO in the graphical programming languages
7) The operators of the instruction list language
8) The elements of the structure text language and
9) The elements of sequential charts

Table 32 shows all the keywords in the IEC standard.

Table 32: All Keywords in the IEC standard

| A |  |  |
| :---: | :---: | :---: |
| ABS <br> ACOS <br> ACTION <br> ADD <br> AND <br> ANDN | ```ANY ANY_BIT ANY_DATE ANY_DERIVED ANY_ELEMENTARY ANY_INT``` | ANY_MAGNITUDE ANY_NUM ANY_REAL ARRAY ASIN AT ATAN |
| B |  |  |
| BOOL | BY | BYTE |
| C |  |  |
| CAL <br> CALC <br> CALCN <br> CASE <br> CD <br> CDT | CLK CONCAT CONFIGURATION CONSTANT COS CTD | CTU CTUD CU CV |
| D |  |  |


| D |  |  |
| :---: | :---: | :---: |
|  | DIV | DWORD |
| DATE_AND_TIME | po |  |
| DELETE | DS |  |
| E |  |  |
|  | END_PROGRAM END_REPEAT END_RESOURCE END_STEP END_STRUCT END_TRANSITION END_TYPE END_VAR END_WHILE | EN ENO EQ $E T$ EXIT EXP EXPT |
| F |  |  |
| $\begin{array}{\|l} \hline \text { FALSE } \\ \text { F_EDGE } \\ \text { F_TRIG } \end{array}$ | $\begin{aligned} & \text { FIND } \\ & \text { FOR } \\ & \text { FROM } \end{aligned}$ | FUNCTION <br> FUNCTION_BLOCK |
| G |  |  |
| GE | GT |  |
| I |  |  |
| $\begin{aligned} & \text { IF } \\ & \text { IN } \end{aligned}$ | $\begin{aligned} & \text { INITIAL_STEP } \\ & \text { INSERT } \end{aligned}$ | INT |
| J |  |  |
| JMP | JMPC | JMPCN |
| L |  |  |
| $\begin{array}{\|l} \hline \hline \mathrm{L} \\ \mathrm{LD} \\ \mathrm{LDN} \\ \mathrm{LE} \end{array}$ | LEFT | $\begin{aligned} & \mathrm{LN} \\ & \text { LOG } \\ & \text { LREAL } \\ & \text { LT } \end{aligned}$ |


| LWORD |  |  |
| :---: | :---: | :---: |
| M |  |  |
| MAX <br> MID <br> MIN | MOD <br> MOVE <br> MUL | MUX |
| N |  |  |
| $\begin{aligned} & \mathrm{N} \\ & \mathrm{NE} \end{aligned}$ | NEG NON_RETAIN | NOT |
| 0 |  |  |
| $\begin{aligned} & \mathrm{OF} \\ & \mathrm{ON} \end{aligned}$ | OR ORN |  |
| P |  |  |
| P <br> PRIORITY | $\begin{aligned} & \text { PROGRAM } \\ & \text { PT } \end{aligned}$ | PV |
| Q |  |  |
| $\begin{aligned} & \hline Q \\ & Q 1 \end{aligned}$ | $2 \mathrm{Lu}$ |  |
| R |  |  |
| $\begin{aligned} & \hline \text { R } \\ & \text { R1 } \\ & \text { R_EDGE } \\ & \text { R_TRIG } \\ & \text { READ_ONLY } \\ & \text { READ_WRITE } \\ & \text { REAL } \end{aligned}$ | RELEASE <br> REPEAT <br> REPLACE <br> RESOURCE <br> RET <br> RETAIN <br> RETC | RETCN <br> RETURN <br> RIGHT <br> ROL <br> ROR <br> RS |
|  | S |  |
| $\begin{aligned} & S \\ & S 1 \\ & S D \\ & S E L \end{aligned}$ | SEMA | SINGLE SINT SL SQRT |


| $\begin{aligned} & \text { SR } \\ & \text { ST } \\ & \text { STEP } \end{aligned}$ | $\left\lvert\, \begin{aligned} & \text { STN } \\ & \text { STRING } \\ & \text { STRUCT } \end{aligned}\right.$ | SUB |
| :---: | :---: | :---: |
| T |  |  |
| T <br> TAN <br> TASK <br> THEN <br> TIME | FIME_OF_DAY | PP PRANSITION PRUE TYPE |
| U |  |  |
| $\begin{aligned} & \hline \text { UDINT } \\ & \text { UINT } \end{aligned}$ | JLINT | USINT |
| V |  |  |
| $\begin{array}{\|l} \hline \text { VAR } \\ \text { VAR_ACCESS } \\ \text { VAR_CONFIG } \end{array}$ | $\begin{array}{\|l} \text { NAR_EXTERNAL } \\ \text { NAR_GLOBAL } \\ \text { NAR_INPUT } \end{array}$ | $\begin{array}{\|l} \text { VAR_IN_OUT } \\ \text { VAR_OUTPUT } \\ \text { VAR_TEMP } \end{array}$ |
| W |  |  |
| $\begin{aligned} & \text { WHILE } \\ & \text { WITH } \end{aligned}$ | TNORD <br> WSTRING |  |
| X |  |  |
| XOR | KORN |  |

## Section 7.3.1.3: Literals

Literals are used to represent the value of a variable (constant). They depend on the data types of these variables. A distinction is made between the following three basic types:

1) Numerical Literals, that give the numeric value of a number as a bit sequence, as well as integer and floating-point numbers.
2) String Literals, that give the value of a string in either single- or double-byte representation. A string literal is delimited by single straight quotation marks ( $\quad ; \mathrm{U}+0027$ ), for example, '' is the empty string literal and 'Automation Engineering!'. If
we wish to use a reserved symbol in a string literal, we must place a dollar sign before the reserved symbol, for example, ' $\$ \$ 45$ ' will give " $\$ 45$ ". As well, there are various nonprintable special characters that can be represented using the dollar sign. Table 33 lists some of the more common special characters.
3) Time Literals, that give the value for time points, durations, and dates.

Table 33: Special Strings

| Dollar-Sign Representation | Representation on the Screen or Printer |
| :---: | :---: |
| $\begin{gathered} \$ \mathrm{nn} \\ \text { \$ } \$ \\ \text { \$', } \quad \text { " } \\ \$ \mathrm{~L}, \quad \$ 1 \\ \$ \mathrm{~N}, \quad \$ \mathrm{n} \\ \$ \mathrm{P}, \quad \mathrm{p} \\ \$ \mathrm{R}, \quad \$ \mathrm{r} \\ \$ \mathrm{~T}, \quad \$ \mathrm{t} \end{gathered}$ | Shows "nn" in hexadecimal in ASCII \$ <br> Line feed (\$0A) <br> New line <br> New page <br> Carriage return ${ }^{20}$ (\$0D) <br> Tab |

The octothorpe is used to give additional information about the literal, for example, 2\# implies a binary representation. The additional information always comes before the octothorpe. Common representations are:

1) Binary Representation: 2\#
2) Hexadecimal Representation: 16\#
3) Duration Representation: T\# or TIME \#
4) Date Representation: D\# or DATE\#
5) Time-of-Day Representation: TOD\# or TIME_OF_DAY\#
6) Date-and-Time Representation: DT\# or DATE_AND_TIME\#
[^15]It is possible to use any defined data structure with the octothorpe. Numerical and time literals may also contain underscores in order to make the presentation more legible. Capitalisation is not important.

Time literals have some special properties. There are multiple different types of time literals: duration, date, time of day, and date and time. For each case, there is a special representation with its own rules.

Duration is represented by T \#. After the octothorpe, the duration is given using the following units:

1) d: Day
2) h: Hours
3) m: Minutes
4) $s$ : Seconds
5) ms: Milliseconds

Each unit is separated by an underscore. ${ }^{21}$ The units must be placed from largest to smallest. The smallest unit can have a decimal value, for example, T\#1m_10s_100.7ms. The highest value can "overflow", for example, the time duration $\mathrm{T} \# 127 \mathrm{~m} \_19 \mathrm{~s}$ is valid and will be automatically converted into the proper representation of $\mathrm{T} \# 2 \mathrm{~h} \mathrm{~m}^{2} 7 \mathrm{~m}_{-} 19 \mathrm{~s}$. A negative value is also possible, for example, T\#-22s_150ms.

Dates are represented using D\#. After the octothorpe comes the date in scientific notation, that is, year-month-day, for example, D\#2017-02-28.

The time of day is represented using TOD\#. After the octothorpe comes the time in the format Hours:Minutes:Seconds.Decimal Part, for example, TOD\#12: 45:25.21. Note that the 24 -hour clock is used.

The date and time is shown using DT\#. After the octothorpe comes the date in scientific notation followed by a dash and the time in the time-of-day format, for example, DT\#2017-05-30-2:30:12.

[^16]
## Section 7.3.1.4: Identifiers

Identifiers are alphanumeric strings that allow the PLC programmer to give individual names to the variables, programmes, and related elements. These include jump and network names, configurations, resources, tasks, runtime programmes, functions, function blocks, access paths, variables, derived data types, structures, transitions, steps, and action blocks. The identifiers must satisfy the following rules:

1) The first element cannot be a number ( $X$ 1Prog).
2) No more than one underscore can be used together ( $X$ A __B [with two underscores]).
3) No delimiters can be used ( $X$ w $34 \$ 23$ ).

Only the first six characters are considered when comparing two identifiers, that is, both TUI_123 and TUI_125 are equivalent. Capitalisation also plays no role, that is, TUI, tui, and TuI are all equivalent.

## Section 7.3.2: Variables

A variable is a representation of a physical memory location on a PLC to which a meaning has been assigned. The variable declaration block is bracketed with VAR_type and VAR_END. It is possible to specify the type of variable. Each variable is declared on its own line with the following key components:

```
Variable_name : Data_type := Initial_value;
```

The components in bold must always be given, while the values of components in cursive are specified by the programmer. The initial value need not be given, but it is always better to do so.

## Section 7.3.3: Data Types

There are two data types: elementary and derived data types.

## Section 7.3.3.1: Elementary Data Type

An elementary data type is defined as a simple data type that is predefined in the IEC standard. The elementary data types are characterised by their data size (number of bits) and the data range. Both values are defined by the standard, except for dates, times, and strings, whose data size and range depend on the implementation. Table 34 shows the elementary data types with their properties (data range and default initial value). The IEC 61131-3 standard defines five
groups of elementary data types that can be referenced by the given general data type given in brackets:

1) Bit Sequence and Boolean (ANY_BIT)
2) Signed and Unsigned Integers (ANY_INT)
3) Floating-Point Numbers (ANY_REAL)
4) Dates and Times (ANY_DATE)
5) Strings and Durations (ANYSTRING, TIME)

The general data types ANY_INT and ANY_REAL can be represented together by the group name ANY_NUM.

Table 34: The elementary data types in the IEC 61131-3 Standard. The initial letters in the data types represent: $\mathrm{D}=$ double, $\mathrm{L}=$ long, $\mathrm{S}=$ short, and $\mathrm{U}=$ unsigned.

| Data Type | Keyword | Bits | Range | Initial Value |
| :---: | :---: | :---: | :--- | :---: |
| BOOL | Boolean | 1 | $\{0,1\}$ | 0 |
| BYTE | Bit sequence 8 | 8 | $[0,16 \#$ FF $]$ | 0 |
| WORD | Bit sequence 16 | 16 | $[0,16 \#$ FFFF $]$ | 0 |
| DWORD | Bit sequence 32 | 32 | $[0,16 \#$ FFFF FFFF $]$ | 0 |
| LWORD | Bit sequence 64 | 64 | $[0,16 \#$ FFFF FFFF FFFF FFFF $]$ | 0 |
| SINT | Short Integer | 8 | $[-128,+127]$ | 0 |
| INT | Integer | 16 | $[-32768,+32767]$ | 0 |
| DINT | Double Integer | 32 | $\left[-2^{31},+2^{31}-1\right]$ | 0 |
| LINT | Long Integer | 64 | $\left[-2^{63},+2^{63}-1\right]$ | 0 |
| USINT | Short Integer | 8 | $[0,+255]$ | 0 |
| UINT | Integer | 16 | $[0,+65535]$ | 0 |
| UDINT | Double Integer | 32 | $\left[0,+2^{32}-1\right]$ | 0 |
| ULINT | Long Integer | 64 | $\left[0,+2^{64}-1\right]$ | 0 |
| REAL | Floating-Point | 8 | see IEC 60559 | 0 |
| LREAL | Long Floating- | 16 | see IEC 60559 | d\#0001-01-01 |
|  | Point |  |  | tod\#00:00:00 |
| DATE | Date | - | - | - |
| TOD | Time of Day | - | - | 0 |


| Data Type | Keyword | Bits | Range | Initial Value |
| :---: | :---: | :---: | :---: | :---: |
| DT | Date with Time |  |  | $\mathrm{dt} \mathrm{\# 0001-01-01-}$ |
|  | of Day |  |  | $00: 00: 00$ |
| TIME | Duration | - | - | $\mathrm{t} \# 0 \mathrm{~s}$ |
| STRING | (Single) String | - | - | $"$ |
| WSTRING | Double (String) | - | - | $" "$ |

For an elementary data type, it is possible to define the initial value and range. The initial value is defined as the value of the variable when it is first used. The range defines what the possible values for the variable are.

## Section 7.3.3.2: Arrays

Arrays are data elements of identical type that are sequentially stored in memory. An array element can be accessed with the help of an array index that lies within the array boundaries. The value of the index indicates which array element is to be accessed. Most PLC systems ensure that array access with an array index outside the array limits results in an error message during runtime. The array is defined using square brackets ([]). The dimensions are separated by commas, e.g.
ARRAY [1...45] OF INT
is a one-dimensional array with 45 elements of data type INT, while
ARRAY [1...50,1...200] OF INT
is a two-dimensional array of data type INT with 50 elements in one dimension and 200 elements in the other dimension.

The elements of an array are accessed using square brackets, for example, TEST [3] takes the third element in the array TEST. The dimensions are separated using commas.

The initial values in the array can be defined using square brackets. When values are repeated, then we can use the format Repeats (Values), to simplify matters, for example, 2 (4) means that we will write the value 4 twice. Thus, the following two array definitions are equivalent:

```
TEST1 : ARRAY [1..5] OF INT := [1, 1, 1, 3, 3];
    TEST1 : ARRAY[1..5] OF INT := [3(1), 2(3)];
```


## Section 7.3.3.3: Data Structures

Using the keywords STRUCT and END_STRUCT, it is possible to define new hierarchical data structures that contain arbitrary elementary or other already defined derived data types as subelements. If a subelement is in turn a structure, a hierarchy of structures is created, for which the lowest structure level is formed of elementary or derived data types.

As in many other programming languages, the components of a data structure are accessed using a period and the component name, for example, VAR.TEST[3], where the component TEST is an array.

## Section 7.3.3.4: Derived Data Types

A derived data type is a user-defined data type that consists of elementary data types, arrays, and data structures. This procedure is called derivation or type definition. In this way, a programmer can define the best data model for the problem at hand. A derived data type is defined by TYPE and END_TYPE. The initial values for the elements of a derived data type is given by $:=$ for example,

```
TYPE
    COLOUR : (red, yellow, green);
    SENSOR : INT;
    MOTOR :
    STRUCT
        REVOLUTIONS : INT := 0;
        LEVEL : REAL := 0;
        MAX : BOOL := FALSE;
        FAILURE : BOOL := FALSE;
        BRAKE : BYTE := 16#FF;
    END_STRUCT;
END_TYPE;
```

where the element COLOUR is an enumeration that can only take the values red, green, and yellow; SENSOR is an INT; MOTOR is a data structure that contains the following elements: REVOLUTION with an initial value of 0 , LEVEL with an initial value of 0 , MAX with an initial
value of FALSE, FAILURE with an initial value of FALSE, and BRAKE with a hexadecimal initial value of FF .

## Section 7.4: Ladder Logic (LL)

The programming language ladder logic (LL) comes from the domain of electromechanical relay systems and describes the flow of electricity through a single network representing the POU. This programming language is primarily used for working with Boolean signals.

The ladder network or diagram consists of two vertical tracks and horizontal rungs connecting the vertical tracks. It is assumed that "electricity" flows from the left-hand track to the right-hand track following the rungs. Thus, the ladder network is always read rung by rung from top to bottom and in a given row from left to right, as long as no other order is provided. It is traditional to label all the left-hand rungs with a number. Normally, the numbers are not sequential but increase in units of 5 or 10 in order that additional future rungs can be easily added. Finally, the ladder network is often visually split into two parts: a left-hand part that shows the computations and a right-hand part that shows the storing or using of the variables. This convention makes reading the ladder network easier.

## Section 7.4.1: Components of Ladder Logic

Table 35 shows the components of ladder logic.

Table 35: Components of Ladder Logic

| Name | Commbolary |  |
| :---: | :---: | :--- |
| Rung | Read from left to right <br> Copies the value from left to right, when the <br> Closed contact | value of the variable VarName is TRUE; <br> otherwise, FALSE is copied. <br> Copies the value from left to right, when the <br> value of the variable VarName is FALSE; |
| otherwise, TRUE is copied. |  |  |


| Name | Symbol | Commentary |
| :---: | :---: | :---: |
| Positive-transition sensing contact | $\frac{\text { VarName }}{\substack{\text { Vam }}}$ | Copies the value from left to right if and only if a FALSE $\rightarrow$ TRUE transition in the variable VarName is detected; otherwise, FALSE is copied. |
| Negative-transition sensing contact | $\begin{aligned} & \text { VarName } \\ & -\|N\| \end{aligned}$ | Copies the value from left to right if and only if a TRUE $\rightarrow$ FALSE transition in the variable VarName is detected; otherwise, FALSE is copied. |
| Coil ${ }^{22}$ | $\begin{aligned} & \text { VarName } \\ & {[]} \end{aligned}$ | Copies the value on the left into the variable VarName. |
| Negated coil ${ }^{22}$ | $\begin{aligned} & \text { VarName } \\ & {[/]} \end{aligned}$ | Copies the negated value of the left into the variable VarName. |
| Set coil ${ }^{22}$ | $\begin{gathered} \text { VarName } \\ {[\mathrm{S}]} \end{gathered}$ | Copies TRUE to variable VarName, if the left link is TRUE; otherwise, no change. |
| Reset coil ${ }^{22}$ | $\begin{gathered} \text { VarName } \\ {[R]} \end{gathered}$ | Copies FALSE to variable VarName, if the left link is TRUE; otherwise, no change. |
| Positive-transition sensing coil ${ }^{22}$ | $\begin{aligned} & \text { VarName } \\ & \hline \end{aligned}$ | Saves TRUE to the variable VarName if and only if a FALSE $\rightarrow$ TRUE transition is detected on the left link; otherwise, no change in the variable. |
| Negative-transition sensing coil ${ }^{22}$ | $\begin{gathered} \text { VarName } \\ {[\mathrm{N}]} \end{gathered}$ | Saves TRUE to the variable VarName if and only if a TRUE $\rightarrow$ FALSE transition is detected on the left link; otherwise, no change in the variable. |

[^17]| Name | Symbol | Commentary |
| :---: | :---: | :---: |
| Set－Reset Block | $S^{\text {SR }}$ 21 <br> $R$  | This block combines the functions of the set and reset coils． |
| Return | －〈RETURN〉 | Exits the POU and returns to the calling POU． |
| Conditional return | －tnw－〈RETURN〉 | If the left link $t n w^{23}$ is TRUE，exit the POU und return to the calling POU；otherwise，no meaning． |
| Jump | $\longmapsto$ NAME | Jump to the network with the given NAME． |
| Conditional jump | tnw $\rightarrow$ NAME | If the left link $\mathrm{tnw}^{23}$ is TRUE，jump to the network with the given NAME． |
| Label | $\stackrel{1}{\text { LABEL }}^{\text {l }}$ | Shows the name for part of a network． |

Figure 96 shows how the typical Boolean operators can be implemented in ladder logic． The AND operator is implemented by placing the two contacts in series，while the OR operator places the two contacts in parallel．This follows from the observation that the electricity flows from left to right．For an AND operation，we need both contacts to be true for the electricity to flow． This implies that both need to be in series．On the other hand，for the OR operation，electricity can flow through either of the two paths．Therefore，the contacts should be placed in parallel．

[^18]

Figure 96: (top) AND and (bottom) OR in ladder logic

## Section 7.4.2: Functions and Ladder Logic

Since ladder logic was originally developed for logical (or Boolean) systems, the implementation and running of complex functions using the simple ladder logic components can be difficult. For this reason, it is possible to define a function block that is programmed using another programming language. This function block always contains two Boolean variables (EN and ENO), as well as all the other required parameters. The Boolean input variable EN (enable in) determines if the function will be called. If EN is TRUE, then the function is called. The Boolean output variable ENO (enable out) determines if the programme completed successfully. It takes the value TRUE, if no errors occurred. Figure 97 shows such an implementation using ladder logic.


Figure 97: Calling a function in ladder logic

## Section 7.4.3: Examples of Using Ladder Logic

Example 41: Ladder Logic for a Boolean Function
Please write the corresponding ladder diagram for the following Boolean function:

$$
Q=X Y+X Z+Y Z
$$

## Solution

When we want to create the ladder logic diagram, we should first convert the Boolean function into a minimal form. In our case, we can easily convert it to

$$
Q=X Y+(X+Y) Z
$$

In the ladder logic diagram, we will require (at least) one row for each term that is separated by a " + ", since each row corresponds to an AND Term. Thus, in our example, we will require 3 rows. All the values are not negated. Thus, the open contact will be used for all values. In order to save the resulting function value in $Q$, we require a coil. Figure 98 shows the ladder logic diagram for this example.


Figure 98: Ladder Logic for Example 41

## Example 42: Ladder Logic for a Recipe

Consider a tank that needs to be filled with two components and then mixed before being sent on its merry way. The recipe is:

1. Once the start button is pressed and both level sensors (L1 and L2) read FALSE, turn on Valves 1 and 2 (V1 and V2). Go to Step 2.
2. Once the fluid reach level sensor 2 (L2), that is, it reads TRUE, close both valves and turn on the mixer (M1) for five minutes. Go to Step 3.
3. Turn off the mixer and wait one minute. Go to Step 4.
4. Open the bottom valve (V3) and let the fluid drain. Go to Step 5.
5. Once the bottom level sensor 1 (L1) reads FALSE, that is, the tank is empty, close the bottom value. Go to Step 1.

Implement this recipe using ladder logic.

## Solution

In order to implement a recipe in ladder logic, we will need to define Boolean variables that keep track of in which step we are. In this example, since we have five steps, it makes sense to define the Boolean step variables as $S 1, S 2, S 3, S 4$, and $S 5 . S 1$ will be initialised as TRUE, while all other variables will be initialised as FALSE. At the end of each step, the current step variable will be reset (to zero) and the next step variable will be set (to one). Turning on and off valves will be implemented using the set and reset coils, while the delay and timer will be implemented using the built-in time function (TON). Rung numbers have been added in multiples of 5. Finally, it can be noted that in this particular recipe each set has a corresponding reset. This is a feature of many recipes that will be made explicit in the PLC programming language sequential function charts. The final programme in ladder logic is shown in Figure 99.


Figure 99: Ladder Logic for Example 42

## Section 7.4.4: Comments

Ladder logic is perfect for large Boolean networks. However, it is not appropriate for sequence or recipe programming. In ladder logic, it is not recommended to use jumps, since these can lead to inconsistencies during runtime.

## Section 7.5: Instruction List (IL)

Instruction list (IL) is a language similar to assembly language. This language is a simple list of instructions, each of which is placed on a separate line. Each instruction has the following form:

Label: Operator/Function Operand(list) Comments
where Label names a given row of instructions and along with the colon can be ignored; the Operator or Function is an operator in instruction list or a function, Operand is zero, one, or more constants or variables for the operator or input parameters for the functions that are separated by commas; and Comments is an optional field that provides additional information about what the given instruction row does. All fields in italics are optional.

## Section 7.5.1: Universal Accumulator

Most assembly languages have a physical accumulator in the processor, that is, a value is loaded into this accumulator, additional values are added, subtracted and then stored. Instruction list also has such an accumulator that is often called the "current result." However, it is not a memory area with permanently defined register lengths. Instead, the statement list compiler ensures that an abstract accumulator of any memory width is available that depends on the data type to be processed. In contrast to other assembly languages, there are no special processor status bits. The result of comparisons is written as a Boolean 0 or 1 in this accumulator. Conditional jumps and calls use the current value stored in the accumulator to evaluate the jump requirements. The current result can be of any general data type, derived data type or function-block type. The data width of the current result (number of bits) does not matter. Instruction list requires that two consecutive operations be type compatible, which means that the data type of the current result must be compatible with the following statement.

Table 36 shows the modifications of the current result by different operator groups. Unchanged means that an instruction passes the result of the previous instruction on to the following one without having made any changes to the value and type. Undefined means that the subsequent operation cannot use the current result. The first command of a function block (which is called using CAL) may only be a load LD, jump JMP, function block call CAL or return RET, since these commands do not require a current result.

The IEC 61131-3 standard itself does not define any operator groups. In fact, the behaviour and evaluation of the current result is only partially described in the standard. With operations such as AND, the type and content of the current result is intuitively clear both before and after execution. However, the standard leaves open the question of how the current result is defined after an unconditional jump.

| Impact of the Operator Group <br> on the Current Result | Abbreviation | Example Operators |
| :--- | :--- | :--- |
| Create | C | LD |
| Process | P | GT |
| Unchanged | U | ST ; JMPC |
|  |  | Unconditional CAL block call, since the <br> Undefined |
|  | - | following instruction must reload the current |
|  |  | result, because it does not have a clearly defined |

## Section 7.5.2: Operators

Table 37 shows the defined operators in instruction list. The modifiers are defined as follows:

- N: Negation,
- (: Nesting levels using brackets,
- C: Conditional performance of an operator when the current result equals TRUE.

The modifiers are written together with the operator, for example, ANDN is a negated AND.

Table 37: Operators in Instruction List

| Operator | Acceptable Modifiers | Operand | Definition |
| :---: | :---: | :---: | :--- |
| LD | N | ANY | Load |
| ST | N | ANY | Save |
| S |  | BOOL | Set |
| R |  | BOOL | Reset |
| AND / \& | $\mathrm{N},($ | ANY | Boolean AND |
| OR | $\mathrm{N},($ | ANY | Boolean OR |
| XOR | $\mathrm{N},($ | ANY | Boolean Exclusive OR |
| ADD | $($ | ANY | Addition |
| SUB | $($ | ANY | Subtraction |
| MUL | $($ | ANY | Multiplication |
| DIV | $($ | ANY | Division |
| GT |  | ANY | $>$ |


| Operator | Acceptable Modifiers | Operand | Definition |
| :---: | :---: | :---: | :--- |
| GE | $($ | ANY | $\geq$ |
| EQ | $($ | ANY | $=$ |
| NE | $($ | ANY | $\neq$ |
| LT | $($ | ANY | $<$ |
| LE | $($ | ANY | $\leq$ |
| JMP | C, ( | LABEL | Jump to LABEL |
| CAL | C, ( | NAME | Call function NAME. |
| RET | C, ( |  | Return from a function call |
| $)$ |  |  |  |
|  |  |  | Take the last deferred instruction |

The current result can be linked with the result of an instruction sequence using the bracket operators. When the modifier " (" appears, the associated operator, the value of the current result and the data type of the current result are cached. The data type and value of the following line are loaded into the current result. When the ")" operator appears, the cached values and data type based on the operator and modifiers are linked with the current result. The result is then stored as the current result. Expressions in brackets can be nested.

## Section 7.5.3: Functions in Instruction List

In instruction list, functions are called by giving their name. The parameters can be passed using two different approaches: actual parameters or formal parameters. With actual parameters, the first parameter of a function is the current result that was previously loaded. Therefore, the first operand after the function name is the second function parameter in this variant. With formal parameters, all parameters are explicitly defined. Table 38 shows the two possibilities.

Table 38: Two Possibilities for Calling the Function LIMIT (MN, IN, MX)

| Possibility 1: Actual Parameters |  |
| :--- | :---: |
| LD | Possibility 2: Formal Parameters |
| LIMIT | 2,3 |
|  | LIMIT ( |
|  | MN $:=1$, |
|  | IN $:=2$, |
|  | MX $:=3$ |
|  |  |
|  |  |

A function has at least one output parameter (function value) that is returned using the current result. Should the function have additional output parameters, these can be returned using parameter assignments. A call without formal parameters takes place in one line. The original order of the output declarations must be observed. In the case of formal parameter assignments, this is done line by line, followed by bracket line. With formal parameter assignment, the output parameters are marked with $=>$, e.g., ENO $=>$ ErrorOut means that ENO is an output value and is saved in the ErrorOut variable. The programming system assigns the function value to a variable with the function name. This name is declared automatically and does not have to be specified separately by the user in the declaration section of the calling block.

## Section 7.5.4: Calling Function Blocks in Instruction List

Function blocks are called using the CAL operator. The IEC 61131-3 standard provides three methods for calling function blocks:

1) Calling a function block with a bracketed list of input and output parameters,
2) Calling a function block with previously loaded and saved input parameters, and
3) Calling a function block implicitly by using the inputs as operators.

Method 3 can only be used for standard function blocks, since the programming system must use the name of the function block inputs as operators. This method cannot be used for the output parameters. Table 39 shows the three methods for calling the function block ZEIT1 (IN, PT) with output variables $Q$ and ET.

Table 39: Three Methods for Calling the Function Block ZEIT1 (IN, PT) with output variables Q and ET.

| Method 1 | Method2 | Method 3 |
| :--- | :--- | :--- |
| CAL | ZEIT1 ( | LD t\#500_ms |
| IN $:=$ Frei, | ST ZEIT1.PT | ST ZEIT1.PT |
| PT $:=$ t\#500_ms, | LD Frei | LD Frei |
| Q => Aus, | ST ZEIT1.IN | IN ZEIT1 |
| ET => WERT | CAL ZEIT1 |  |
| ) | LD ZEIT1.Q | LD ZEIT1.Q |


| Method 1 | Method 2 | Method 3 |
| :--- | :--- | :--- |
|  | ST Aus | ST Aus |
|  | LD ZEIT1.ET | LD ZEIT1.ET |
|  | ST WERT | ST WERT |

## Section 7.5.5: Examples

Example 43: Example of the Computation of the Current Result
For the given instructions, follow how instruction list uses and updates the different registers. Values in grey are the default initial values.

| Instruction | (* Comments *) | CR | X1 | X2 | W1 | FB1.i1 | FB1.01 |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  | Initial values: | "STR" | 0 | 1 | 10 | 1 | 103 |
| LDN X1 | Load the negated value in X1 | 1 | 0 | 1 | 10 | 1 | 103 |
| AND X2 | $\begin{aligned} & \mathrm{CR} \text { AND } \mathrm{X} 2 \rightarrow \\ & \mathrm{CR} \end{aligned}$ | 1 | 0 | 1 | 10 | 1 | 103 |
| S X1 | If $C R=1$ <br> then: $1 \rightarrow X 1$ | 1 | 1 | 1 | 10 | 1 | 103 |
| R X2 | If $C R=1$ <br> then: $0 \rightarrow X 2$ | 1 | 1 | 0 | 10 | 1 | 103 |
| JMPCN <br> Lab1 | Jump to Lab1 if $C R=0$ | 1 | 1 | 0 | 10 | 1 | 103 |
| LD W1 | Load the value in $W 1 \rightarrow$ $C R$ | 10 | 1 | 0 | 10 | 1 | 103 |
| MUL W1 | $\mathrm{CR} * \mathrm{Wl} \rightarrow \mathrm{CR}$ | 100 | 1 | 0 | 10 | 1 | 103 |
| SQRT | Function: <br> sqrt(CR) $\rightarrow$ CR | 10 | 1 | 0 | 10 | 1 | 103 |
| $\begin{aligned} & \text { ST } \\ & \text { FB1.il } \end{aligned}$ | $\begin{aligned} & \text { Save } \\ & \text { FB1.i1 } \end{aligned} \quad \rightarrow$ | 10 | 1 | 0 | 10 | 10 | 103 |
| CAL FB1 | FB1: internal computation | 10 | 1 | 0 | 10 | 10 | 112 |


| Instruction | (* Comments *) | CR | X1 | X2 | W1 | FB1.i1 | FB1.01 |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| $\begin{aligned} & \text { LD } \\ & \text { FB1. } 01 \end{aligned}$ | Load the <br> output o1 <br> from FB1 $\rightarrow$ CR | 112 | 1 | 0 | 10 | 10 | 112 |
| ST W1 | Save CR $\rightarrow$ W1 | 112 | 1 | 0 | 112 | 10 | 112 |
| GT 90 | $\underset{\mathrm{CR}}{(\mathrm{CR}>90) ? \rightarrow}$ | 1 | 1 | 0 | 112 | 10 | 112 |

Example 44: Writing the Instruction List Programme
Please write the instruction list programme for the following function:

$$
Q=X Y+X Z+Y Z
$$

## Solution

The simplest approach to writing the instruction list programme is to go from left to write and write the corresponding instruction. The programme is then

```
LD X
AND Y
OR( X
AND Z
)
OR( Y
AND Z
)
ST Q
```


## Section 7.5.6: Comments

Instruction list is perfect for large Boolean networks. However, like with ladder logic, it is inappropriate for sequence or recipe programming. Instruction list has no complex programming flow concepts, such as loops. Finally, in instruction list, jumps should be avoided as they can lead to inconsistencies during runtime.

## Section 7.6: Function-Block Language (FB)

The function-block language is a graphical model with Boolean operators and complex functionality represented using blocks. Motivated by electronic circuits, the signals flow between the blocks and are converted from Boolean inputs into Boolean outputs. Although this language is based on electrical circuits, the current standard has adopted different representations for the blocks symbolising AND and OR.

## Section 7.6.1: Blocks used in the Function-Block Language

Table 40 shows the blocks used in the function-block language.

Table 40: Blocks used in the function-block language

| Name | Symbol | Comments |
| :---: | :---: | :---: |
| AND |  |  |
| OR |  |  |
| Negation | 0 | Negation is placed where the signal enters or leaves a block. |
| Link | $\bullet$ | When two connections coincide and the value is to be transferred. |
| General Block | NAME | There are as many input and output ports as required by the block. |
| Rising Edge | > | This is placed within the function block next to the corresponding variable in order to show a rising edge transition. |


| Name | Symbol | Comments |
| :---: | :---: | :---: |
| Falling <br> Edge | $<$ | This is placed within the function block next to the corresponding variable in order to show a falling edge transition. |
| Return | -<Return> | Leaves a POU and returns to the calling POU. |
| Conditional Return | -snw-<RETURN〉 | If the left connection $s n w^{24}$ is TRUE, then exit the POU and return to the calling POU; otherwise ignored. |
| Jump | $\longrightarrow$ NAME | Jump to the network with the given identifier NAME. |
| Conditional Jump | $\rightarrow$ snw $\rightarrow$ NAME | If the left connection $\mathrm{Snw}^{24}$ is TRUE, jump to the network with the given identifier NAME; otherwise ignored. |

Inputs and outputs are defined upon first use. Figure 100 shows an example of a function block diagram with common components. The exact method for labelling the variables is not specified in the standard; any reasonable method can be used.

[^19]

Figure 100: Diagram using the Function-Block Language

## Section 7.6.2: Feedback Variable

Function block language allows an output to be used as the new input of a network. Such a variable is called a feedback variable. The first time such a situation is encountered the initial value is used; afterwards, the last value is used. Figure 101 shows an example with feedback.


Figure 101: Feedback in the Function-Block Language

## Section 7.6.3: Example

## Example 45: Creating the Diagram using the Function-Block Language

Please create the diagram using the function-block language for the following Boolean function:

$$
Q=X Y+X Z+Y Z
$$

## Solution

Figure 102 shows the solution.


Figure 102: The Function $Q$ in the Function-Block Language

## Section 7.6.4: Comments

Function block language should be used in cases where there is signal flow between different components, for example, in closed-loop control. This language contains no complex programme flow concepts, such as loops.

## Section 7.7: Structured Text (ST)

Structured text (ST) is a further text-based programming language in the IEC 61131-3 standard. It is referred to as a high-level programming language, since assembly-language commands are not used in structured text, but powerful constructs are built using more abstract commands. In structured text, the solution to a particular programming task is broken down into individual steps (instructions). An instruction is used to calculate and assign values, control the flow of commands, or to call or exit a POU.

## Section 7.7.1: Commands in Structured Text

Table 41 shows the commands in structured text. Each command must be separated by a semicolon, that is, unlike previous text-based languages, multiple commands can be placed on a single line.

The IF command has many different forms. The obligatory components are the IF, THEN, and END_IF commands. The IF command opens the IF block, while the END_IF closes it. The other possibilities are the ELSE and ELSIF commands. The ELSE block is only performed if all previous expressions have been evaluated as FALSE. The ELSIF block is only performed if all previous expressions have been evaluated as FALSE and the corresponding Boolean expression is TRUE. This command can be repeated as often as desired. This gives the following possibilities:

1) IF expression THEN code block; END_IF;
2) IF expression THEN code block; ELSE code block; END_IF;
3) IF expression THEN code block; ELSIF expression THEN code block; ELS IF expression THEN code block; ELSE code block; END_IF;

The FOR command also has many different possibilities. The obligatory parts are FOR, TO, DO, and END_FOR. The FOR command opens this code block and the END_FOR command closes it. The TO command specifies the end value and the DO command closes the declaration line of the FOR command. The optional part is the BY command that gives the step value. If no step value is given using the BY command, then it is assumed that the step is one. The two possibilities for the FOR command are:

1) FOR Counter $:=$ expression $T O$ expression $B Y$ expression $D O$ code block; END_FOR;
2) FOR Counter := expression $T O$ expression DO code block; END_FOR;

Table 41: Commands in Structured Text

| Keyword | Description | Example | Comments |
| :---: | :--- | :--- | :--- |
| $:=$ | Assignment | $\mathrm{d}:=10 ;$ | Assigns the value on the right to <br> the variable on the left. |
| $:=,=>$ | FBNAME ( | Calls another POU of type <br> function block and assigns the |  |
| Function Blocks | Part $:=10$, | required parameters: $:=$ for <br> RETURN | Return |


| Keyword | Description | Example | Comments |
| :---: | :---: | :---: | :---: |
| $\begin{gathered} \text { IF ... THEN } \\ \text { ELSE, } \\ \text { ELSIF } \\ \text { END_IF } \end{gathered}$ | Branching | IF A>100 THEN <br> f: =1; <br> ELSIF d=e THEN <br> f:=2; <br> ELSE <br> f:=4; <br> END_IF <br> CASE f OF | Selection of alternatives using Boolean logic. |
| $\begin{gathered} \text { CASE ... OF } \\ \text { ELSE } \\ \text { END_CASE } \end{gathered}$ | Multiple Selection | $\begin{aligned} & 1: \mathrm{f}:=11 ; \\ & 2: \mathrm{f}:=14 ; \\ & \text { ELSE f:=-1; } \\ & \text { END_CASE; } \\ & \text { FOR h:=1 TO } 10 \end{aligned}$ | Selects a code block based on the value of the expression $f$. |
| $\begin{gathered} \text { FOR ... TO ... } \\ \text { BY ... DO } \\ \text { END_FOR } \end{gathered}$ | FOR Loop | BY 2 DO <br> f[h/2]:=h; <br> END_FOR; <br> WHILE m>1 DO | Repeated running of a code block with start and end conditions. |
| $\begin{gathered} \text { WHILE ... DO } \\ \text { END_DO } \end{gathered}$ <br> RFPFAT | WHILE Loop | $\mathrm{n}:=\mathrm{n} / 2 \text {; }$ <br> END_WHILE; <br> REPEAT | Repeated running of a code block with end conditions. |
| REPEAT ... <br> UNTIL END_REPEAT | REPEAT Loop | $\begin{aligned} & i:=i * j ; \\ & \text { UNTIL i>10000 } \\ & \text { END_REPEAT; } \end{aligned}$ | Repeated running of a code block with end conditions. |
| EXIT | Breaking a loop Command delimiter | EXIT; | Immediate exit from a loop <br> Shows the end of a command. |

## Section 7.7.2: Operators in Structured Text

Table 42 shows the priority of the operators in structured text.

| Operator | Description | Priority |
| :---: | :---: | :---: |
| (...) | Brackets | Highest |
| Function (...) | Function Evaluation |  |
| ** | Exponentiation |  |
| -, NOT | Negation, Boolean Complement |  |
| *, / | Multiplication, Division |  |
| MOD | Modulo |  |
| +, - | Addition, Subtraction |  |
| $>,<, \leq, \geq$ | Comparison Operators |  |
| $=$ | Equality |  |
| <> | Inequality ( $\ddagger$ ) |  |
| AND, \& | Boolean AND |  |
| XOR | Boolean exclusive OR |  |
| OR | Boolean OR | Lowest |

## Section 7.7.3: Calling Function Blocks in Structured Text

In structured text, function blocks are called by their name with the required parameters placed in brackets. The $:=$ is used to assign the value of the actual parameters, while $=>$ is used to assign any output variables. Since the assignments must be explicit, the order of the assignments is irrelevant. If a parameter is not initialised during the call, then the initial value or the last value will be used.

## Section 7.7.4: Example

## Example 46: Structured Text

Please write the structured text programme for the following system. As shown in Figure 103, a tank can be filled using valves. The weight of the tank is determined using a scale. The function block monitors the tank weight to determine if the tank is full, empty, or in between the two states. The block gives a single command that can have four values:

1: Fill the tank;
2: Stop filling the tank;
3: Start the stirrer;
4: Empty the tank.

As necessary, the appropriate values are opened or closed to control the tank level. The stirrer can only work if the tank is full; otherwise, the command is ignored.


Figure 103: Tank System for the Structured-Text Example
Solution

```
(*Tank States*)
TYPE T_STATE:(FULL,NOT_EMPTY,EMPTY); END_TYPE;
(*Valve States*)
TYPE T_VALVE:(ON,OFF); END_TYPE;
FUNCTION_BLOCK Weightcontrol
VAR_IN
    Command: INT;
    Weight: Real;
    Full_weight,Empty_weight: Real; (*Same data type declared on
one line*)
END_VAR
VAR_OUTPUT
    V1: T_VALVE := OFF;
    V2: T_VALVE := OFF;
    Speed: REAL := 0.0;
```

```
END_VAR
VAR (*Internal Variables*)
    State: T_STATE := EMPTY;
END_VAR
(*Determine the Tank State: Compare the full and empty weights *)
IF Weight >= Full_weight THEN
    State := FULL;
ELSIF Weight <= Empty_weight THEN
    State := EMPTY;
ELSE
    State := NOT_EMPTY;
END_IF
(*Implementation of Commands: 1-Fill, 2-Stop, 3-Stir, 4-Empty*)
CASE Command OF
    1: V2 := OFF;
    V1 := SELECT(G := State=FULL,INO := ON,IN1 := OFF);
    (*ON only if G is false*)
    2: V2 := OFF;
    V1 := OFF;
    4: V1 := OFF;
    V2 := ON;
END_CASE;
(*Stirrer Speed*)
Speed := SELECT(G := Command=3; IN0 := 0.0; IN1 := 100.0);
END_FUNCTION_BLOCK
```


## Section 7.7.5: Comments

Structured text has the following advantages (especially in comparison with instruction list) are:

- Compact formulation of the programming tasks,
- Clear programme structure, and
- Powerful constructs to control the flow of commands.

However, it has the following disadvantages:

- The conversion of the programme into machine code cannot be directly influenced since it is performed using a compiler.
- The higher abstraction level brings with it a loss of efficiency, that is, the translated programme is longer and slower.


## Section 7.8: Sequential-Function-Chart Language (SFC)

The sequential-function-chart language (SFC) is the second graphical method for programming a PLC. It consists of a sequence of steps and transitions that implement predefined tasks and provide a visual overview of the process.

## Section 7.8.1: Steps and Transitions

A step is either active or inactive. It consists of a number of instructions that are performed as long as the given step is active. A transition using a Boolean expression determines when a step becomes inactive. Connections with a predefined direction describe which step or steps should be activated next.

Steps are shown using a rectangle. Figure 104 shows two possible forms for showing a step. The general step block, which is shown in Figure 104 (left), is a simple rectangle with a single border. The initial step block, which is shown in Figure 104 (right), is a simple rectangle with double border.


Figure 104: Steps in Sequential Function Charts: (left) general step and (right) initial step
The transition conditions are marked with a horizontal line with identifiers. The Boolean description, often called a guard, shows the requirements for a transition to occur. The guard can be written in any of the other PLC languages. Figure 105 shows some possibilities. The most frequently used PLC languages for describing transitions is structured text, function block language, or ladder logic. Often, the guard is ignored.



Structured Text

Function-Block Language

Ladder Logic

## Instruction List

Figure 105: Transitions Conditions in Different PLC Languages
When the POU is called, the specially marked initialisation step is made active. All the assigned instructions will be performed. When the transition conditions become TRUE, the initialisation step will be made inactive and the next step will be activated. When a transition
occurs, the active attribute (often called a token) is passed from the current active step to its successor. This active attribute wanders through the individual steps, multiplying for parallel branches, and coming together again. It may not be completely lost nor is uncontrolled distribution, which occurs when multiple tokens are present in a single step, possible.

## Section 7.8.2: Action Blocks

An action block shows the details of what must be accomplished in a given step. Each action block is always assigned to a particular step. It is not necessary to include the action block. No more than two action blocks can be associated with a given step.

Each action block has four components, which are shown in Figure 106:
a) Qualifier, that consists of an acceptable tag.
b) Action Name, that briefly describes the action.
c) Indicator Variable, that provides which PLC variable is to be used.
d) Process Description, that describes in an appropriate language the action (optional).

Each component has a specified location in the action block. The qualifier must be selected from a small group of abbreviations that describe how the step is to be performed. Table 43 shows the possible qualifiers.


Figure 106: Components of an Action Block: a: Qualifier; b: Action Name; c: Indicator Variable; d: Process Description
Table 43: Qualifiers in Sequential Function Charts

| Qualifier | Short Name | Description |
| :---: | :--- | :--- |
| N | not saved | Means that as soon as the step has ended, the action is <br> stopped. |
| S | saved | Means that the action continues with the given values <br> until it is stopped. |
| R | reset | Means that a saved action is stopped. |
| (time) length | Means that the action lasts for the specified time $T$. |  |


| Qualifier | Short Name | Description |
| :---: | :--- | :--- |
| D | delay | Means that the action is delayed by the specified time <br> before being implemented. <br> Means that the action occurs for a very short period of <br> time. |
| DS | delayed and saved | Means that the action is first delayed and then saved. <br> SD |
| saved and delayed that the action is first saved and, then after the |  |  |
| SL | Melay, if the step is still active, implemented. |  |
| saved and time | Means that a time-limited action is saved. |  |
| limited |  |  |

## Section 7.8.3: Sequential Function Charts

As the name suggests, in the sequential-function-chart language, a chart is designed to show how the individual components are linked together. This chart is built using the following rules:

1) A vertical link connects two steps.
2) An arrow is used to clarify the direction in which steps occur.
3) An action block is connected with the step block using a straight horizontal line.

There are two special types of connections: alternative and parallel paths. An alternative path is shown using a single horizontal line across all possible alternative ways. This horizontal line is placed at the beginning and end of the region corresponding to the alternative paths. Only one of the paths must end at which point all the other paths will be stopped. Figure 107 shows an alternative path. Alternative paths are normally selected from left to right. This is shown in Figure 108. In this case, the paths will be selected in order from S_2a, S_2b, and then S_2c. This means that first the transition condition for S_2a will be tested. If it is TRUE, this path will be taken and all other paths ignored. If it is FALSE, then $S \_2 b$ and $S \_2 c$ (in this order) will be tested. When a user defined order is provided, as, for example, in Figure 109, then this order is followed. In this example, it means that first $S_{-} 2 b$ will be tested, followed by $S_{-} 2 c$ and finally $S_{-} 2 a$.

Parallel paths are shown by a doubled horizontal line across all possible parallel paths. This horizontal line is placed at the beginning and end of the region corresponding to the parallel paths. All parallel paths must end before the process can continue. Figure 110 shows an example of a parallel path.


Figure 107: Alternative Paths in Sequential Function Charts


Figure 108: Usual Decision Order for Alternative Paths in Sequential Function Charts


Figure 109: User-Defined Decision Order for Alternative Paths in Sequential Function Charts


Figure 110: Parallel Paths in Sequential Function Charts

## Section 7.8.4: Example

## Example 47: Creating a Sequential Function Chart

Consider a chemical reactor as shown in Figure 111. It is desired to design a PLC to control the process. The objective is to draw the sequential function chart for the process. The desired process can be described as follows:

When the START button has been pressed, the reactor status sensor S1 confirms that the reactor is empty and that the temperature sensor S 4 and the pressure sensor S 5 give no error messages, then valve Y1 should be opened until level sensor S2 reaches the desired value, that is, it returns a 1. Then, the motor for agitator M should be turned on and valve Y2 opened. When the level sensor S 3 reaches the desired value, valve Y2 should be closed. After waiting 5 s , heater H is turned on until the temperature sensor S 4 reaches the desired value. During the heating procedure, should the pressure sensor S 5 return a high-pressure alarm in the reactor, then the pressure release valve Y 4 should be opened until the alarm is lowered. After the heating procedure has been completed, the agitator runs for 10 more seconds at which point the valve Y3 is opened. Once the reactor is empty, that is, the reactor status given by S 1 returns 1 , valve Y 3 is closed and the process can be restarted.


Figure 111: Schematic of the Reactor

## Solution

Before the solution to the problem is provided, it is useful to consider the general procedure for solving such problems. This procedure can be written as

1. Define all the variables and their values, especially for the Boolean variables.
2. Write the process description as a series of steps. Note that a single sentence can be associated with multiple steps or multiple sentences can be combined into a single step.
3. Draw the sequential function chart.

## Defining the Variables

The variables are defined as follows.

| Input Variable | Symbol | Data Type | Logical Values |  | Address |
| :---: | :---: | :---: | :---: | :---: | :---: |
| Start Button | START | BOOL | Pressed | START = 1 | E 0.0 |
| Reactor Status | S1 | BOOL | Tank empty | S1 $=1$ | E 0.1 |
| Level Sensor 1 | S2 | BOOL | Level 1 reached | $\mathrm{S} 2=1$ | E 0.2 |
| Level Sensor 2 | S3 | BOOL | Level 2 reached | S3 $=1$ | E 0.3 |
| Temperature Sensor | S4 | BOOL | Temperature reached | $\mathrm{S} 4=1$ | E 0.4 |


| Pressure Sensor | S5 | BOOL | High-Pressure <br> Alarm | S5 = 1 | E 0.5 |
| :---: | :---: | :---: | :---: | :---: | :---: |
| Valve 1 | Y1 | BOOL | Valve open | $\mathrm{Y} 1=1$ | A 4.1 |
| Valve 2 | Y2 | BOOL | Valve open | $\mathrm{Y} 2=1$ | A 4.2 |
| Release Valve | Y3 | BOOL | Valve open | $\mathrm{Y} 3=1$ | A 4.3 |
| Pressure Release Valve | Y4 | BOOL | Valve open | $\mathrm{Y} 4=1$ | A 4.4 |
| Heater | H | BOOL | Heater on | $\mathrm{H}=1$ | A 4.5 |
| Agitator Motor | M | BOOL | Motor on | $\mathrm{M}=1$ | A 4.6 |

Figure 112 shows the sequential function chart along with the corresponding text, so that is clear how each of the components were created.


Figure 112: Sequential Function Chart

## Section 7.8.5: Validity of a Sequential Function Chart

With sequential function charts, it is easy to create complex networks. However, the question remains if these networks are truly valid and result in proper operation.

Although there exists a method to determine the validity of a sequential function chart, it not guaranteed that it will find all valid networks. The proposed method only determines if the network is valid. If the method fails, there is no guarantee that the chart is truly invalid. The procedure is:

1) Replace all step-transition-step by a single step.

2) Replace all transition-step-transition by a single transition.

3) Replace two parallel transitions by a single transition.

4) Replace two parallel steps by a single step.

5) Remove any self-transitions.


The validity is then determined using the following rule. A sequential function chart is valid when the chart can be reduced to an implementable single step.

Example 48: Determining the Validity of a Network
Determine the validity of the following sequential function chart shown in Figure 113.


Figure 113: Sequential Function Chart for Checking Its Validity

## Solution

The method is followed until no further reductions can be produced. The first step is to apply Rule \#1 to give Figure 114.


Figure 114: First Reduction
Then, Rules \#4 and 5 can be applied to give Figure 115.


Figure 115: Second Reduction
Then Rule \#2 is applied twice to give Figure 116.


Figure 116: Third Reduction
The last step involves applying Rule \#3 to give Figure 117.


Figure 117: Fourth and Final Reduction
The reduced chart is obviously valid, since it contains only a single step and a single transition. Thus, we can conclude that the proposed sequential function chart is in fact valid.

## Section 7.9: Chapter Problems

Problems at the end of the chapter consist of three different types: (a) Basic Concepts (True/False), which seek to test the reader's comprehension of the key concepts in the chapter; (b) Short Exercises, which seek to test the reader's ability to compute the required parameters for a simple data set using simple or no technological aids. This section also includes proofs of theorems; and (c) Computational Exercises, which require not only a solid comprehension of the basic material, but also the use of appropriate software to easily manipulate the given data sets.

## Section 7.9.1: Basic Concepts

Determine if the following statements are true or false and state why this is the case.

1) Global variables can only be used in a single resource.
2) A nonpre-emptive task must always be completed before another task can be started.
3) 1 Prog 12 is a valid function name.
4) TUI_124 and TUI_12456 are equal.
5) $\mathrm{T} \# 4 \mathrm{~d} 4.2 \mathrm{~h}$ represents 4 days and 4.2 hours.
6) LE means less than.
7) A task with a priority value of 3 has the highest priority.
8) In the function-block language, a double arrow denotes a jump between two parts of the function-block network.
9) In the function-block language, we can use $\&$.
10) In ladder logic, a coil can be used to save a value.
11) In ladder logic, we can create feedback loops.
12)In instruction list, the statement SN HIPPO is a valid one.
12) In instruction list, the brackets () delay the implementation of an instruction.
14)In structured text $O R$ has a higher priority compared to AND.
13) In structured text, we can use the CASE command.
14) In sequential function charts, $N$ means that as soon as the step is ended, the action is stopped.
15) In sequential function charts, $R$ means that an action is delayed by the given time.
16) In sequential function charts, it is possible to easily determine the validity of the result chart.
17) Ladder logic is a great idea to create complex, high-level PLC programmes.
18) Instruction list is designed for simple, optimised PLC programmes.

## Section 7.9.2: Short Questions

These questions should be solved using only a simple, nonprogrammable, nongraphical calculator combined with pen and paper.
21)Is the sequential function chart in Figure 118 valid?

22) Give the sequential function charts for the controlling the following systems:
a. Stirring process: Figure 119 shows the P\&ID diagram of a stirred tank B with flow controller US2. The binary input signals to US2 are the high-level alarm L, the start signal $S$ and the end signal E. The binary output signals from US2 are V1 for switching the input valve, V 2 for switching the drain valve, and R for switching the stirrer motor. The following sequence is required: After the start signal reaches 1 ( $\mathrm{S}=1$ ), a liquid is to be dosed by opening the input valve (with $\mathrm{V} 1=1$ ) and turning the stirrer on $(R=1)$ until the level $L$ reaches its upper limit $(L=1)$. Then, valves are closed and the stirrer should alternately between being switched on for 1 min and switched off for 1 min . This switching on and off should, regardless of the current cycle, be ended immediately when the end button is pressed $(\mathrm{E}=1)$. Then the drain valve should be opened for 5 minutes. Then everything should be switched into the idle state, that is, $\mathrm{V} 1=\mathrm{V} 2=\mathrm{R}=0$. When $\mathrm{S}=1$, the process is started from the beginning.


Figure 119: Stirring Process
b. Washing Machine: The control for a simple washing machine, shown in Figure 120 with an explanation of the variables in Table 44, should implement the following sequence. After the input signal START has a value of 1 , valve V 1 should be opened until the binary level sensor L1 reports that the desired water level has been reached. Then, the MOTOR of the washing drum should be switched on alternately for a period of time $T$ _on and switched off for a period of time $T$ _off. During this alternating switching on and off, the electrical HEATING is first switched on until the water temperature has reached a certain value, which is reported by the binary sensor T 2 . Then, the process should wait a certain minimum waiting time $T$ _wait before any further heating occurs. Once the motor has stopped running, the alternative switching of the motor is also stopped. Then, the PUMP is switched on and valve V 2 is opened at the same time. Once the level sensor L0 reports an empty drum, the initial state is reached, and the process can be repeated.


Figure 120: Washing Machine

Table 44: I/O for the Washing Machine

| I/O | Name | Description for a Boolean Value 1 |
| :---: | :---: | :--- |
| Output | V1 | Valve V1 is open. |
| Output | V2 | Valve V2 is open. |
| Output | HEATING | Heating is on. |
| Output | MOTOR | Motor is on. |
| Output | PUMP | Pump is on. |
| Input | START | Start button has been pressed. |
| Input | L0 | Level $L \leq L_{\text {min }}$ |
| Input | L1 | Level $L \geq L_{\text {max }}$ |
| Input | T2 | Temperature $T \geq T_{\text {desired }}$ |

23) In the examples below, various Boolean functions of the form $Q=f(\mathrm{X}, \mathrm{Y}, \mathrm{Z})$ are given. For each, please provide a) the function in its sum-of-products and product-of-sums forms, b) the reduced function found using a Karnaugh diagram; and c) the PLC programme using instruction list, ladder logic, and function-block language for the minimised function.
a. The function shown using ladder logic in Figure 121.
b. The function that implements a majority decision rule for a 2-out-of-3 redundancy system, that is, the output $Q$ is 1 if and only if at least two of the inputs are 1 .
c. The function shown by the truth table in Table 45.
d. The function shown by the truth table in Table 46.
e. The function $Q=f(\mathrm{~B}, \mathrm{G}, \mathrm{M})$ is described as follows:
i. Three sensors in a sorting system measure the properties of parts and deliver the following binary signals to the inputs of a PLC:
1. $\mathbf{B}$, where $\mathrm{B}=1$ means a hole is present,
2. $\mathbf{G}$, where $\mathrm{G}=1$ means green paint is present, and
3. $\mathbf{M}$, where $\mathbf{M}=1$ means a metallic material is present.
ii. Parts produced without errors have the following properties:
4. Either metallic and green (with or without a hole)
5. Nonmetallic, not green, and with a hole.
iii. The output $Q$ of the PLC should be 1 when the part is to be rejected, i.e., if it does not have the above properties of a faultlessly produced part.


Figure 121: Ladder Logic
Table 45: Truth Table I

| $X$ | 0 | 0 | 0 | 0 | 1 | 1 | 1 | 1 |
| :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- |
| $Y$ | 0 | 0 | 1 | 1 | 0 | 0 | 1 | 1 |
| $Z$ | 0 | 1 | 0 | 1 | 0 | 1 | 0 | 1 |
| $Q$ | 0 | 1 | 1 | 0 | 1 | 1 | 1 | 1 |

Table 46: Truth Table II

| $X$ | 0 | 0 | 0 | 0 | 1 | 1 | 1 | 1 |
| :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- |
| $Y$ | 0 | 0 | 1 | 1 | 0 | 0 | 1 | 1 |
| $Z$ | 0 | 1 | 0 | 1 | 0 | 1 | 0 | 1 |
| $Q$ | 0 | 0 | 1 | 1 | 1 | 0 | 1 | 1 |

## Section 7.9.3: Computational Exercises

The following problems should be solved with the help of a computer and appropriate software packages, such as CodeSys.
24) Using CodeSys, write the PLC programme to control a filling station.

# Chapter 8: Safety in the Automation Industry 

As automation becomes more widespread, there is a corresponding need to consider the safety of those who come in contact with the automated systems. Furthermore, automated systems need to take into consideration safety constraints themselves when they implement various actions. For these reasons, it is helpful to briefly review the different safety regulations. Obviously, it is always your engineering obligation to make sure that you have checked the most recent and most relevant regulations and laws applying to your specific situation.

In general, safety can be divided into two parts: physical and digital. Physical safety considers the steps and regulations required to ensure that the plant and its surroundings are safe for the workers, visitors, and the environment. Environmental protection considers not only hazardous substances, but also the emission of electromagnetic radiation or noise. Digital safety considers the steps required to ensure that the communication networks and the associated devices cannot be hacked and used to create an unsafe physical situation. Furthermore, digital safety must ensure that no sensitive information can be stolen or used in an inappropriate manner. This aspect is very important given the increasing legislative interest in making sure that confidential personal information remains confidential.

## Section 8.1: Safety of the Physical System

Before delving into the specifics of physical safety, it is useful to consider some basic principles. A hazard is a source of danger, while a risk is the degree or extent of peril that a hazard could create. Hazards arise from such things as moving objects, stored energy, and explosions. Essentially, hazards primarily arise from the release of latent (or hidden) energy. These include:

1) Kinetic Energy, which is the energy resulting from moving objects, can present hazards in any rotating equipment, such as pumps and turbines, vehicles, and conveyor belts.
2) Potential Energy, which is the energy resulting when an object unexpectedly falls. Such releases of energy are common in structural failures, such as building collapses.
3) Work, which is energy stored in springs, electrical circuits, and other devices. When released, this can lead to catastrophic failure and impressive damage. A common electrical example is a short circuit.
4) Heat, which is energy crossing a boundary. Heat exchange is always based on a temperature difference between two points in space. Since heat always influences the environment, touching hot surfaces, for example, can lead to burns. The same situation applies to touching very cold surfaces, which can also lead to burns or frostbite.

In order to increase the safety of a physical system, it can be useful to consider the following steps when designing the overall system:

1) Minimise: Avoid using (or use as little as possible) any hazardous substances.
2) Substitute: Replace hazardous substances by ones that are less hazardous but with similar properties.
3) Moderate: Replace extreme operating conditions by ones that are less severe.
4) Simplify: Create designs or processes that are less complex and simpler.
5) Isolate: Create situations where the effect of a hazard has minimal impact on the overall system, for example, physically separating the office space from the production area.
The above ideas can be framed within the inherently safer predesign (ISPD) approach, which consists of the following 4 steps:
6) Identify: Determine what the hazards are and how they can impact the process.
7) Eradicate: Remove the effects of as many of the hazards as possible, for example, by designing fail-safe systems, that is, should a system encounter a failure it ends up in an inherently safe mode that cannot cause further problems. A common example of this is the design of control valves to fail open or closed depending on the process conditions, for example, a control valve for a critical cooling jacket should fail open so that the system remains cooled even if the valve has failed.
8) Minimise, simplify, and moderate: If a hazard cannot be eradicated, then try and limit its effects on the system.
9) Isolate: Create structures that will limit the amount of damage that can occur, for example, by segregating hazardous operations.

Another approach to this situation is to perform a hazard-and-operability study (HAZOP) to determine where the given hazards lie and how best to mitigate them. The HAZOP procedure
consists of the following four steps: determine, using a systematic search, possible faults; find the causes for the faults; estimate the effects of the faults on the system; and suggest appropriate countermeasures. A HAZOP requires that a detailed plan of the process be available. Thus, such analysis is often performed at a later stage of design after the overall process design has already been completed. This implies that, due to the high cost or structural changes required, it could be too late to make substantive changes at this point. Thus, HAZOPs are often used to provide an understanding of the hazards present and their risk.

Finally, it can be mentioned that when designing safety systems, it is useful to consider redundancy, that is, having more than one channel or way to achieve a given task. As well, this implies that different channels should be used for different tasks (control, alarm, and monitoring), so that there is less risk of catastrophic failure. Thus, for example, alarms and monitoring should not occur on the same channel. If it should fail, alarm signals can no longer be transmitted in addition to the less important monitoring signals. There are two main ways to implement redundancy. In the first approach, often called homogeneous redundancy, the same task is implemented by multiple (identical or similar) devices, for example, the temperature is measured using three temperature sensors at the same point. However, this approach carries the risk that systematic errors in the sensors or devices are propagated into the system. In the second approach, often called heterogeneous redundancy or redundancy by diversity, the same task is implemented using different devices, for example, three separate path-planning computer systems that use different computational algorithms could be implemented and the decision taken by majority vote, that is, if two or more algorithms give the same result it is followed. Selecting the appropriate approach depends on the requirements and standards of the process.

## Section 8.1.1: Quantifying Risk and the Safety Integrity Level

In order to properly understand and implement a safety system, it is necessary to quantify the risk in some manner. One approach is to define the risk $R$ as

$$
\begin{equation*}
R=C D \tag{249}
\end{equation*}
$$

where $C$ is the impact of the hazard and $D$ is the frequency of the hazard. The frequency of the hazard is defined as

$$
\begin{equation*}
D=F P W \tag{250}
\end{equation*}
$$

where $F$ is the occurrence frequency, $P$ is the probability that the hazard cannot be mitigated, and $W$ is the probability that without any safety system in place an undesired state will be reached. Since it makes little sense to define the risk with an exact value, it is common to instead associate a given level to the individual variables and look at the combination of levels.

For $C$, the following levels are defined:

- C1: Minor injuries
- C2: Major or permanent injuries of one or more people or a single death
- C3: Up to 5 deaths
- C4: More than 5 deaths.

For $F$, two levels are defined:

- F1: During the course of the day, people are in the danger zone less than or equal to $10 \%$ of the time.
- F2: During the course of the day, people are in the danger zone more than $10 \%$ of the time.

For $P$, two levels are defined:

- P1: It is possible to mitigate the hazard (one should provide the appropriate steps to do so).
- P2: It is not possible to mitigate the hazard.

For $W$, three levels are defined:

- W1: The undesired state will occur less than once every 10 years.
- W2: The undesired state will occur less than once per year.
- W3: The undesired state will occur more often than once per year.

The above levels are then combined to give an overall safety integrity level (SIL). There are four safety integrity levels labelled SIL1 to SIL4, where the larger number reflects a more hazardous level. The relationship between the different parameters is shown in Figure 122.


Figure 122: Relationship between the parameters and the safety integrity levels

## Section 8.2: Safety Regulations

Safety regulations in Germany (and most of Europe and even the world) come in many different flavours and legal force. In Germany, standards are developed by the German Institute for Standardisation (DE: Deutsches Institute für Normung, commonly abbreviated as DIN). DIN publishes standards not only for safety but for all sorts of other aspects, such as formatting of letters, transliteration, and engineering. All standards developed or accepted by DIN are prefixed by the letters $D I N$ followed by a number. Additional letters can be added to show the general applicability of the standards. Using DIN EN implies that we are dealing with a German edition of a European standard, while DIN ISO implies that we are dealing with a German edition of an international standard developed by the International Standards Organisation (ISO). Obviously, it is possible to have a DIN EN ISO label which would denote a German edition of a European and international standard. In general, when there are multiple prefixes, this implies that we are dealing with the same standard under different guises, for example, a DIN ISO \#\#\#\# standard would be the same as the corresponding ISO \#\#\#\#. Examples of different DIN standards are:

1) DIN 31635 which standardises the transliteration of the Arabic script used to write Arabic, Ottoman Turkish, Iranian, Kurdish, Urdu, and Pasto
2) DIN EN ISO 216 which standardises writing paper and certain classes of printed matter, trimmed sizes, A and B series, and indication of machine direction
3) DIN ISO 509 which covers the production of technical drawings, relief grooves, and types and dimensions of them.
4) DIN EN 772-7 which covers clay masonry units, namely the determination of water absorption of clay masonry damp proof course units by boiling water.
Other countries use similar systems, for example, Austria prefixes its codes with ÖNORM while the United States of America uses the ANSI system.

Safety regulations in Germany (and most of Europe) for machines can be split into three types:

- Type A Regulations: These regulations cover the basic safety concepts, design principles, and general aspects of machines. An example of a Type A regulation is the EN ISO 12100 standard that provides the general principles of design.
- Type B Regulations: These regulations are baseline safety standards and requirements for protective equipment. There are two subtypes: B1-Regulations that cover specific safety aspects and B2-Regulations that cover the standards for protective equipment required for machines. An example of a Type B1 regulation is the EN ISO 13855 standard that provides the guidelines for the arrangement of protective devices, while a Type B2 regulation would be the EN 953 standard that considers fixed guards for machines.
- Type C Regulations: These regulations are the specific safety standards for a particular machine or group of machines. An example of a Type C regulation would be the EN 693 standard that provides the standards for hydraulic presses.


## Section 8.3: Digital Safety

In today's world, where many components of a plant are now linked digitally using networks, it is imperative that the digital system be secured. The main topic in digital safety focuses on preventing intruders from entering the digital system and making unauthorised changes to the system, that is, digital safety must prevent hacking from occurring. Hacking can expose a
company to various threats including loss of proprietary information, damage to the process by illegal changing of process operating conditions, and legal issues due to distribution of confidential personal information.

Digital security can be achieved by creating strong passwords, changing the passwords on a regular basis, and checking for any intruders. Strong antivirus software should also be implemented. Multiple interacting networks should be considered to provide additional levels of security. Finally, the workers should be educated in proper digital safety, as the digital system is only as strong as its weakest link. People placing unknown USB-drivers into work computers can easily introduce unwanted viruses into the company network.

## Section 8.4: Chapter Problems

Problems at the end of the chapter consist of three different types: (a) Basic Concepts (True/False), which seek to test the reader's comprehension of the key concepts in the chapter; (b) Short Exercises, which seek to test the reader's ability to compute the required parameters for a simple data set using simple or no technological aids. This section also includes proofs of theorems; and (c) Computational Exercises, which require not only a solid comprehension of the basic material, but also the use of appropriate software to easily manipulate the given data sets.

## Section 8.4.1: Basic Concepts

Determine if the following statements are true or false and state why this is the case.

1) Process safety is never a topic of concern.
2) A hazard is something that has a high likelihood of causing harm.
3) A short circuit is an example of a hazard caused by the release of stored energy or work.
4) Collapsing buildings are an example of hazards posed by kinetic energy.
5) Hazards that cannot be eliminated should be isolated.
6) Designing complex processes is a good risk mitigation strategy.
7) The fail-safe principle states that a failing system should always end up in a safe mode.
8) Using the fail-safe principle, valves should always fail closed.
9) A HAZOP seeks to identify the risks and how best to mitigate them.
10) Redundancy means that a single temperature sensor is used for process monitoring and process control.
11) Selecting a course of action based on two-out-of-three voting is an example of redundancy by diversity.
12) A risk with an occurrence frequency greater than $10 \%$ is always an SIL 4 risk.
13) A risk with a hazard impact level of $C 3$ implies that it can cause only minor injuries.
14) A risk with a mitigation level of P 2 implies that it cannot be mitigated.
15) A risk with a $W$ level of W 3 implies that the undesired state will occur more often than once per year.
16) In Germany, Type A regulations cover the detailed requirements for the safety of a specific machine.
17) An example of a $B 2$ regulation would be the design of protective guards for machines.
18) An example of a $C$ regulation would be the standards for the design of chemical reactors.
19) Using simple passwords and minimal digital security is a good idea for a critical chemical process.

## Section 8.4.2: Short Questions

These questions should be solved using only a simple, nonprogrammable, nongraphical calculator combined with pen and paper.
20)Determine the safety integrity levels for the following risks:
a. A risk evaluated as belonging to $\mathrm{C} 2, \mathrm{~F} 1, \mathrm{~W} 2$, and P 2 .
b. A risk evaluated as belonging to $\mathrm{C} 3, \mathrm{~F} 2, \mathrm{P} 1$, and W 1 .
21)Determine some of the safety issues involved with the production of the following chemicals. Based on the principles of safety considered here, would you recommend that these chemicals be produced by a new start-up that has no chemical engineering background.
a. Acrylonitrile $\left(\mathrm{CH}_{2} \mathrm{CHCN}\right)$, which is produced from propene, air, and ammonia.
b. Propene $\left(\mathrm{C}_{3} \mathrm{H}_{6}\right)$, which is produced from ethene and 2-butene.
22)Perform a HAZOP on the compressor unit shown in Figure 123.


Figure 123: P\&ID for a Compressor Unit

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## Appendix I: Partial Fractioning

When solving equations in the frequency domain and it is desired to convert them back into the time domain, it may be necessary to perform partial fractioning to obtain a solution. Although there exist many different approaches, the following is one simple method that will allow the final result to be obtained easily. Consider a rational function of the form:

$$
\begin{equation*}
\frac{N(s)}{D(s)}=\frac{N(s)}{\prod_{i=1}^{n_{i}}\left(\alpha_{i} s+\beta_{i}\right)^{n_{i}} \prod_{j=1}^{n_{q}}\left(\alpha_{j} s^{2}+\beta_{j} s+\gamma_{j}\right)^{n_{j}}} \tag{251}
\end{equation*}
$$

where $n_{l}$ represents the number of distinct linear terms, $n_{q}$ the number of distinct irreducible quadratics (those that have imaginary roots as their solution), and $\alpha, \beta$, and $\gamma$ are known constants. Let $n$ be the overall order of the system. In order to perform partial fractioning, write the following fraction depending on the form of the root:

1) For each linear term $(\alpha s+\beta)^{n}$, put the term, $\sum_{k=1}^{n} \frac{B_{k}}{(\alpha s+\beta)^{k}}$.
2) For each irreducible quadratic term, $\left(\alpha s^{2}+\beta s+\gamma\right)^{m}$, put the term, $\sum_{k=1}^{m} \frac{A_{k}+B_{k} s}{\left(\alpha s^{2}+\beta s+\gamma\right)^{k}}$.

Once the form of the partial fractioning solution has been obtained, it is necessary to solve for the unknown parameters. First, cross-multiply, so that the denominators are the same. Then, it is necessary to solve for the unknown parameters by equating the unknown side with the known, $N(s)$. The easiest way to solve this is using the following approach:

1) For each linear term, set $s=-\beta / \alpha$ to obtain $B_{n}$ of the linear terms. This will reduce the equation to the form $B_{n} \prod_{\substack{i=1 \\ i \neq n}}^{n_{1}}\left(\alpha_{i} s+\beta_{i}\right)^{n_{i}} \prod_{j=1}^{n_{4}}\left(\alpha_{j} s^{2}+\beta_{j} s+\gamma_{j}\right)^{n_{j}}=N(s)$ evaluated at the given root.
2) For each quadratic term, set $s$ equal to the imaginary roots. This will also reduce the equation to a simpler form and allow for $A_{n}$ and $B_{n}$ to be solved.
For the remaining terms, create a system of equations by selecting different values of $s$ and evaluating the known values, so that the remaining unknowns can be solved. You will need $n-$ $n_{l}-2 n_{q}$ equations in order to find the remaining $n-n_{l}-2 n_{q}$ terms.

## Example 49: Partial Fractioning

Consider the following fraction

$$
\begin{equation*}
\frac{3 s+1}{(s+2)(s+1)^{2}\left(s^{2}+1\right)} \tag{252}
\end{equation*}
$$

for which we wish to determine the partial fraction form.

## Solution

First, we need to write the general form, that is, for each of the terms in the denominator, we will write the corresponding partial fraction using the rules above. This will give

$$
\begin{equation*}
\frac{A}{s+2}+\frac{B}{s+1}+\frac{C}{(s+1)^{2}}+\frac{D s+E}{s^{2}+1} \tag{253}
\end{equation*}
$$

It can be noted that for the first term $s+2$, as well as the last term $s^{2}+1$, there will only be a single component, since its exponent is one. For the middle term, $(s+1)^{2}$, there will be two terms, since the exponent is two. For the linear terms, we set a simple constant term in the numerator, while for the quadratic term, we include a linear term in the numerator.

Next, we need to determine the values of the constants in the numerator. Before doing this, let us cross-multiply and determine the general form of the numerator

$$
\begin{equation*}
A(s+1)^{2}\left(s^{2}+1\right)+B(s+2)(s+1)\left(s^{2}+1\right)+C(s+2)\left(s^{2}+1\right)+(D s+E)(s+2)(s+1)^{2}=3 s+1 \tag{254}
\end{equation*}
$$

As was previously mentioned, we can see that setting $s=-1$ or -2 will cancel every term but one, allowing us to effortless compute that term. This gives for $s=-1$

$$
\begin{align*}
C(-1+2)\left((-1)^{2}+1\right) & =3(-1)+1 \\
C & =\frac{-2}{2}=-1 \tag{255}
\end{align*}
$$

Similarly, for $s=-2$, we get

$$
\begin{align*}
A(-2+1)^{2}\left((-2)^{2}+1\right) & =3(-2)+1 \\
A & =\frac{-5}{5}=-1 \tag{256}
\end{align*}
$$

For the quadratic term, we can set $s= \pm j$, which will give us a linear system of equations in two unknown ( $D$ and $E$ ) that we can then solve to obtain the value of $D$ and $E$. This gives

$$
\left\{\begin{array}{c}
(D j+E)(j+2)(j+1)^{2}=3 j+1  \tag{257}\\
(-D j+E)(-j+2)(-j+1)^{2}=-3 j+1
\end{array}\right.
$$

Placing all the constant terms on the right gives

$$
\left\{\begin{align*}
D j+E & =0.5-0.5 j  \tag{258}\\
-D j+E & =0.5+0.5 j
\end{align*}\right.
$$

Adding the two equations together gives that $E=(0.5+0.5) / 2=0.5$. From this, it follows that $D$ $=-0.5$. It can be noted that these two equations will always be complex conjugates of each other and one can use this fact to solve them without necessarily computing both components.

The remaining term $B$ can be found by selecting an arbitrary value of $s$ (that has not already been used) and solving Equation (254). The known constants are inserted as required. Setting $s=$ 0 , we get

$$
\begin{align*}
-1(1)^{2}(1)+B(2)(1)(1)-1(2)(1)+(0.5)(2)(1)^{2} & =1 \\
B & =\frac{3}{2}=1.5 \tag{259}
\end{align*}
$$

Obviously, we have the correct answer if the numerator given by Equation (254) holds. This can be used to check the solution obtained.


[^0]:    ${ }^{1}$ This word is ultimately derived from the Greek word av̉tóg, meaning self, and an unattested root, which comes from the proto-Indo-European word *méntis $\sim$ mntéis, meaning thought (which in turn gave us such words as mind in English)

[^1]:    ${ }^{2}$ In this textbook, the two words will not be distinguished. Some authors consider the process to be the overall way something is done, while a system represents the actual physical realisation of the process. Since in many cases, these words are used interchangeably, the textbook will primarily use system.

[^2]:    ${ }^{3}$ Often also called a state, but this term is avoided since it has another meaning in control and process analysis.

[^3]:    ${ }^{4}$ A psi is a unit of pressure in the imperial system of measurement. It is an abbreviation for pounds (force) per square inch. The conversion factor is $1 \mathrm{psi}=6.894757 \mathrm{kPa}$. The $g$ represents gauge or the value above atmospheric pressure.

[^4]:    ${ }^{5}$ Equivalently, the real component of all the poles must be less than 0 or the system is stable.

[^5]:    ${ }^{7}$ The roots are expressed in terms of $z$. Equivalently, the system is stable.

[^6]:    ${ }^{8}$ In automation engineering, we often deal with deviation variables that are defined as $\tilde{x}=x-x_{s s}$, where $x_{s s}$ is some steady-state value. Since we assume that the process is initially in steady state, this means that the initial conditions will always be zero.
    ${ }^{9}$ The inverse exists since the $\mathcal{A}$-matrix only contains numeric entries.

[^7]:    ${ }^{10}$ In some programmes, an "astronomical" arctangent function is defined as $\arctan 2(y, x)$, which will return the value in the correct quadrant based on the signs of $y$ and $x$. If available, it should be used.

[^8]:    ${ }^{11}$ In the discrete domain with a transfer function, it is very important to take into consideration what variable is being used and how the stability condition is being phrased. Often, especially in control, stability is discussed in terms of $z^{-1}$, in which case the above rules need to be inverted, that is, stability implies poles outside the unit circle. For consistency, stability in this book will be discussed in terms of $z$.

[^9]:    $12 \cosh$ is the hyperbolic cosine function defined as $1 / 2\left(\mathrm{e}^{x}+\mathrm{e}^{-x}\right)$ and $\sinh$ is the hyperbolic sine function defined as $1 / 2\left(\mathrm{e}^{x}-\mathrm{e}^{-x}\right)$.

[^10]:    ${ }^{13}$ Often it will also be used for a time domain representation by ignoring the composition and summation rules and using slightly different models. In such cases, they are better called process flow diagrams, which are described in Section 4.2.

[^11]:    ${ }^{14}$ The older German standard DIN 19227 matches closely the ISA or North American standard. Differences include how the different sensors or functions are denoted and minor details regarding the proper location for various pieces of additional information. This book will follow the new standard without necessarily making any comments about the other possibilities in order to avoid confusion.
    ${ }^{15}$ This represents a general pipe.

[^12]:    ${ }^{16}$ PCE is an abbreviation for "Process Control Engineering".

[^13]:    ${ }^{17}$ Officially, this is a free variable to be set as needed. Practically, it is often used for conductivity.

[^14]:    18 continuous domain: $\operatorname{Re}(\lambda)<0$; discrete domain: $\|\lambda\|<1$
    ${ }^{19}$ Often, the estimated state is replaced by the true state value. As it will be shown, this makes sense in many cases since we have full information about the states by actually measuring them.

[^15]:    ${ }^{20}$ This term comes from the days of typewriters, where the writing head with the ink ribbon was attached to a carriage that had to be pushed at the end of each line to the beginning of the new line.

[^16]:    ${ }^{21}$ The underscore is not obligatory to separate the units, but it does help readability of the text.

[^17]:    ${ }^{22}$ The value on the left is always transferred to the right-hand side.

[^18]:    ${ }^{23}$ tnw represents a Boolean variable that determines if the given link should be performed．

[^19]:    ${ }^{24}$ snw represents a subnetwork that returns a Boolean value.

