



Norwegian University of
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Nonlinear Model Predictive Control of Gravity Separators

Melissa Florine Dlima

Chemical Engineering

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Supervisor: Sigurd Skogestad, IKP

Co-supervisor: Christoph Backi, IKP

Norwegian University of Science and Technology
Department of Chemical Engineering

Abstract

For economical reasons, it is necessary to operate separators such that the the purity specification is maximized while having the levels and the system pressure controlled. In this thesis, a nonlinear model predictive control is applied to a three phase subsea gravity separator system.

The main modeling principle used to describe the oil droplets rising and settling of the droplets is given by Stokes law. Nonlinear dynamic equations were derived for the water level, total liquid level and for the pressure of the gas in the system based on the inflow and outflow dynamics, and were implemented in MATLAB as script files. The equations were nonlinear due to the geometry of the separator. These nonlinear models were formulated as a semi-explicit DAE system.

The NMPC optimization problem used in this report is solved using direct collocation method using CasADi software within the MATLAB programming environment. Simulation studies have been carried out for four cases, namely setpoint tracking, disturbance rejection, sensitivity to measurement noise and the optimal water level investigation. The performance of the closed loop Nonlinear Model Predictive Control (NMPC) has been studied based on the simulation results.

Simulation results were carried out to analyze the effect of measurement noise on the performance of the controller. Based on this analysis, the performance of the system for remaining cases was studied. The controller was analyzed for the setpoint tracking scenario where step change in water level was introduced to the system. Additionally, the performance of the controller to reject the disturbances was studied. Finally, the optimal level of water in the separator was investigated based on maximization of the total volume of oil entering its native phase.

Preface

This master thesis was written for the process systems engineering group at the Norwegian University of Science and Technology. This work was carried out during the Autumn of 2017 at the Department of Chemical Engineering.

I would like to express my deepest gratitude to my supervisor professor Sigurd Skogestad, for giving me an opportunity to work on this project and also for his support and useful discussions during my work.

I am also extremely thankful to my co-supervisor Dr. Christoph J. Backi for his constant guidance throughout the project, for being available to me even while being away on conferences, for providing the technical background on which this project is based and for reviewing the thesis. I would also like to thank Ph.D candidates Dinesh Krishnamoorthy, Tamal Das and Post Doctoral candidate Eka Suwartadi for their key inputs in understanding CasADi. I am grateful for the opportunities given by the process systems engineering group to participate in the meetings and discussions.

Finally, I would like to thank my family and friends without whom this journey would not have been possible. I am especially grateful to my parents who believed in me and wanted the best for me. I am also thankful to my sister, Megan for encouraging me throughout this experience.

Declaration of Compliance

I declare that this thesis is an independent work in agreement with the exam regulations of the Norwegian University of Science and Technology (NTNU).

Trondheim, Norway

June 12, 2017

Melissa Florine Dlima

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List of Symbols

Latin Letters

Symbols	Description	Unit
A_L	Cross sectional area of liquid in the separator	m^2
A_O	Cross sectional area of water in the separator	m^2
A_p	Reference area of the droplet	m^2
A_W	Area of water in the separator	m^2
b	Time invariant parameters	-
C_d	Drag Coefficient	-
d	Disturbance vector	-
F_b	Buoyant force	N
F_d	Drag force	N
F_g	Gravitational force	N
g	Acceleration due to gravity	$m.s^{-2}$
h_L	Liquid level	m
h_L^{SP}	Liquid level setpoint	m
h_O	Oil level	m
h_W	Water level	m
h_W^{SP}	Water level setpoint	m
l_p	Droplet position	m
N_{Re}	Reynolds number	-
n	Initial number of droplets	-

n_G	Amount of gas	<i>moles</i>
p	System pressure	<i>bar</i>
p^{SP}	System pressure set point	<i>bar</i>
$q_{G,in}$	Gas inflow	$m^3 s^{-1}$
$q_{G,out}$	Gas outflow	$m^3 s^{-1}$
$q_{L,in}$	Liquid inflow	$m^3 s^{-1}$
$q_{L,out}$	Liquid outflow	$m^3 s^{-1}$
$q_{W,in}$	Water inflow	$m^3 s^{-1}$
$q_{W,out}$	Water outflow	$m^3 s^{-1}$
R	Universal gas constant	$J mol^{-1} K^{-1}$
r	Radius	<i>m</i>
T	Temperature	<i>K</i>
T_c	Control horizon	<i>s</i>
T_p	Prediction horizon	<i>s</i>
u	Vector of manipulated variables	-
u_{opt}	Optimal control sequence	-
V_G	Volume of gas in the separator	m^3
V_L	Volume of liquid in the separator	m^3
V_{OoW}	Total oil volume leaving water phase	m^3
V_{Sep}	Volume of the separator	m^3
V_{WoO}	Total water volume leaving oil phase	m^3
v_h	Horizontal velocity	$m s^{-1}$
v_v	Vertical velocity	$m s^{-1}$
x	Vector of state variables	-

Greek Letters

Symbols	Description	Unit
α	Water cut	-
β	Oil cut	-
η_O	Efficiency of oil	-
η_W	Efficiency of water	-
μ_O	Viscosity of oil	$kg(ms)^{-1}$
μ_W	Viscosity of water	$kg(ms)^{-1}$
ϕ_{oo}	Fraction of oil entering the oil phase	-
ϕ_{ow}	Fraction of oil entering the water phase	-
ϕ_{wo}	Fraction of water entering the oil phase	-
ϕ_{ww}	Fraction of water entering the water phase	-
ρ_G	Density of gas	kgm^{-3}
ρ_O	Density of oil	kgm^{-3}
ρ_W	Density of water	kgm^{-3}

Abbreviations

CasADi	Computer algebra system for Automatic Differentiation
CV	Controlled Variable
DAE	Differential Algebraic Equations
DMS	Direct Multiple Shooting
DOF	Degree of Freedom
DSS	Direct Single Shooting
HJB	Hamilton- Jacobi- Bellman
IPOPT	Interior Point OPTimizer
KKT	Karush-Kuhn-Tucker conditions
LP	Linear Programming
LQG	Linear Quadratic Gaussian
LTI	Linear Time Invariant
MINLP	Mixed Integer Nonlinear Programming
MPC	Model Predictive Control
NLP	Nonlinear Programming
NMPC	Nonlinear Model Predictive Control
OCP	Optimal Control Problem
ODE	Ordinary Differential Equation
PID	Proportional Integral Derivative
PMP	Pontryagins Maximum principle
QP	Quadratic Programming
RHC	Receding Horizon Control
SQP	Sequential Quadratic Programming

Introduction

The recent interest in subsea processing is mainly driven by the advantages of reduced pumping cost, water management cost and reduced environmental impact due to water re-injection (Moraes et al., 2013). Also, the subsea processing makes it possible to use smaller or no platforms at all, due to the fact that the equipment used for processing is placed on the seabed rather than topside (Albuquerque et al., 2013).

The majority of oil wells produce a mixture of mainly gas, oil and water. Conventionally, to separate this mixture, a series of separators is used placed starting with a rough separation of gas-liquid mixture in three phase gravity separators to separate the mixtures into oil, water and the gas phase. The produced water that has been separated can be used for re-injection into the reservoir, which helps to increase the pressure and ultimately the recovery. With this advantage, many subsea separators have been implemented around the world. In addition to the increased production, the companies need to comply with strict requirements concerning the disposal of produced waste water. There are many regulations for the maximum

allowable concentration of oil in the water that could be discharged into the sea which are set in terms of daily maximum and monthly average (Prescott, 2012). For example, the maximum limit for hydrocarbon discharge in North Sea is set at 30ppm (Durdevic et al., 2017).

Three phase horizontal gravity separators are considered in this thesis. The separation is based on density difference between the single phases as well as gravitational forces. A proper control scheme allows to separate the mixtures based on maximization of oil production, minimization of oil in the produced water or to control the levels at their desired setpoints.

For the last two decades, predictive control has been found to be a very good controller design scheme mainly due to the high performance controllers that can be used for multivariable processes and also that the constraints in the process are handled in a straightforward and organized way (Nunes, 2001).

According to Stokke et al. (1994), a study was carried out on Statoil's Statfjord A crude oil production platform, where Model Predictive Control (MPC) was applied to control the water, oil and the system pressure. In this process, due to the higher production levels and large variations in the oil levels, experienced operators had to adjust several PI controllers simultaneously. When the PI controllers are implemented in a system, it is inappropriate to manipulate more than one PI controller at a time. Therefore, in order to solve this problem, a Model Predictive Control layer is placed on top of the PI control loops, so that it is possible to optimally and dynamically control the setpoints of the PI controllers.

There has not been much work done on Model Predictive Control applied to the gravity separator system. Nunes (2001) used a polynomial operator technique to study the stability and performance of predictive control applied to a gravity separator system. Here, the nonlinear system was linearized around a steady state

and proves that the linear model predictive control is a good control method for oil -water-gas separators. Stokke et al. (1994) studied both MPC and LQG (Linear Quadratic Gaussian) to control the levels in an oil-water-gas separator train. It was found that MPC was far superior to LQG controller since the LQG controller was sensitive to noise.

The remainder of this thesis is structured as follows. Chapter 2 outlines majorly the description of the process, introduction to optimization and optimal control, methods to solve optimal control problems and the control structure hierarchy in chemical plants. Chapter 3 presents the control oriented modeling approach. Chapter 4 outlines the theory of Model Predictive Control (MPC). Chapter 5 describes the simulation results and chapter 6 concludes the thesis with an outlook on future work.

Background

2.1 Particle movement through the fluid

To understand the separation process, it is important to understand the principle of emulsions. An emulsion is a dispersion of one liquid in another liquid, where both liquids are immiscible. The phase in which the droplets are present is the continuous phase and the droplets are termed as the dispersed phase. The main principle of gravity separation is the separation of droplets based on density differences and gravitational forces. The forces exerted on the droplets are the drag force, buoyant force and the gravitational forces as shown in Figure 2.1.

The drag force also known as the frictional force will be exerted on the droplet in the viscous fluid. This force is always in the direction opposite to the velocity of the droplet and is given by

$$F_d = \frac{C_D v^2 \rho A_p}{2}, \quad (2.1)$$

where C_D is the drag coefficient, v is the relative velocity of the droplet with

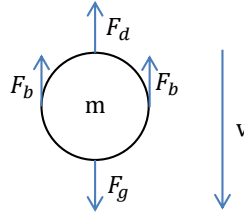


Figure 2.1: Forces acting on a droplet settling through a continuous phase.

respect to the surrounding fluid, ρ is the density of the surrounding fluid and A_p is the reference area of the droplet.

The droplet experiences an upward force which opposes the weight of the particle and is termed the buoyant force. It comes from the pressure exerted on the droplet by the fluid and is given by

$$F_b = V_p \rho g, \quad (2.2)$$

where V_p is the volume of the droplet and g is the acceleration due to gravity.

In addition to the drag and the buoyant force, the particle is exposed to the gravitational force given by

$$F_g = V_p \rho_p g, \quad (2.3)$$

where ρ_p is the density of the droplet.

The drag coefficient C_D depends on the Reynolds number and it can be defined for several flow regimes. For laminar flow, Stokes' law is applied and the drag coefficient is given by

$$C_D = \frac{24}{N_{Re}}, \quad (2.4)$$

where the Reynolds number N_{Re} is given by

$$N_{Re} = \frac{d_p v \rho}{\mu}, \quad (2.5)$$

where d_p is the diameter of the droplet and μ is the viscosity of the medium in which the droplet is present.

At equilibrium, gravitational force on the particle is equal to the drag force and the buoyant force. Examining equations (2.1), (2.2), (2.3) for the forces acting on the particle, yields the terminal velocity of the particle in the laminar flow, assuming Stokes' law valid, and is given by

$$v = gd_p^2 \frac{(\rho_p - \rho)}{18\mu}. \quad (2.6)$$

2.2 Horizontal Gravity Separator Description

In this thesis, a horizontal gravity separator is considered for bulk separation of gas, oil and water. Many different separator types are used in industry, differing in shape and internal devices. Thereby, the three phase horizontal gravity separator as presented in Figure 2.2 is one of the most common types being employed (Mendes et al., 2012).

Figure 2.2 shows volumetric liquid and the gas inflows $q_{L,in}$ and $q_{G,in}$ respectively. The liquid which is a mixture of oil and water enters the separation chamber and separates out into the continuous water and the oil layers based on the difference in density and the gravitational forces as introduced in section 2.1. The oil flows over a weir plate to the oil chamber where the valve u_O is manipulated to extract the oil from the separator vessel. Also, the valves u_W and u_G are used to control

the outflows of water and gas from the separator. The control of the levels h_W , h_L and the system pressure p is done using the valves u_W , u_O and u_G respectively.

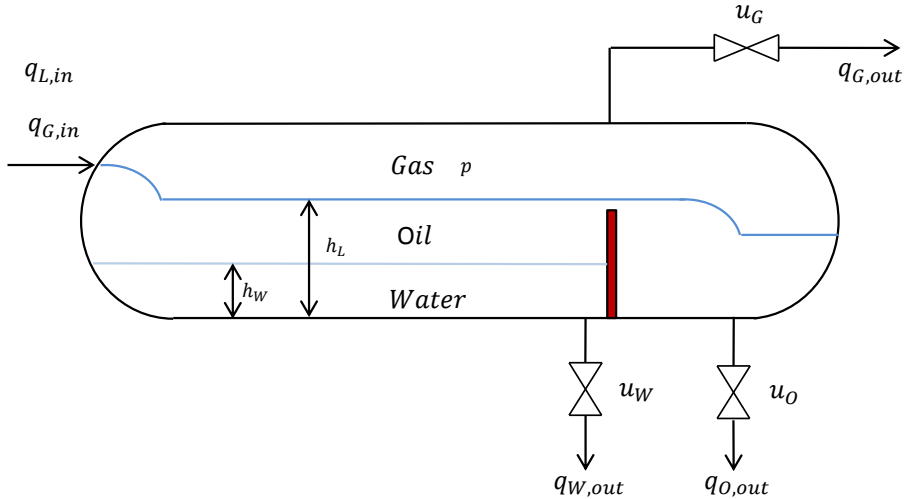


Figure 2.2: Schematic of the three phase horizontal gravity separator.

2.3 Introduction to Optimization

Mathematical optimization is an important tool for the analysis of physical system. This tool has been extensively used for the selection of best elements from a set of available elements. To analyze the system of study, an objective must be identified which could be the maximization of a company's profits or minimization of an energy consumption in a chemical plant. In either of the cases, the objective is largely determined by certain characteristics of the system namely the variables or unknowns. The aim of optimization is to find those unknown values or variables that optimize the objective. Usually, the variables are constrained or restricted to some bounds. The process of deciding the objective of the given problem, along with variables and bounds on the variables is termed modeling of the given system.

After the formulation of the model, an appropriate optimization algorithm has to be chosen to solve the problem of interest.

2.3.1 Mathematical Formulation

Optimization is the maximization or minimization of an objective function subject to variables which are constrained by bounds. Optimization models represent the problem choices as decision variables and attempt to find the variables that maximize or minimize the objective function value subject to constraints on variables. The optimization problem can be formulated as follows:

$$\min_x f(x) \tag{2.7a}$$

$$\text{subject to } g_i(x) = 0 \quad i = 1, \dots, m \tag{2.7b}$$

$$h_j(x) \leq 0 \quad j = 1, \dots, p \tag{2.7c}$$

$$x_{min} \leq x \leq x_{max} \tag{2.7d}$$

where $f(x)$ is the objective function, $g(x)$ are the set of equality constraints and $h(x)$ are the set of inequality constraints, x are the decision variables constrained by lower and upper bounds x_{min} and x_{max} respectively.

If the goal is to minimize the objective function, then the minimization problem can be written as

$$\min_x f(x) \Leftrightarrow \max_x -f(x)$$

Now the minimization problem given by (2.7) can be formulated as the maximiza-

tion problem

$$\max_x -f(x) \tag{2.8a}$$

$$\text{subject to } g_i(x) = 0 \quad i = 1, \dots, m \tag{2.8b}$$

$$h_i(x) \leq 0 \quad i = 1, \dots, p \tag{2.8c}$$

$$x_{min} \leq x \leq x_{max} \tag{2.8d}$$

Optimization problems can be of many classes comprehensively presented in the book by Nocedal and Wright (2006). In this thesis, a brief introduction to a few optimization problem types will be discussed.

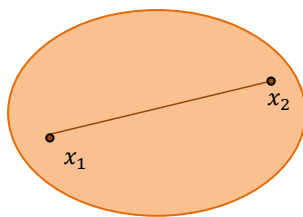
Optimization problems formulated in the general form (2.8) can be classified based on the linearity, nonlinearity, convexity of the objective function and constraints. It also depends on how large or small the number of variables in the optimization problem.

Constrained and unconstrained optimization: In case of unconstrained optimization problems, the general form (2.8) does not have any equality or inequality constraints, meaning Equations (2.8b) - (2.8d) are absent. These problems consider maximizing or minimizing the objective function value with no restriction on the decision variable values. However, most of the problems in engineering applications are constrained.

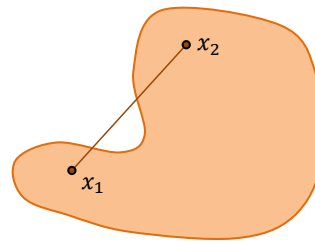
Local and global optimization: Many optimization problems attempt to find a local solution, which is a point at which the objective function value is smaller than the neighbouring points. A Global solution is the point at which the objective function value is smaller than all feasible points.

Convex and non-convex optimization: The optimization problem is said to be con-

convex if the objective function is a convex function, all the equality constraints are linear and inequality constraints are concave. A set is defined as a convex set if a line segment joining any two points in that set lies within that set as shown in Figure 2.3a. The function which is defined in an interval is said to be convex if the straight line segment joining any two points in that interval is as shown in Figure 2.4a.

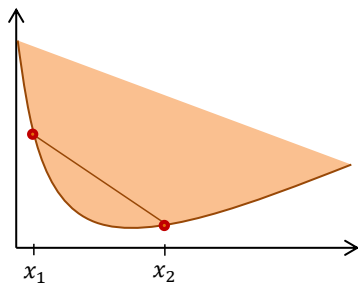


(a) A convex set

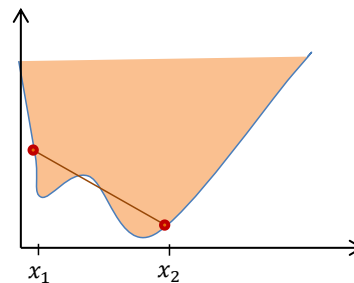


(b) A nonconvex set

Figure 2.3: The figure on the left shows convex set and the figure to the right indicates nonconvex set



(a) A convex function



(b) A nonconvex function

Figure 2.4: The figure on the left shows a convex function and the figure to the right indicates nonconvex function

In addition to the above optimization problem types, some specific optimization problems will be discussed for the completeness of the report.

Linear Programming (LP): If the objective function $f(x)$ defined by (2.8) is linear

and all the constraints are linear, then the formulation is said to be a linear programming problem. LP problems are always convex since the objective function is linear and the constraints are linear with the feasible set being convex. The LP problem is defined by

$$\min_{x \in R^n} c^T x \quad (2.9a)$$

$$\text{subject to } g_i(x) = a_i^T x - b_i = 0, \quad i = 1, \dots, m \quad (2.9b)$$

$$h_j(x) = a_j^T x - b_j \leq 0 \quad j = 1, \dots, p \quad (2.9c)$$

Quadratic Programming (QP): The mathematical optimization problem is said to be quadratic programming if the objective function $f(x)$ is quadratic and the constraints are linear. The convexity of a QP problem depends on the positive semi-definiteness of the matrix Q in (2.10a). The QP problem is defined by

$$\min_{x \in R^n} x^T Q x + c^T x \quad (2.10a)$$

$$\text{subject to } g_i(x) = a_i^T x - b_i = 0, \quad i = 1, \dots, m \quad (2.10b)$$

$$h_j(x) = a_j^T x - b_j \leq 0 \quad j = 1, \dots, p \quad (2.10c)$$

Nonlinear Programming (NLP): A general optimization problem is to select n decision variables $x_1, x_2, x_3, \dots, x_n$ from a feasible region such that a given objective function is minimized or maximized. The problem becomes a nonlinear programming problem if any of the objective or constraints are defined by nonlinear func-

tions of variables. A general nonlinear minimization problem is given by

$$\min_{x,u} f(x) \tag{2.11a}$$

$$\text{subject to } f_i(x) \leq 0, \quad i = 1, 2, \dots, m \tag{2.11b}$$

$$h_j(x) = 0, \quad j = 1, 2, \dots, p \tag{2.11c}$$

The most promising approaches to solve nonlinear optimization problems are active-set SQP (Sequential quadratic programming) methods and the interior-point methods.

SQP is one of the most widely used methods to solve nonlinear programming problems. It is an iterative procedure, where at every major iteration, a QP sub problem is solved. The solution of the QP problem is used to construct a new iterate. The construction of new iterates is done such that the series of iterates converge to a local minimum.

Interior point methods solve the problem (2.11) by applying Newtons method to a series of modified KKT (Karush–Kuhn–Tucker) conditions. In the hierarchy of convex optimization algorithms, interior point methods form the third level. The simplest or the first level includes problems whose KKT conditions are linear equations that can be solved analytically. Next level is Newtons method, which reduces the problem into a sequence of equality constrained quadratic problems. Interior point methods solve the optimization problem by reducing it to a sequence of linear equality constrained problems. Barrier method and primal-dual interior point method are frequently used interior point techniques. The goal of interior point algorithms is to approximately formulate the inequality constrained problem as an equality constrained problem where Newtons method can be used.

There are several software packages that implement nonlinear interior-point methods such as LOQO, KNITRO/DIRECT, IPOPT, BARNLP, whereas active set packages are MINOS, SNOPT, KNITRO/ACTIVE etc (Nocedal and Wright, 2006).

2.4 Optimal Control Problem

An optimal control problem is a constrained optimization problem with a dynamic system as one of its constraint. There are three parts in an optimal control problem formulation namely an objective function, constraint functions and a process model. The general optimal control problem for a class of dynamic optimization problems can be formulated as

$$\min_{x,u} J = E(x(t_f), b, t_f) + \int_{t_0}^{t_f} L(x(t), u(t), b, t) dt \quad (2.12a)$$

$$\text{subject to} \quad \dot{x} - f(x(t), u(t), b, t) = 0, \quad (2.12b)$$

$$g(x, u, b, t) = 0, \quad (2.12c)$$

$$x_0 = x(t_0), \quad (2.12d)$$

$$h(x, u, b, t) \leq 0 \quad \forall t \in [t_0, t_f] \quad (2.12e)$$

The objective function also known as performance index represents the mathematical expression that has to be either maximized or minimized. The constraint function and the plant models both determine a search domain for the optimization procedure.

The objective function given in (2.12a) is in Bolza form where $L(\cdot)$ is known as the Lagrange form and $E(\cdot)$ is known as the Mayer form. Here, t represents the time

variable, $x(t) \in \mathbb{R}^{n_x}$ and $u(t) \in \mathbb{R}^{n_u}$ denote a vector of state variables and control decision variables respectively which are going to be optimized. Parameters are the time independent decision variables denoted by $b \in \mathbb{R}^{n_b}$.

The process model given by (2.12b) and (2.12c) is also known as system equation constraints which represent an additional set of equality constraints and gives a mathematical representation of phenomena taking place in the observed system. (2.12b), (2.12c) represent ODE models and algebraic equations respectively. Both set of equations combined present a set of differential algebraic equations (DAEs).

Different equality and inequality constraints may bound the values of control and state variables to safeguard the environment, to maintain certain setpoints in the control loop etc. Equation (2.12e) is termed path constraint, which is used to determine a particular value of the state or decision variables at specific time instants. They are also known as the terminal constraints, when for instance state at the final time t_f is restricted to a particular value.

2.4.1 Methods to solve optimal control problem

There are three fundamental approaches to solve a constrained optimal control problem as presented in the form (2.12a). These methods are outlined as follows.

- 1) Dynamic Programming, Hamilton- Jacobi- Bellman (HJB)
- 2) Indirect methods, Calculus of variations, Euler Lagrange differential equation, Pontryagin's Maximum principle (PMP)
- 3) Direct methods which are based on parameterizing optimal control problem into a finite dimensional nonlinear programming problem (Kaya, 2007; Geiger, 2009).

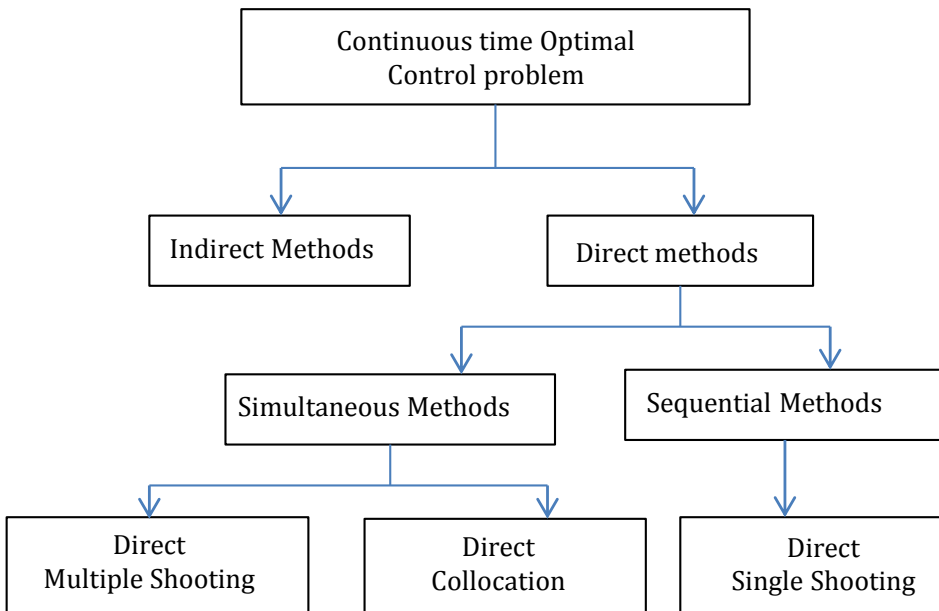


Figure 2.5: Classification of methods to solve optimal control problem

Indirect Method

Indirect methods make use of optimal conditions of the infinite optimal control problem to obtain a boundary value problem. These optimality conditions are discretized and then the boundary value problem is solved. Hence they are termed "optimize and then discretize" (Diehl et al., 2006).

Direct Method

Contrary to the indirect methods, direct methods can be used to solve the optimal control problems without the need to derive necessary conditions for optimality. The principle of direct method is to discretize the original infinite dimensional optimal control problem into a nonlinear program (NLP) with finite dimensions. Therefore these methods are also termed "discretize and then optimize" methods. The finite dimensional NLP is solved numerically using standard solution algo-

rithms. The two well known approaches within this group of methods are sequential and simultaneous approaches. Direct single shooting is a class of sequential methods while direct multiple shooting and direct collocation represents the simultaneous methods.

The following methods give an overview of the direct methods to solve the infinite dimensional optimal control problems. Direct methods can be used both for ODE and DAE models. For the sake of simplicity in understanding, the optimal control problem subject to an ODE model is briefly explained. However, this thesis is based on a system of semi-explicit differential algebraic equations (DAE) as will be seen in section 3.3, and the direct method used to solve this is described in section 4.4.

Direct Single Shooting (DSS)

This is the basic direct method for solving optimal control problems. Here the model simulation and optimization are done sequentially. It is a control parametrization method where the control is approximated and the control trajectory is parametrized using a piecewise smooth approximation, typically piecewise linear functions and the ODE or the DAE constraints (2.12b) are solved using standard numerical integration solvers. The parametrization of the control trajectory is done in the time horizon $[t_0, t_f]$, where the horizon is divided into set of gridpoints as $t_0 = 0 < t_1 < \dots < t_N = t_f$.

$$u(t) = q_i \quad \text{for } t \in [t_i, t_{i+1}], \quad i = 0, 1, \dots, N - 1 \quad (2.13)$$

In general, the control parametrization is represented as $u(t; q)$, where the term after the semicolon indicate the dependence of the discretized control parameters on every subinterval. Using numerical integration, the states are obtained as a function of several control parameters and they are denoted as $x(t; q)$. Finally,

an NLP solver is used to obtain the optimal control sequence. This is called the sequential approach for solving the optimal control problem since the ODE/DAE solvers integrate the differential equations and the NLP solver finds optimal control parameters sequentially.

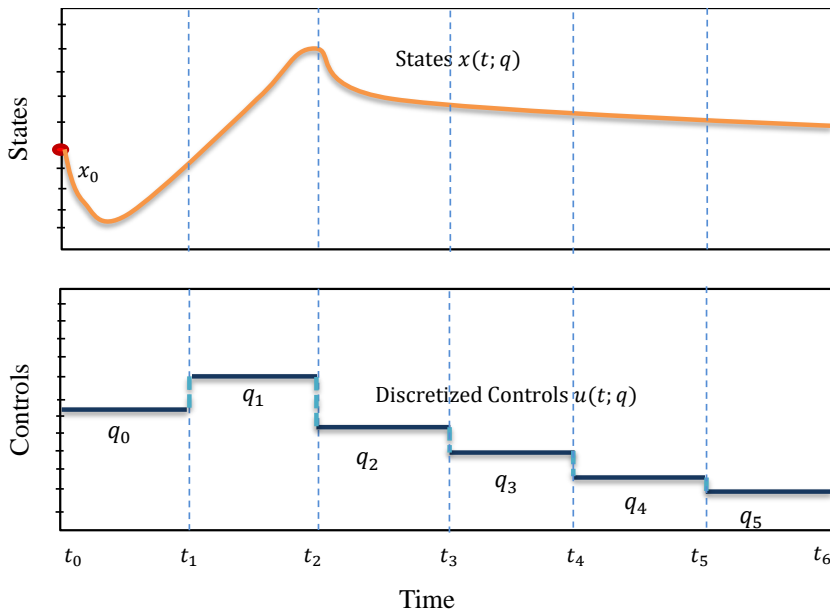


Figure 2.6: Schematic of Direct Single Shooting.

Direct Multiple Shooting (DMS)

In a direct multiple shooting method, the time interval $[t_0, t_f]$ is divided into several subintervals as presented earlier and a direct shooting method is used over each subinterval. On every subinterval the control trajectory is piecewise discretized as in (2.13) and the ODE (2.14a) is solved numerically on each interval independently, starting with an initial value (2.14b) at every gridpoint to obtain multiple

trajectory segments as shown in Figure 2.7.

$$\dot{x}_i(t) - f(x_i(t), q_i) = 0, \quad (2.14a)$$

$$x_i(t_i) = s_i \quad (2.14b)$$

Solving the above set of initial value problems (2.14) at every gridpoint, trajectory segments equal to the number of subintervals are obtained and they are denoted as $x_i(t; s_i, q_i)$.

The set of control parametrization coefficients and the initial values of state beginning at each subinterval as denoted by v remain unknown in the optimization problem and can be solved using NLP methods like Sequential Quadratic Programming (SQP) methods or Interior Point methods as described in section 2.3.1.

$$v = [s_0, q_0, s_1, q_1, \dots, s_5, q_5, s_6]$$

In order to enforce continuity of the state trajectories, additional constraints are added at the interface of each subinterval where it is desired to drive the difference between $x(t_i^-)$ and $x(t_i^+)$ to zero.

Numerical integration of the IVP over the time interval $[t_i, t_{i+1}]$ provides the state x at t_{i+1} denoted by $x_i(t_{i+1}; s_i, q_i)$ as shown in Figure 2.7. The trajectory is physically meaningful when the shooting gaps at every subintervals are closed as shown in Figure 2.7 meaning

$$x_i(t_{i+1}; s_i, q_i) - s_{i+1} = 0$$

The variables s_i and q_i indicate that the trajectory segments depend on the initial

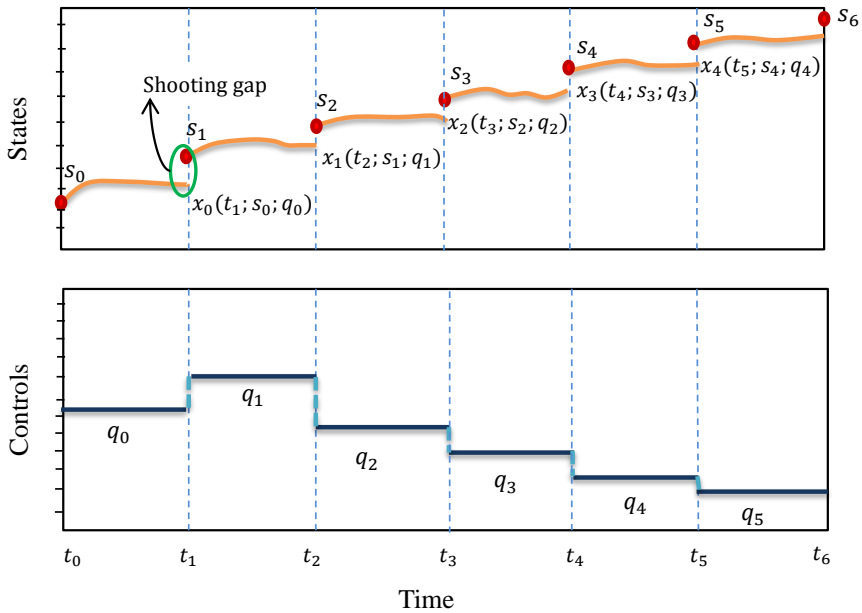


Figure 2.7: Schematic of Direct Multiple Shooting. Note the discontinuities of states in the intermediate stage of optimization, where the equality constraints are not yet satisfied.

values at the intervals and the corresponding control inputs. The shooting gaps are introduced as additional constraints in the optimization problem.

After the convergence of the NLP solver, the shooting gap constraints are fulfilled as indicated in Figure 2.8.

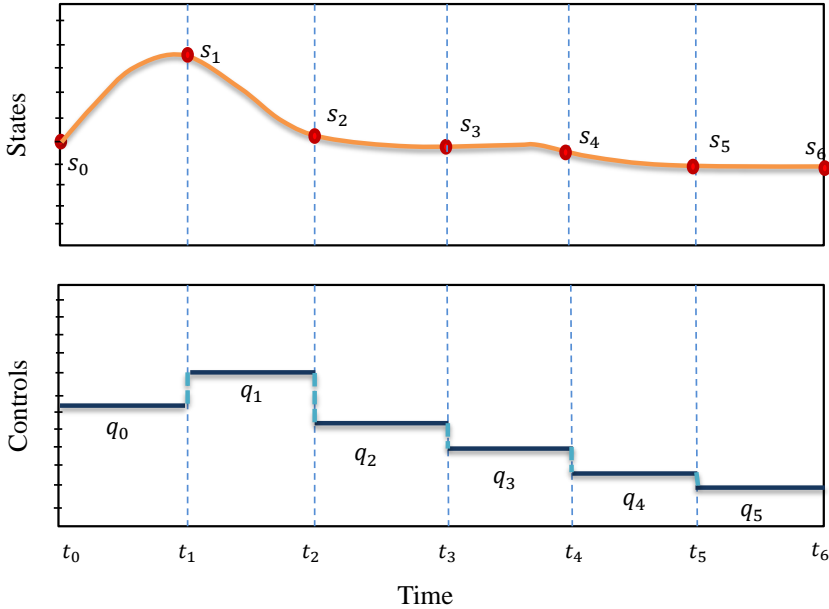


Figure 2.8: Schematic of Direct Multiple Shooting where the shooting gap constraints are satisfied after the convergence of the NLP solver (Kirches, 2010)

Direct Collocation

A direct collocation method is a state and control parametrization method where the states are approximated by cubic polynomials and the controls are expressed as piecewise linear function. The discretization is performed on a grid of N intervals with c intermediate collocation points in each interval. For example, in interval $[t_0, t_1]$ three intermediate collocation points are indicated as shown in Figure 2.9. There are different schemes to choose the collocation points depending on factors like stability and order of convergence (Magnusson and Åkesson, 2015).

$$\dot{x}(t) = f(x(t), u(t)), \quad (2.15a)$$

$$x(t_0) = x_0 \quad (2.15b)$$

An infinite ODE system is approximated using a collocation scheme, for instance the Lagrange collocation method, to obtain a polynomial approximation on each interval.

At every collocation point, the state derivative of the approximated Lagrange polynomial is evaluated and is denoted as \dot{x}_c . The value of the system dynamics (2.15a) is also computed at the collocation points $f(x_c, u(t))$. The main idea is to minimize the error between the state derivative from the system dynamics and the derivative of polynomial approximation at the collocation points as shown in Figure 2.9.

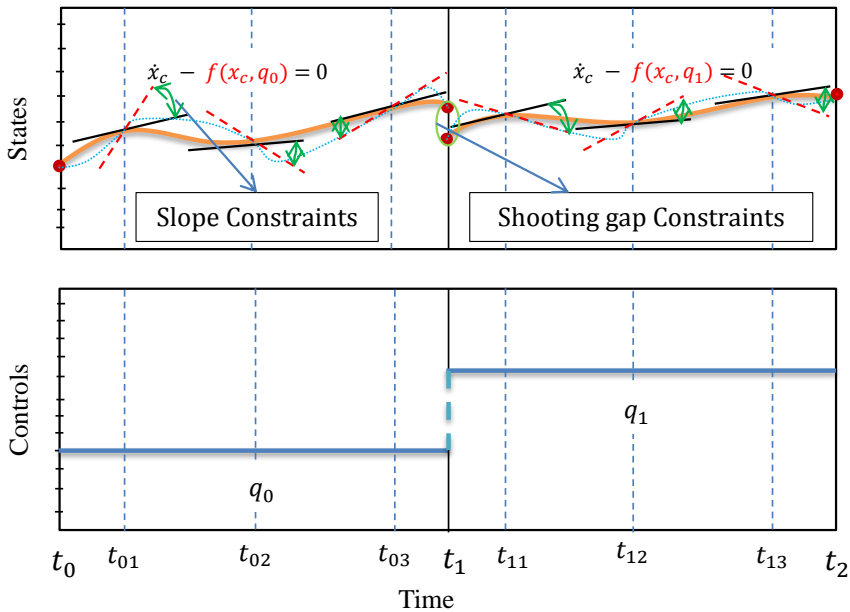


Figure 2.9: Schematic of the Direct Collocation method. The state derivative of the approximated polynomial (black line) and the value of the system dynamics (red dotted line). The polynomial approximation (orange curve) and the true solution of the system dynamics (blue dotted curve). The control input is piecewise constant over the interval $[t_k, t_{k+1}]$

The equality constraints $\dot{x}_c - f(x_c, u(t)) = 0$ are introduced into the optimization algorithm as slope constraints. In addition to this, shooting gap constraints as

discussed before for the multiple shooting method are also introduced into the optimization problem. Figure 2.9 illustrates the concept of the direct collocation method.

However, the optimal control presented above are open loop techniques which do not consider measurement noise or errors in the model. Figure 2.10 illustrates open loop control. To take the disturbances affecting the process into account, there has to be a closed loop system as shown in Figure 2.11, where the estimated states or the actual measurements are fed back to the optimizer. This is known as Model Predictive Control (MPC) or receding horizon control and is described in detail in Section 4.1.

2.4.2 Open loop vs. closed loop optimal control

In open loop optimal control, an optimal control law which is of the form $u^{opt}(t) = \omega(t; t_0, x_0)$ obtained for a specific initial value of the state $x(t_0) = x_0$, for $t \in [t_0, t_f]$.

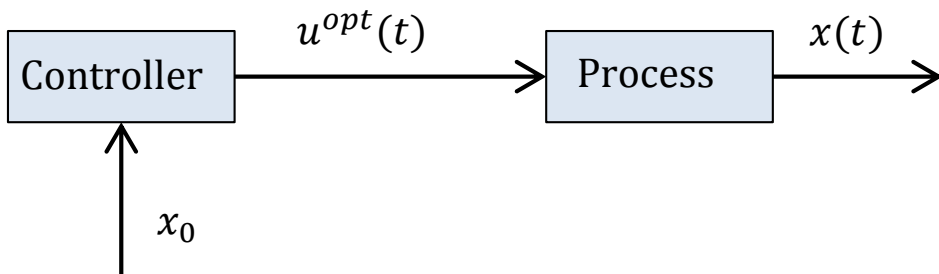


Figure 2.10: Schematic of open loop optimal control

In case of closed loop optimal control, a feedback control law is of the form $u^{opt}(t) = \omega(t; x(t))$, wherein the optimal control actions $u^{opt}(t)$ are calculated

corresponding to the current state $x(t)$, for $t \in [t_0, t_f]$.

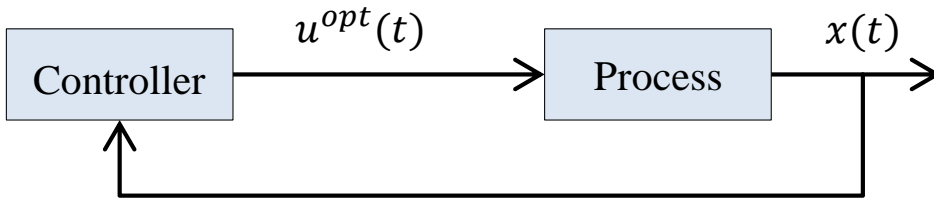


Figure 2.11: Schematic of closed loop optimal control

2.5 Software Package

In this thesis, the optimization software package CasADi (Computer algebra system for Automatic Differentiation) is used to solve Optimal Control Problems (OCP) and nonlinear model predictive control (NMPC) for a gravity separator system. CasADi solves the OCPs using direct and indirect methods both for ODE and DAE systems. Details on the installation and the syntax of CasADi can be found in the user guide by Andersson (2013).

CasADi is a tool used to efficiently compute the derivatives using automatic differentiation and is specifically developed for dynamic optimization problems. It makes use of two different representations for symbolic expressions. The two most fundamental classes are the SX and the MX symbolics. SX is the scalar representation which involves the unary or binary elementary operations. MX is the matrix representation where the elementary operations are multiple-input and multiple-output. The MX symbolic representation is more efficient for high level operations like matrix multiplication and function calls. On the contrary, the SX

representation provides faster computations by reducing the additional symbolical simplifications (Magnusson and Åkesson, 2015). The mixing of these two representations makes it possible to implement a code that is both computationally fast and generic since the SX expressions have much lower overhead per operation. Thus the SX expressions are considered for use in low level operations, for example while declaring the variables, while the MX expressions should be used for constraint function of an NLP. However, it is important to note that performing any other operation like multiplication of the two symbolic representations for instance cannot be done.

2.6 Plantwide Control by Timescale Separation

The term "plantwide control" is used in the area of process control and it is defined as the selection of the control structure (Larsson and Skogestad, 2000). It is very important to have the right control selection since the chemical or the process plants are largely complex. As seen in Figure 2.12, the layers constituting scheduling, site-wide optimization and local optimization deal with economic optimization whereas the layers supervisory control and regulatory control are concerned with setpoint tracking defined by the layer above. The two main objectives for control is the longer term economics and short term stability. Based on these two objectives, the control layer is divided into a "slow" *supervisory (economic) layer* and a "fast" *regulatory (stabilization) layer*.

The objectives of the supervisory layer is to perform advanced economic control actions (meaning to control CV_1 at their desired setpoints and to provide setpoints CV_2^{SP} to the regulatory layer) and avoid saturation of the regulatory layer. The implementation of the supervisory control is done either by using model predictive

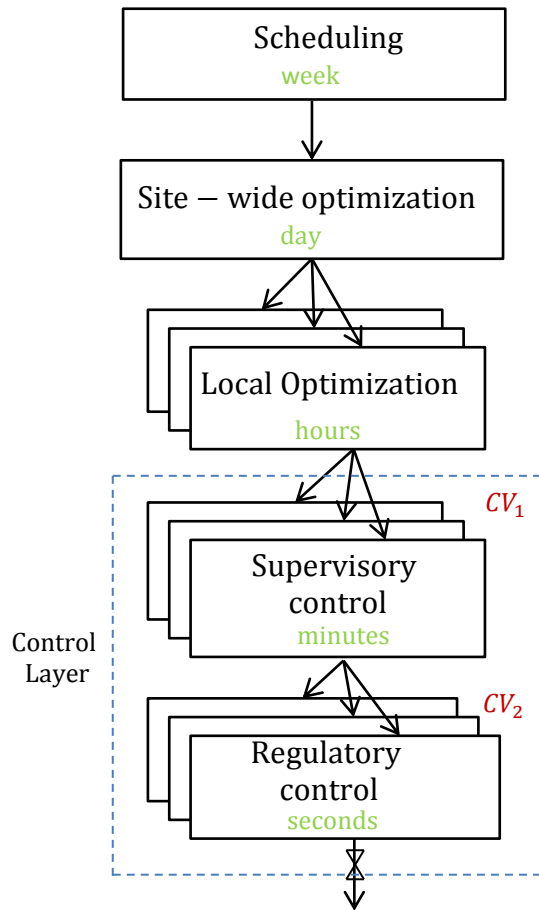


Figure 2.12: Typical control hierarchy in a chemical plant (Larsson and Skogestad, 2000).

control (MPC) or by making use of advanced control based on simpler elements like cascade control, selectors, ratio control, split range control, decoupler and valve position control. The advantage of having these simple elements is that the implementation can take place in existing basic control structure. However, it is not easily understood by the engineer who has not designed it. Hence, model based solutions (MPC) are usually preferred since the design can be easily modified.

The main objective of the regulatory layer is to keep the plant or the process at the steady state, by using simple single loop PID controllers. This layer tracks the

setpoints given by the supervisory layer (MPC).

Chapter 3

Modeling of the process

The models presented in this chapter are based upon the work of Backi and Skogstad (2017).

3.1 Assumptions

- Absence of dense - packed layer in the separator which means that the droplets will either rise or fall from the interface to disperse in the continuous phases.
- Plug flow model for both the oil and the water phase.
- Absence of coalescence and breakage of the droplets.
- *Stokes' law* is applied for the rising and the settling velocity of the water and oil droplets.
- There are neither water nor oil droplets dispersed in the gas phase and no gas droplets dispersed in the water and oil phases respectively.

3.2 Mathematical Models Description

This section describes the models for the horizontal gravity separator. The simplified schematic of the gravity separator is illustrated in Figure 3.1. The liquid enters the separator with volumetric flow rate $q_{L,in}$ and the gas flow rate into the separator is $q_{G,in}$. The separator is divided into two sections, namely the separation section and an oil chamber. The incoming liquid is a mixture of water and oil and the volume fraction of water in the liquid is termed water cut α , whereas the volume fraction of oil is termed oil cut $\beta = 1 - \alpha$. As the liquid enters the separator, some of the water droplets disperse into the bulk oil phase which is denoted by ϕ_{wo} . It follows that out of inflowing water, the fraction of water entering the water continuous phase is $\phi_{ww} = 1 - \phi_{wo}$. Similarly, some of the oil droplets will disperse into the bulk water phase which is denoted by ϕ_{ow} . Therefore, the fraction of the inflowing oil entering the continuous oil phase is $\phi_{oo} = 1 - \phi_{ow}$. The oil rich phase, flows over the weir into the oil chamber and the flow is controlled by the valve u_O . Similarly, the water rich phase is released before the weir and the flow is manipulated using valve u_W .

The objective is to control the levels of water, oil and the system pressure and to calculate the amount of dispersed phase in the bulk phase outlets. The variables in the model are classified as state variables $x = [h_L \ h_W \ p]$, manipulated variables $u = [q_{W,out} \ q_{O,out} \ q_{G,out}]$ and the disturbance variable $d = [q_{L,in} \ q_{G,in}]$.

3.2.1 Inventory Models

In this section, non-linear dynamic models are derived representing the change in water level, total liquid level, which includes the oil and the water level, as well as the system pressure. Figure 3.1 illustrates the simplified schematic of the

separation process indicating the water level, the total liquid level, the inlet and the outlet sections.

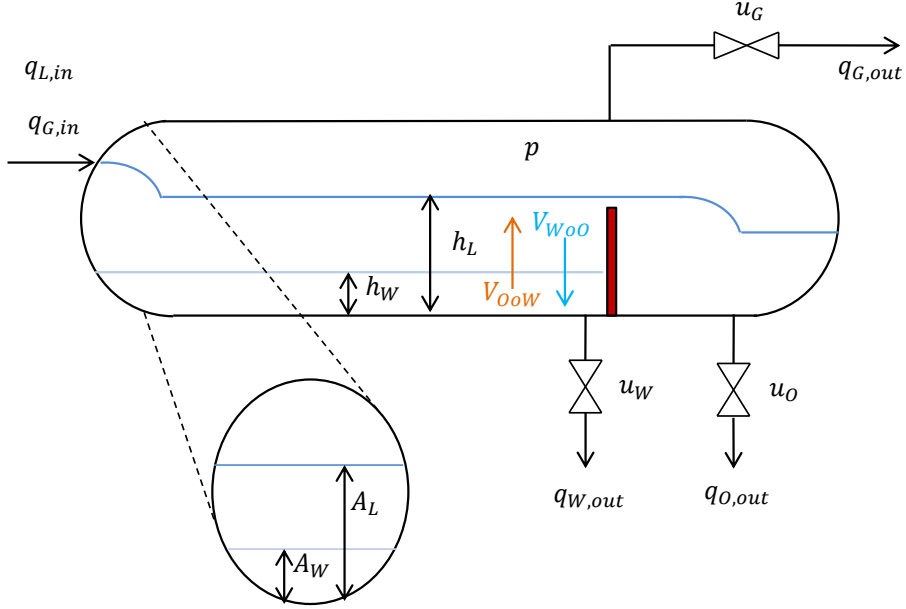


Figure 3.1: Schematic of the three phase horizontal gravity separator indicating heights of inventory.

Rate of change of water level

The rate of change of water volume is given as

$$\frac{dV_W}{dt} = q_{W,in} - q_{W,out} - \frac{V_{OoW}}{t_h^{water}} + \frac{V_{Woo}}{t_h^{oil}}. \quad (3.1)$$

The inflowing water contains dispersed oil droplets, and hence $q_{W,in}$ is given by

$$q_{W,in} = q_{L,in}(\alpha\phi_{ww} + \beta\phi_{ow}).$$

Hence, the rate of change of water volume is

$$\frac{dV_W}{dt} = q_{L,in}(\alpha\phi_{ww} + \beta\phi_{ow}) - q_{W,out} - \frac{V_{OoW}}{t_h^{water}} + \frac{V_{WoO}}{t_h^{oil}}.$$

where V_{OoW} denotes the total volume of oil droplets leaving the water phase into their continuous phase (subscript **OoW** : **O**il **o**ut of **W**ater), V_{WoO} denotes the total volume of water droplets leaving the oil phase and entering their continuous phase (subscript **WoO**: **W**ater **o**ut of **O**il) as shown in Figure 3.1, t_h^{water} and t_h^{oil} denotes the horizontal residence times of water and oil phase respectively.

The volume of water contained in the separator is $V_W = A_W L$, where L is the length of the separator, A_W is the cross sectional area of the circular segment containing water with water height h_W as indicated in Figure 3.1.

Therefore the change in area is given by

$$\frac{dA_W}{dt} = \frac{1}{L} \frac{dV_W}{dt}.$$

The cross sectional area of the circular segment filled with water A_W is calculated as

$$A_W = r^2 \left[\cos^{-1} \left(\frac{r - h_W}{r} \right) - (r - h_W) \sqrt{2rh_W - h_W^2} \right]. \quad (3.2)$$

In order to obtain the differential equation for the height of water level, the time derivative of (3.2) is given by

$$\frac{dA_W}{dt} = 2\sqrt{2rh_W - h_W^2} \frac{dh_W}{dt}. \quad (3.3)$$

The differential equation for the height of water level is

$$\frac{dh_W}{dt} = \frac{1}{2L\sqrt{2rh_W - h_W^2}} \frac{dV_W}{dt}. \quad (3.4)$$

Rate of change of total liquid level The rate of change of liquid volume in the separator is given by

$$\frac{dV_L}{dt} = q_{L,in} - q_{L,out}. \quad (3.5)$$

The total outflow is the amount of oil and water leaving the separator. Therefore the equation is rewritten as

$$\frac{dV_L}{dt} = q_{L,in} - q_{W,out} - q_{O,out}.$$

The volume of the segment in which the liquid is contained is given by $V_L = A_L L$ where A_L is the cross sectional area of the cylindrical segment containing total liquid.

The rate of change of liquid cross sectional area is

$$\frac{dA_L}{dt} = \frac{1}{L}(q_{L,in} - q_{W,out} - q_{O,out}).$$

The cross sectional area of the circular segment containing the total liquid A_L is calculated as

$$A_L = r^2 \left[\cos^{-1} \left(\frac{r - h_L}{r} \right) - (r - h_L) \sqrt{2rh_L - h_L^2} \right]. \quad (3.6)$$

In order to obtain the rate of change of liquid level, (3.6) is differentiated with time,

$$\frac{dA_L}{dt} = 2\sqrt{2rh_L - h_L^2} \frac{dh_L}{dt}. \quad (3.7)$$

Therefore, the differential equation for height of the liquid is

$$\frac{dh_L}{dt} = \frac{1}{2L\sqrt{2rh_L - h_L^2}} \frac{dV_L}{dt}. \quad (3.8)$$

Rate of change of system pressure:

As mentioned previously in Section 3.1, the gas phase of the separator is modeled by taking the ideal gas law into account. The ideal gas law is given by

$$\bar{p}V_G = n_GRT, \quad (3.9)$$

and $V_G = V_{Sep} - V_L$, where V_{Sep} is the total volume of the separator and V_G is the volume of the separator containing the gas phase both measured in cubic metres, \bar{p} is the system pressure measured in Pascal, n_G is the number of gas moles, R is the universal gas constant, T is the temperature measured in Kelvin.

The ideal gas law given by (3.9) is differentiated, yielding

$$V_G \frac{d\bar{p}}{dt} + \bar{p} \frac{dV_G}{dt} = RT \frac{dn_G}{dt} + n_G R \frac{dT}{dt}.$$

Assuming isothermal compression, the above equation reduces to

$$V_G \frac{d\bar{p}}{dt} + \bar{p} \frac{dV_G}{dt} = RT \frac{dn_G}{dt}.$$

Since the volume of the gas phase in the separator is inversely proportional to the total liquid phase volume, volume change of the gas phase is given by

$$\frac{dV_G}{dt} = -\frac{dV_L}{dt} = q_{L,in} - q_{W,out} - q_{O,out}. \quad (3.10)$$

The component balance for the gas phase is given by

$$\frac{dn_G}{dt} = -\frac{\rho_G}{M_G}(q_{G,in} - q_{G,out}). \quad (3.11)$$

Using (3.10) and (3.11) the gas pressure dynamics is written as

$$V_G \frac{d\bar{p}}{dt} = RT \frac{\rho_G}{M_G}(q_{G,in} - q_{G,out}) + \bar{p}(q_{L,in} - q_{L,out}), \quad (3.12)$$

and $\bar{p} = p \cdot 10^5$, where p is measured in bars.

The scaled form of (3.12) is thus given by

$$\frac{dp}{dt} = \left(\frac{RT}{V_G} \frac{\rho_G}{M_G}(q_{G,in} - q_{G,out}) + \left(\frac{p * 10^5}{V_G} \right)(q_{L,in} - q_{L,out}) \right) \cdot 10^{-5}. \quad (3.13)$$

3.2.2 Droplet size distribution

As stated in section 3.1, it is assumed that there is absence of coalescence or breakage of the droplets in the separator. Therefore a fixed droplet size distribution is assumed. Droplet sizes of 10 classes were defined as shown in Table 3.1. It is observed from the Figure 3.2, that the droplets are approximately normally distributed.

Table 3.1: Droplet size distribution.

Droplet size class	Droplet size μm	Number of droplets
1	50	1e8
2	100	5e8
3	150	1e9
4	200	5e9
5	250	1e10
6	300	1e10
7	350	5e9
8	400	1e9
9	450	5e8
10	500	1e8

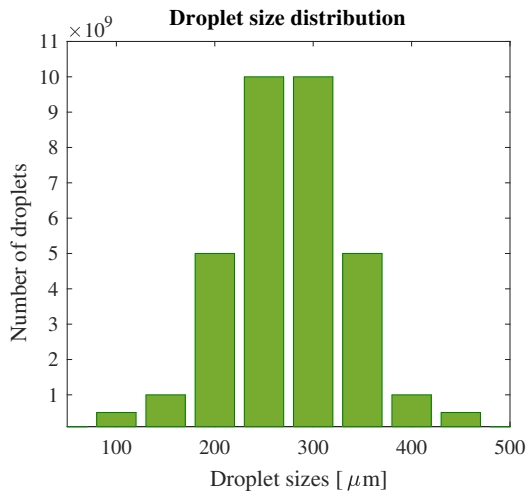


Figure 3.2: Droplet size distribution for 10 droplet classes

3.2.3 Horizontal Velocities

As the mixture enters the gravity separator, it is divided into the water continuous phase and the oil continuous phase, each of these phases containing dispersed oil and water droplets, respectively. As the continuous phase moves in the horizontal direction towards the outlet, the dispersed droplets move relative to their continuous phase. Therefore, the horizontal velocity of the droplets is assumed to be equal to the continuous phase in which they are present.

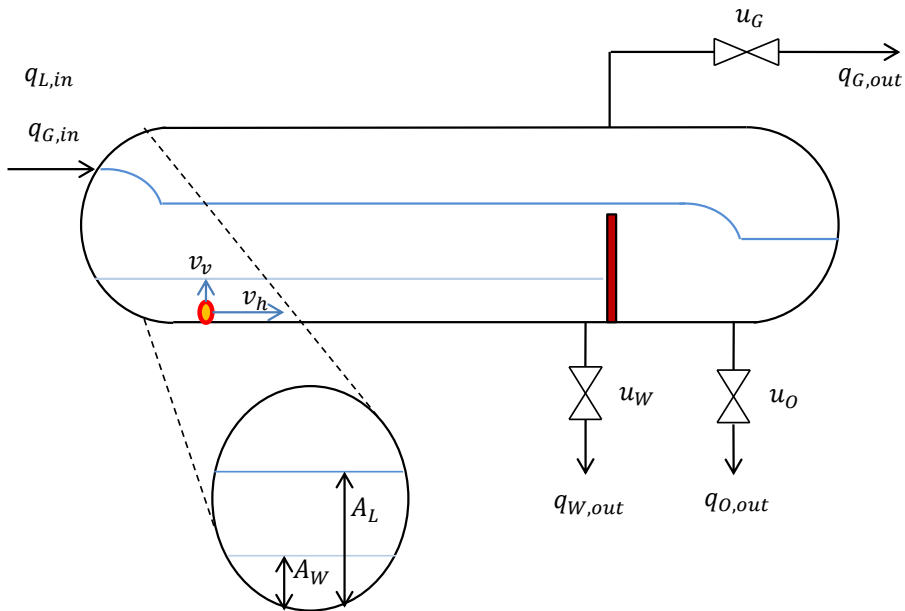


Figure 3.3: Schematic of the three phase gravity separator indicating velocity components on the droplets.

Horizontal velocity of water droplets:

$$v_h^{water} = \frac{q_{O,in}}{A_O}, \quad (3.14)$$

where $q_{O,in}$ is the flow rate of the oil into the oil continuous phase and A_O is the

area of the circular segment containing the oil phase in which the water droplets are being dispersed.

The area of the segment containing oil is obtained using (3.2) and (3.6) and is given by $A_O = A_L - A_W$.

The total flow rate of the oil $q_{O,in}$ is the sum of the oil entering the oil continuous phase and the oil dispersed into the water phase

$$q_{O,in} = q_{L,in}(\beta\phi_{oo} + \alpha\phi_{wo}).$$

Horizontal velocity of oil droplets:

$$v_h^{oil} = \frac{q_{W,in}}{A_W}, \quad (3.15)$$

where $q_{W,in}$ is the total flow rate of the water into the water continuous phase and the water dispersed into the oil phase and A_W is the area of the circular segment containing the water phase in which the oil droplets are being dispersed.

The area of the segment containing water A_W is given in (3.2).

The flow rate of liquid into the water continuous phase $q_{W,in}$ is the amount of water entering the water phase and the amount of dispersed oil entering the water phase

$$q_{W,in} = q_{L,in}(\alpha\phi_{ww} + \beta\phi_{ow}).$$

3.2.4 Vertical Velocities

The terminal velocity for the dispersed particles is given by *Stokes' Law*

$$v_v = \frac{gd_p^2(\rho_p - \rho)}{18\mu}, \quad (3.16)$$

where d_p is the diameter of the particle for every droplet class as mentioned in section 3.2.2, ρ_p is the density of the particle, ρ is the density of the continuous phase and μ is the viscosity of the continuous phase. The particles can be either water or oil droplets dispersed in their continuous phases.

Stokes' Law assumes that the particles are spherical in shape, the flow of the droplet is laminar and that the droplet movement is not hindered by the other droplets or the surface of the separator.

3.2.5 Horizontal and vertical residence time

In order to understand if the dispersed phase is being separated or is carried along with the bulk phase, it is important to understand the horizontal and the vertical residence times. The horizontal residence time for the droplets in their respective bulk phase can be defined as the amount of time the droplets spend in the separator before reaching the outlets.

The horizontal residence time is given by the equation:

$$t_h = \frac{L}{v_h^{oil,water}}. \quad (3.17)$$

The dispersed particles are considered to be separated from the continuous phase, if they rise or settle vertically to reach the interface well within the time it takes for the dispersed phase to move horizontally along the length of the separator.

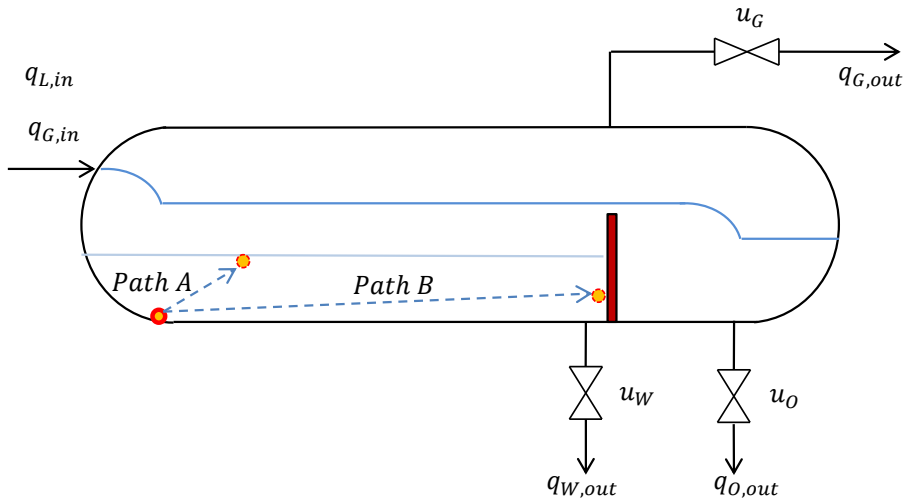


Figure 3.4: Illustration of the horizontal and the vertical residence times.

The vertical residence time is derived from the terminal velocity

$$t_v = \frac{h_{W,O}}{v_v}, \quad (3.18)$$

where $h_{W,O}$ is the level of water and oil layer ($h_O = h_L - h_W$), and v_v is the vertical velocity of the droplets (see section 3.2.1 and 3.2.4).

Comparing these residence time comparisons, the number of droplets entering their respective homogeneous phases can be calculated. Consequently, the total volume of the droplets leaving the bulk phases can be calculated. It is understood that if the time taken by the droplets to reach the interface is less than the horizontal residence time, the droplets will reach their homogeneous phase before reaching the outlet. This is illustrated by path A in Figure 3.4, and if the time taken by the droplets to reach their outlet is lesser than the vertical residence time, the droplets will be present at their outlet and is illustrated by path B in Figure 3.4.

3.2.6 Separation Efficiency

In section 3.2.5, the time taken by the droplets to reach their native phase or remain dispersed in their continuous phase was described. Based on this, the number and volume of oil droplets entering their native phase or remaining in the continuous water phase was estimated using a code scripted in MATLAB (see Appendix E.3) and similar calculations were made for the water droplets (see Appendix E.4). Separation efficiency is an important factor in order to measure the performance of separators. Separation efficiency of oil is defined as the fraction of initially dispersed oil being removed from the water phase and is given by

$$\eta_O = 1 - \frac{\text{Volume of oil droplets present in water outlet}}{\text{Initial volume of oil dispersed}}. \quad (3.19)$$

Similarly, the separation efficiency of water is defined as the fraction of initially dispersed water being removed from the oil phase and is given by

$$\eta_W = 1 - \frac{\text{Volume of water droplets present in the oil outlet}}{\text{Initial volume of water dispersed}}. \quad (3.20)$$

This means an efficiency value of 98% indicates that 2% of initially dispersed volume is contained at the outlet.

3.3 Formulation of Models in the DAE form

$$\frac{dh_W}{dt} = \frac{1}{2L\sqrt{2rh_W - h_W^2}}(q_{W,in} - q_{W,out} - \frac{V_{OoW}}{t_h^{water}} + \frac{V_{WoO}}{t_h^{oil}}) \quad (3.21a)$$

$$\frac{dh_L}{dt} = \frac{1}{2L\sqrt{2rh_L - h_L^2}}(q_{L,in} - q_{W,out} - q_{O,out}) \quad (3.21b)$$

$$\frac{dp}{dt} = \left(\frac{RT}{V_G} \frac{\rho_G}{M_G} (q_{G,in} - q_{G,out}) + \left(\frac{p * 10^5}{V_G} \right) (q_{L,in} - q_{L,out}) \right) \cdot 10^{-5} \quad (3.21c)$$

In section 3.2.5, we understood that if the vertical residence time of the droplets is less than the horizontal residence time, then the droplets will reach the oil water interface before they reach the outlet and if the vertical residence time of the droplets is greater than the horizontal residence time, then fractions of the single droplets size classes will remain in the bulk as dispersed phase.

Oil droplets in the water phase:

The system switches between these two models and since there are 10 droplet classes, the system takes the form

If $(t_h^{oil} \geq t_v^{oil}(i))$ [Oil droplets move to the interface]

$$l_p = h_W, \quad (3.22a)$$

$$V_{OoW}(i) = n(i)V(i) \quad (3.22b)$$

Else $(t_h^{oil} < t_v^{oil}(i))$ [Oil droplets remain dispersed]

$$l_p = t_h^{oil} v_v^{oil}(i), \quad (3.23a)$$

$$V_{OoW}(i) = \frac{l_p}{h_W} n(i)V(i) \quad (3.23b)$$

Water droplets in the oil phase:

If ($t_h^{water} \geq t_v^{water}(i)$) [Water droplets move to the interface]

$$l_p = h_O, \quad (3.24a)$$

$$V_{WoO}(i) = n(i)V(i) \quad (3.24b)$$

Else ($t_h^{water} < t_v^{water}(i)$) [Water droplets remain dispersed]

$$l_p = t_h^{water} v_v^{water}(i), \quad (3.25a)$$

$$V_{WoO}(i) = \frac{l_p}{h_O} n(i)V(i) \quad (3.25b)$$

where $i \in [1, \dots, 10]$, $n(i)$ is the number of particles in each particle class and $V(i)$ is the volume of particles in each particle class, $h_O = h_L - h_W$, $V_{OoW}(i)$ is the volume of oil leaving the water phase for particle class i and $V_{WoO}(i)$ is the volume of water leaving the oil phase for particle class i and l_p is the position of the droplets at the end of the separator.

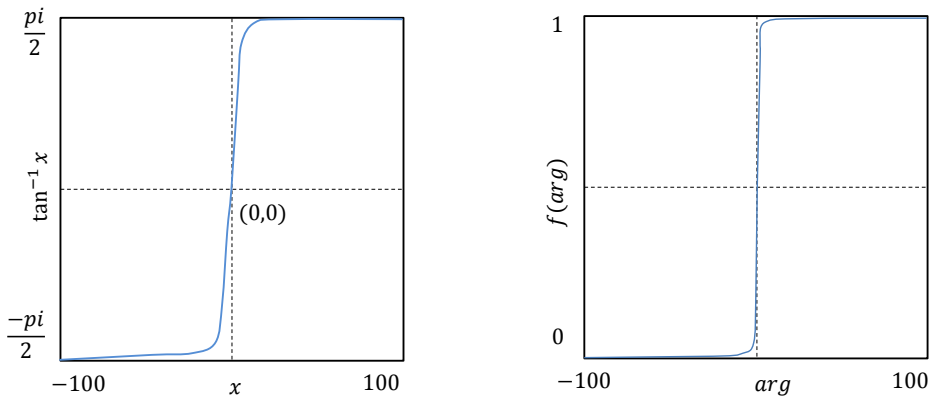
These models (3.21), (3.22), (3.23), (3.24) and (3.25) are collectively termed hybrid DAEs (Najafi and Nikoukhah, 2006) or discontinuous DAEs (Powell et al., 2016; Agrawal et al., 2003).

The if/else logic in the model is represented using greater than or equal to (\geq) and a less than or equal to (\leq) operators. The if/else formulation is presented as a set of continuous algebraic equations, in such a way that only binary solutions (0 or 1) are obtained for certain variables at the solution.

The logical expressions could be introduced into the optimization problems using mixed integer programming and solved using mixed integer programming al-

gorithms. However, combining these algorithms and introducing them into the CasADi environment seemed challenging. Also, converting dynamic optimization problems into MINLP (Mixed Integer Nonlinear Programming) problems does not seem like an easy task (Powell et al., 2016). Therefore, the variables which activate the change in the dynamics of the process should be formulated such that it will have reasonable solutions at either zero or one. This kind of formulation allows logical disjunctions in the dynamic optimization problems to be solved without making use of mixed integer programming.

Therefore, logical statements are embedded as sets of algebraic equations thus maintaining mathematical continuity. There are many operators like sigmoid function or arctan functions that can be used for this purpose. In this thesis, arctan functions are used to express the functions which are discontinuous in nature.



(a) Plot of $f(x) = \arctan(x)$ function where function value ranges from $[-\pi/2, \pi/2]$

(b) Plot of $f(arg) = \frac{\arctan(b \cdot arg) + \frac{\pi}{2}}{\pi}$ where the function value lies in $[0, 1]$ and b is the stiffness factor

Figure 3.5: Plot of arctan function on the left in the range $[-\pi/2, \pi/2]$ converted to a function on the right that takes binary values.

Therefore, the argument function which is the difference in horizontal and vertical

residence times from Equations (3.22) and (3.23) is written as

$$arg_{oil} = t_h^{oil} - t_v^{oil}(i) \quad (3.26)$$

and the argument function from Equations (3.24), (3.25) is written as

$$arg_{water} = t_h^{water} - t_v^{water}(i) \quad (3.27)$$

Since, only the binary values are necessary to represent the if/else statements, the function is rewritten as Equation (3.28) such that either 0 or 1 is obtained for the variables at that solution. The value for b in $f(arg)$ can be adjusted to make the curve smooth or stiff.

$$f(arg) = \frac{\arctan(b \cdot arg) + \frac{\pi}{2}}{\pi} \quad (3.28)$$

where arg stands for either arg_{water} or arg_{oil} .

Hence the algebraic equation is written in the form $f_i(z_i) = 0$ where z_i is the algebraic state given by the particle position l_p for 10 particle classes $i = [1, \dots, 10]$. Depending on the argument function (3.28), if it is greater than or equal to zero, one set of 'if' statement is executed and if it is less than zero, the other set of 'if' statement will be executed causing $f(x)$ to take binary values. This can therefore be written as continuous set of algebraic equations

$$f_i(z_{oil,i}) = z_{oil,i} - (f(arg_{oil})h_W) + (1 - f(arg_{oil}))t_h^{oil}v_v^{oil}(i) \quad (3.29)$$

$$f_i(z_{water,i}) = z_{water,i} - (f(arg_{water})h_O) + (1 - f(arg_{water}))t_h^{water}v_v^{water}(i) \quad (3.30)$$

In this way, the volume of oil leaving the water phase is calculated for the individual particle class and then summed up to obtain the total oil volume leaving the water phase.

$$V_{OoW}(i) = \frac{z_{oil,i}}{h_W}n(i)V(i) \quad (3.31)$$

Therefore the total volume of oil leaving the water phase is given by

$$V_{OoW} = \sum V_{OoW}(i) \quad (3.32)$$

Similarly, the volume of water leaving the oil phase is calculated for the individual particle class and then summed up to obtain the total water volume leaving the oil phase.

$$V_{WoO}(i) = \frac{z_{water,i}}{h_O}n(i)V(i) \quad (3.33)$$

and the total volume of water leaving the oil phase is given by

$$V_{WoO} = \sum V_{WoO}(i) \quad (3.34)$$

The differential equations (3.21) and the set of algebraic equations (3.29), (3.30) for gravity separator are collectively termed semi-explicit DAE (Differential Algebraic Equations)

All the models discussed above have been implemented as separate script files (see Appendix E.1 - E.4) and the plant model is solved using ode15s solver.

Model Predictive Control Theory

4.1 Introduction

Model Predictive Control (MPC) is a control strategy that calculates control inputs by solving constrained optimal control problem over a finite time horizon.

Currently, model predictive control also referred to as receding horizon control (RHC) is the most attractive feedback strategy particularly for linear processes subject to constraints (Findeisen and Allgöwer, 2002). MPC uses a plant model to predict the output trajectories. Many real systems are nonlinear in nature. In these cases, if the plant is approximated by a Linear Time Invariant (LTI) model, it is insufficient to accurately capture the nonlinearities in the model.

4.2 MPC Elements

There are three main stages which constitute the method of model predictive control strategy. First, the future values of the output variables are predicted based on the process model and current measurements. Second, the optimization problem is solved using a suitable algorithm. Finally, for a particular time instant, the manipulated variables are applied to the process which minimizes the cost or the objective function. This procedure is repeated at subsequent intervals for the entire prediction horizon.

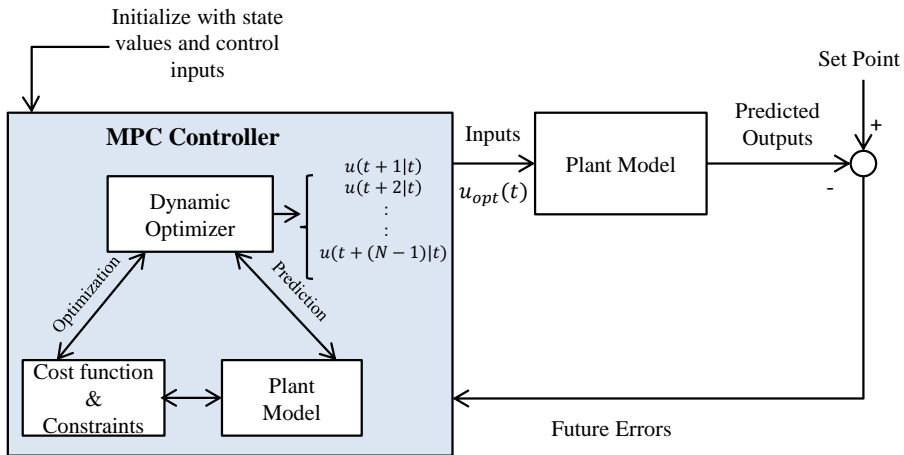


Figure 4.1: Schematic of basic MPC structure

The elements of the Model Predictive Controller are shown in Figure 4.1. The plant model in the figure represents the models of the gravity separator system. In the MPC controller mechanism, the model predicts the process output in the future and the optimization calculates the sequence of optimal control inputs by optimizing a determined set of reference criteria. The criterion, also known as the objective or cost function, can take the form of a quadratic function of the errors between the

predicted outputs and the setpoint or the reference value. These errors will be assigned a weighting factor depending on the relative importance of the manipulated variables, control variables or the constraint violations.

In short, the MPC consists of

1. At a given sampling instant i , given the initial values of the control inputs and the states, the system output for the entire prediction horizon N is obtained $y((i+k)|i)$ along with future control signals $u((i+k)|i)$.
2. Optimize the determined set of reference criterion such that the process is kept close to the setpoint
3. The first control signal calculated is applied to the plant $u(i|i)$, while the others are rejected
4. For the subsequent sampling instants, step 1 is repeated with the new value of the state obtained from the plant.

This is illustrated in Figure 4.2 where the system is simulated for a total time of t_f seconds. Along the prediction horizon T_p , the optimal control sequence is calculated for N sampling instants, that is $N = \frac{T_p}{h}$ where h is the sampling period. Assuming that the sampling period is 1 second each, the optimization problem is solved at every sampling instant and only the first part of the optimal control sequence is applied to the plant denoted by u_{opt} . Towards the end of the simulation, a set of optimal control inputs is obtained $u_{opt} = [u_0, u_1, u_2, \dots, u_{t_f-1}]$.

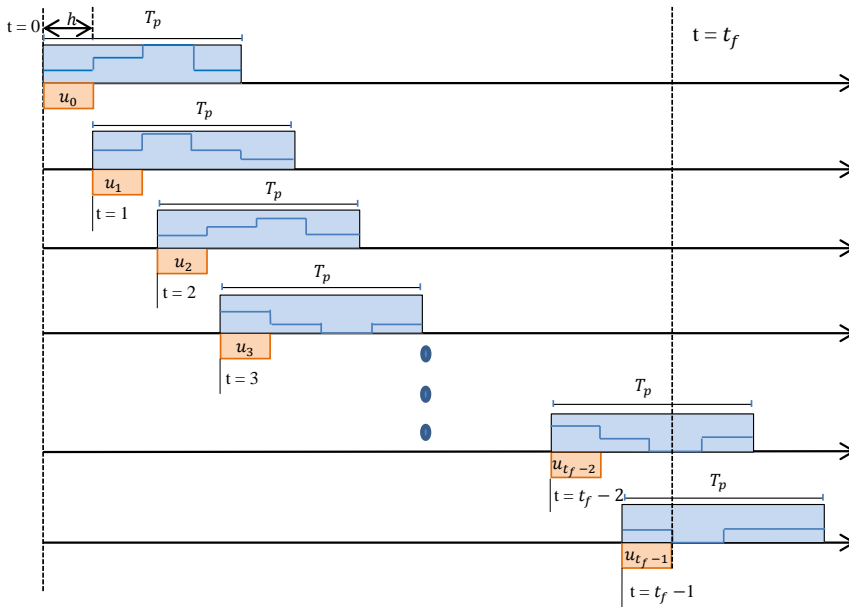


Figure 4.2: Receding horizon strategy for a total simulation time of t_f seconds

In the next section, the formulation of NMPC problem and also the detailed method to solve the optimization problem will be described.

4.3 Formulation of NMPC Optimization Problem

Figure 4.3 illustrates the basic principle of Model Predictive Control. Based on the measurements obtained at time t , the controller predicts the trajectory of the system dynamics over the prediction horizon T_p and determines the control input over the control horizon T_c such that the cost function or the performance criteria is optimized. For simplicity, the prediction horizon T_p is taken to be equal to the control horizon T_c . The time difference h between the measurements can vary, however it is assumed that the measurements take place between fixed sampling

instants h . At time $t + h$, a new measurement is obtained and by making use of it, the entire procedure of prediction and optimization is repeated to obtain a new optimal control input and the prediction horizon is moved forward.

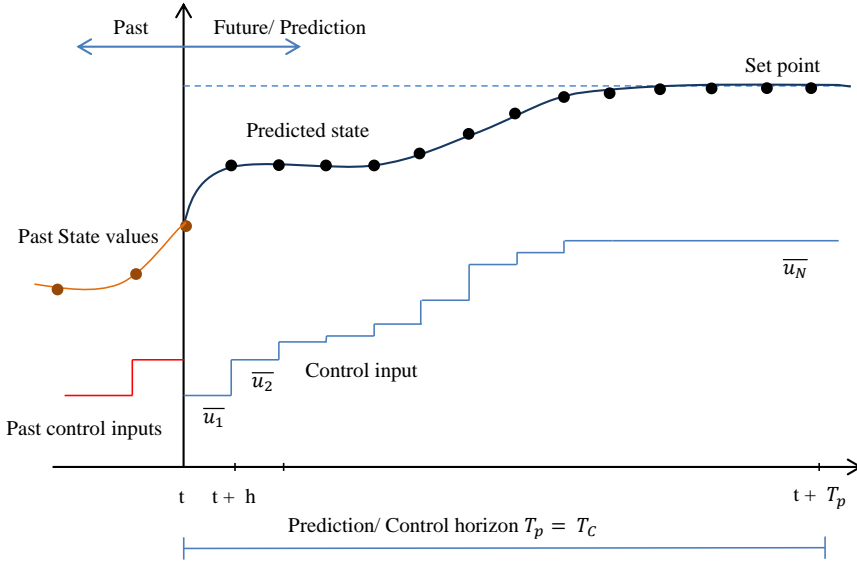


Figure 4.3: Illustration of the predictive control algorithm (Findeisen and Allgöwer, 2002)

As seen from section 3.3, the gravity separator system is modelled as a semi-explicit index-1 DAE (differential algebraic equation) of the form

$$\dot{x} = f(x, z, u, b), \quad (4.1a)$$

$$0 = g(x, z, u, b) \quad (4.1b)$$

where x is the vector of differential states, z is the vector of algebraic states, the vector of control inputs is given by u , and b are the time invariant parameters.

We assume that the state of the system x is measurable. In NMPC, the control input applied to the process in the interval $[t, t+h]$ can be obtained by the repeated

solution of finite horizon optimal control problem of the form

$$\min_{x,z,u} \int_{t_0}^{t_f} L(t, x(t), z(t), u(t), b) dt + \sum_{k=0}^{N-1} \Delta u_k^T R \Delta u_k \quad (4.2a)$$

$$\text{subject to} \quad \dot{x}(t) = f(t, x(t), z(t), u(t), b), \quad (4.2b)$$

$$0 = g(t, x(t), z(t), u(t), b), \quad (4.2c)$$

$$x(t_0) = x_0, \quad (4.2d)$$

$$x^{low} \leq x(t) \leq x^{high}, \quad (4.2e)$$

$$z^{low} \leq z(t) \leq z^{high}, \quad (4.2f)$$

$$u^{low} \leq u(t) \leq u^{high}, \quad (4.2g)$$

$$-\Delta u^{high} \leq \Delta u(t) \leq \Delta u^{high} \quad (4.2h)$$

The objective function is the typical optimal control Lagrange term as discussed in section 2.4 together with a quadratic penalty on Δu_k where R is the diagonal weighting matrix. The optimization variables are the discrete control inputs u , and the trajectories x and z .

The Lagrange term introduced in the objective function is in its continuous form and has to be discretized using the direct collocation algorithm. Whereas, the control inputs are introduced within the direct collocation algorithm as discrete variables. Therefore, in the above formulation (4.2a) the continuous Lagrange term is augmented with the differences of the piecewise constant control inputs, Δu which are discretized with N sampling instants in the time interval $[t_0 \ t_f]$.

4.4 Direct Collocation Method to solve NMPC Optimization Problem

Section 2.4.1 briefly described the methods used in general to solve the optimal control problem. Among the direct methods seen earlier, direct collocation method is employed in this thesis to solve the optimal control problem (4.7b). The basic idea is to discretize the differential equations and transcribe the optimization problem into a finite dimensional NLP, which is later solved by IPOPT. The third order direct collocation gives a polynomial approximation of (4.1) as shown in Figure 4.4.

The optimization time horizon is divided into N elements such that each interval is in $[t_k \ t_{k+1}]$ where $k \in [0, \dots, N - 1]$. The trajectories x and z are approximated using low order polynomials in the interval $[t_k \ t_{k+1}]$ known as collocation polynomials. The collocation polynomials are formulated by choosing K collocation points and these are kept same for all elements. In this thesis, Lagrange interpolating polynomials are being used as the collocation polynomials given by

$$P_{k,i}(t) = \prod_{j=0, j \neq i}^K \frac{t - t_{k,j}}{t_{k,i} - t_{k,j}} \quad (4.3)$$

with the property

$$P_{k,i}(t_{k,l}) = \begin{cases} 1 & \text{if } l = i \\ 0 & \text{otherwise} \end{cases}$$

Let $t_{k,i}$ be the collocation points in the interval $[t_k \ t_{k+1}]$ and $i \in [0, \dots, K]$, where $(x_{k,i} \ z_{k,i} \ u_{k,i})$ is the simplified notation of $(x(t_{k,i}) \ z(t_{k,i}) \ u(t_{k,i}))$. The collo-

cation polynomials are then formulated by interpolating the values $(x_{k,i}, z_{k,i})$ in that time interval. The interpolation yields

$$x(x_k, t) = \sum_{i=0}^K x_{k,i} P_{k,i}(t) \quad (4.4a)$$

$$z(z_k, t) = \sum_{i=1}^K z_{k,i} P_{k,i}(t) \quad (4.4b)$$

where $x_{k,i}$ are the parameters and $P_{k,i}$ are the Lagrange polynomials. Additionally, the degrees of freedom (DOF) for the differential states is $K + 1$ whereas it is K for the algebraic states. The reason for an additional DOF in case of the differential states is to ensure continuity so that the shooting gaps are closed. Algebraic states can be discontinuous and therefore do not require an additional DOF.

In order to ensure integration accuracy, Radau collocation points are used. The Radau collocation points are obtained from Radau quadrature, which is similar to the Gaussian quadrature. However, one collocation point is clearly defined at one end of the time interval, rather than having all points completely inside the time interval like the Legendre collocation points (Powell et al., 2016).

The control inputs are assumed to be piecewise constant in the interval $[t_k, t_{k+1}]$ and given by

$$u(t) = u_k \quad (4.5)$$

The polynomials for \dot{x} are obtained by differentiating the polynomials for x

$$\dot{x}_{k,i} = \frac{1}{(t_{k+1} - t_k)} \sum_{r=0}^K \dot{P}_{k,l}(t_{k,i}) x_{k,r} = \frac{1}{(t_{k+1} - t_k)} \sum_{r=0}^K \alpha_{r,i} x_{k,r} \quad (4.6)$$

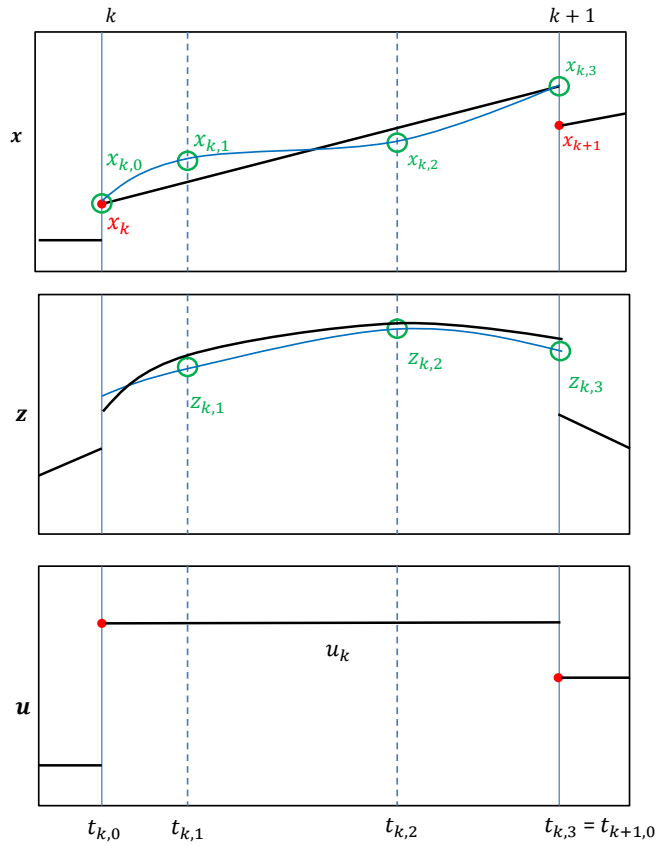


Figure 4.4: Schematic representation of third order direct collocation in the single sampling interval $[t_k, t_{k+1}]$. The control input is piecewise constant in the interval $[t_k, t_{k+1}]$.

4.5 Finite Dimensional NLP

The variables in the transcribed NLP are chosen such that it contains the variable values at all collocation points $\dot{x}_{k,i}, x_{k,i}, z_{k,i}$, the control input values u_k the initial

value of the state $x_{k,0}$ and the parameters b . The transformation results in the NLP

$$\min_{x_{k,i}, z_{k,i}, u_k} \sum_{k=0}^{N-1} \sum_{i=0}^K \omega_i L(t_{k,i}, x_{k,i}, z_{k,i}, u_k, b) + \sum_{k=0}^{N-1} \Delta u_k^T R \Delta u_k \quad (4.7a)$$

$$\text{subject to } \dot{x}_{k,i} - f(t_{k,i}, x_{k,i}, z_{k,i}, u_k, b) = 0, \quad (4.7b)$$

$$0 = g(t_{k,i}, x_{k,i}, z_{k,i}, u_k, b), \quad (4.7c)$$

$$x_{1,0} = x_0, \quad (4.7d)$$

$$x^{low} \leq x_{k,i} \leq x^{high} \quad (4.7e)$$

$$z^{low} \leq z_{k,i} \leq z^{high} \quad (4.7f)$$

$$u^{low} \leq u_{k,i} \leq u^{high} \quad (4.7g)$$

$$-\Delta u^{high} \leq \Delta u_k \leq \Delta u^{high} \quad (4.7h)$$

$$0 = x_{k,K} - x_{k+1,0} \quad (4.7i)$$

$$\dot{x}_{k,i} = \frac{1}{(t_{k+1} - t_k)} \sum_{r=0}^K \alpha_{r,i} x_{k,r} \quad (4.7j)$$

The Lagrange term in (4.2a), which is in the continuous form is transcribed into discrete form in (4.7a) using Gauss quadrature within each element such that the integral is approximated by a summation using quadrature weights ω_i (see Appendix B). The states need to be continuous with respect to time and hence the shooting gap constraint or the continuity constraint is introduced given by (4.7i). Equation (4.7j) is introduced to specify the relation between \dot{x} and x . The weights $\alpha_{r,i}$ are the result of differentiating Lagrange polynomials for the states.

4.6 Continuous Logic in an NMPC problem

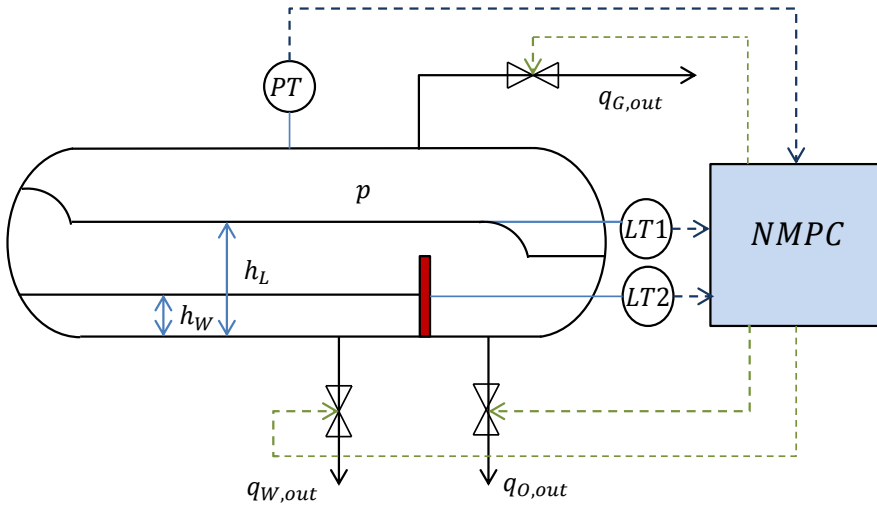


Figure 4.5: Control structure for a gravity separator

The MPC problem attempts to minimize the deviations from the reference value for the water level h_W , liquid level h_L and system pressure p , without making large control moves. Therefore, a quadratic performance index is employed, where the deviations are squared at the end of each time interval, and are weighted accordingly, and are finally summed up to produce an objective function to be minimized. This generates an optimization problem which is subject to system models and the constraints.

As discussed in section 2.6, the MPC provides setpoints to the regulatory layer (PI controllers), and this layer manipulates the control valves to keep the plant at steady state values. However, in this thesis the regulatory layer is not considered, and the NMPC is directly used to manipulate the control valves using the optimal control inputs obtained after solving the optimization problem.

Simulation Results and Discussion

In this thesis, the dynamic optimization problem is transformed into a discrete NLP problem using direct collocation method where radau collocation points are used. This problem has been implemented in CasADi using MATLAB as the modeling environment. Once the NLP problem is created, it is solved using the IPOPT plugin, which is an open source primal- dual interior point method incorporated in CasADi installations (Andersson, 2013). At every iteration, the first value of the obtained optimal input is sent to the plant model. The plant model has been implemented in MATLAB as separate m-files or script files. From the plant model, the states computed are sent to the optimizer and these values will be the starting point for the next iteration. The simulation results in this section are described based on the Gullfaks-A production rates (Laleh et al., 2013; Backi and Skogestad, 2017). The simulation parameters are presented in Appendix A.

As it has been discussed in section 3.3, the models are expressed as semi-explicit DAEs, described by the differential equations and two sets of algebraic equations each for the oil droplets dispersed in the continuous water phase and the water

droplets dispersed in the continuous oil phase. However, while implementing both sets of algebraic equations in the CasADi environment, an error message was given by the IPOPT plugin "Converged to a point of local infeasibility. Problem may be infeasible". Since, the main interest was to minimize the total amount of oil present in the water phase, simulations were carried out using the differential equations 3.21 and considering only the algebraic equations 3.29 which represent the oil droplets present in the water phase, therefore discarding the other set of algebraic equations that represent the water droplets in the oil phase 3.30.

Table 5.1: Setpoint values for the controlled variables for different cases

Parameters	Values	Values
	[Case 1, III, IV]	[Case II]
h_W^{SP}	1.2	1.2, 1.4, 1.6, 1.8
h_L^{SP}	2.5	2.5
p^{SP}	68.7	68.7

Case I: Comparative study of the effect of noise on the system control behavior

In real processes, the controller has to handle noisy measurements. It is important that the controller should not be very sensitive to the noise, because there is a possibility that the control valves will be subjected to continuous wear and tear and can be worn out easily. Therefore, a random noise was added to the states both of high order magnitude and low order magnitude to understand the effect of noise on the system behavior. The NMPC formulation for this case is written as follows, and the setpoint values are displayed in Table 5.1

$$\begin{aligned}
\min_{q_{W,out}, q_{O,out}, q_{G,out}} & \sum_{k=0}^{N-1} (h_W(k) - h_W^{SP})^2 + (h_L(k) - h_L^{SP})^2 + (p(k) - p^{SP})^2 \\
& + \sum_{k=0}^{N-1} (q_{W,out}(k) - q_{W,out}(k-1))^2 \\
& + \sum_{k=0}^{N-1} (q_{O,out}(k) - q_{O,out}(k-1))^2 \\
& + \sum_{k=0}^{N-1} (q_{G,out}(k) - q_{G,out}(k-1))^2 \tag{5.1a}
\end{aligned}$$

$$\text{Subject to } (3.21), \tag{5.1b}$$

$$(3.29), \tag{5.1c}$$

$$0.9 \leq h_W \leq 1.9, \tag{5.1d}$$

$$2.2 \leq h_L \leq 3.3, \tag{5.1e}$$

$$50 \leq p \leq 100, \tag{5.1f}$$

$$-0.05 \leq q_{W,out}(k) - q_{W,out}(k-1) \leq 0.05, \tag{5.1g}$$

$$-0.05 \leq q_{O,out}(k) - q_{O,out}(k-1) \leq 0.05, \tag{5.1h}$$

$$-0.05 \leq q_{G,out}(k) - q_{G,out}(k-1) \leq 0.05 \tag{5.1i}$$

$$h_W(0), h_L(0), p(0) \text{ given} \tag{5.1j}$$

$$q_{W,out}(-1), q_{O,out}(-1), q_{G,out}(-1) \text{ given} \tag{5.1k}$$

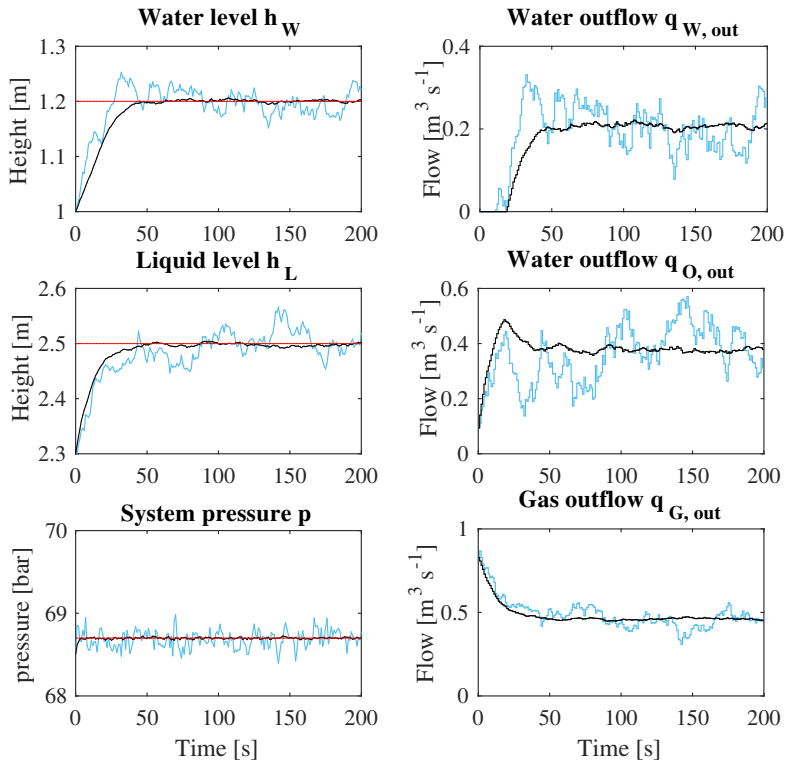


Figure 5.1: Right: Manipulated variables for low magnitude noise addition (black) and high magnitude noise addition (blue) Left: State variables for low magnitude noise addition (black) and high magnitude noise addition (blue) and the setpoint values (red)

For the purpose of generating noise, `randn` function from MATLAB toolbox, which draws a random scalar from a standard normal distribution is used. In Figure 5.1, a low magnitude measurement noise is added to the states, which is of the order of 10^{-4} on the water and the liquid level, and 10^{-3} on the system pressure. Additionally, a high measurement noise is added to the states, which is of the order of 10^{-3} on water level and the liquid level, and 10^{-2} on the system pressure. It is observed that the controller is very sensitive to increase in measurement noise thereby causing the manipulated variables to be noisy as well.

In this thesis, all simulations for case II, case III and case IV have been carried out for low magnitude measurement noise.

CASE II: Setpoint Tracking

The objective is to track setpoints, where the output has to converge to a given reference value as close as possible while fulfilling the constraints. Here, the water level h_W , total liquid level h_L and the system pressure p are considered for tracking. The NMPC formulation for this case is given by (5.1), where the setpoints for water, liquid levels and the system pressure are displayed in Table 5.1.

In Figure 5.2, the plots on the left hand side show the variables to be controlled or the state variables h_W , h_L and p . The three plots on the right hand side show their corresponding manipulated variables $q_{W,out}$, $q_{O,out}$ and $q_{G,out}$. A series of stepwise setpoint changes were employed to the water level. Setpoint changes of +0.2 m were given at every 200 s of simulation time, with the initial water level of 1 m. It is observed from the Figure that the controlled variables are kept at their set-point values by the NMPC controllers where the set-point values are displayed in red.

It is also seen that, when a step is given to the water level at every 200 s, the controller tries to close the control valve, therefore the water outflow drops to zero. Ideally, there should be a gradual drop in water outflow, since the rate of change of flow constraint has been implemented in the code (see Appendix E.5, line). However, there is a sudden fall to zero at the point where the step change in level is introduced, which means that the rate of change of flow constraint is violated at those time instants. Whereas, the implementation of the rate of change constraints work well at those time instants where no step change is introduced. Similar observation is made for the outlet oil flow rate, where there is a rise in the oil outflow which is as expected, however the rise is not gradual where the step

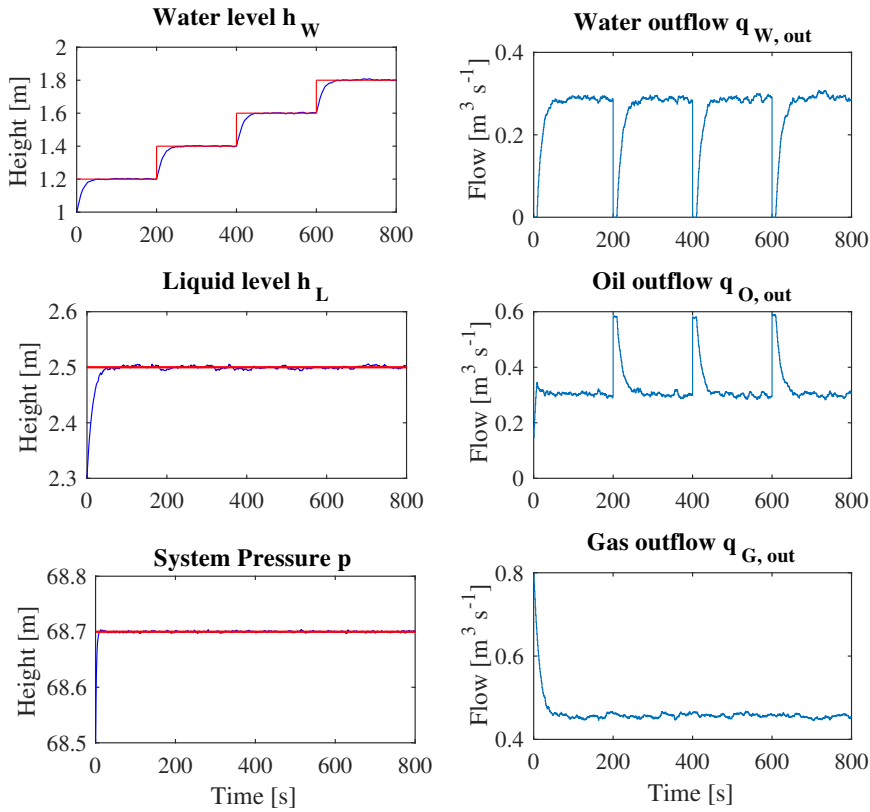


Figure 5.2: Setpoint tracking of gravity separator controlled with NMPC. State variables h_W , h_L and p on the left with their respective manipulated variables $q_{W,out}$, $q_{O,out}$ and $q_{G,out}$ on the right

change in the water level is given, therefore violating the constraints (5.1h) at those time instants (see Appendix D). In this formulation, equal weights were given to both the states as well as the change in manipulated variables. The reason being that it was equally important to keep the states at their reference values, and also to avoid the violation of control move constraints.

CASE III: Disturbance Rejection

As it has been discussed in chapter 3, the separator levels and the system pressure are often disturbed by varying inflow rate into the separator system. Usually, the separator is disturbed by the liquid inflow $q_{L,in}$ and the gas inflow $q_{G,in}$. The optimization problem given by (5.1) is solved and the aim of the controller is to maintain the levels and the system pressure at their desired reference values, (see Table 5.1) despite the disturbances due to the inflows to the system.

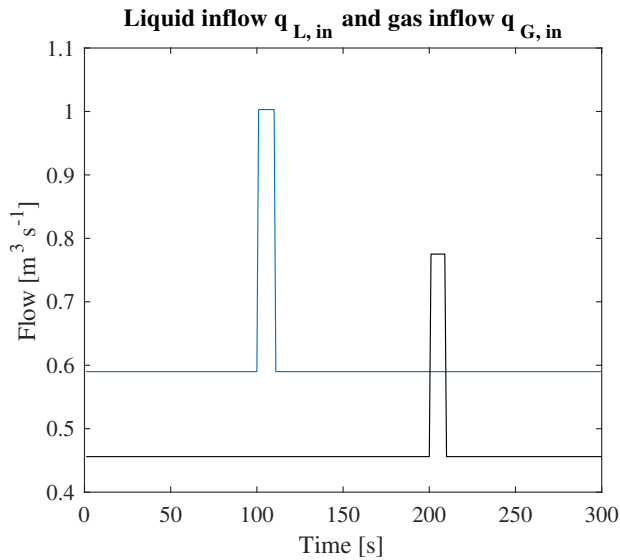


Figure 5.3: Pulse disturbance in liquid inflow $q_{L,in}$ (blue) and disturbance in gas inflow $q_{G,in}$ (black)

Figure 5.4 shows the response of the levels and system pressure of the gravity separator subject to pulse disturbances in the liquid inflow rate $q_{L,in}$ and the gas inflow rate $q_{G,in}$ one at a time. At $t = 100$ s, a sudden increase in liquid inflow rate is applied, therefore increasing from the steady state value $q_{L,in} = 0.59 m^3 s^{-1}$ to $1.003 m^3 s^{-1}$ and then at $t = 110$ s back to the initial value as shown in the Figure 5.3. It can be seen that, with the sudden increase in the liquid inflow, the water

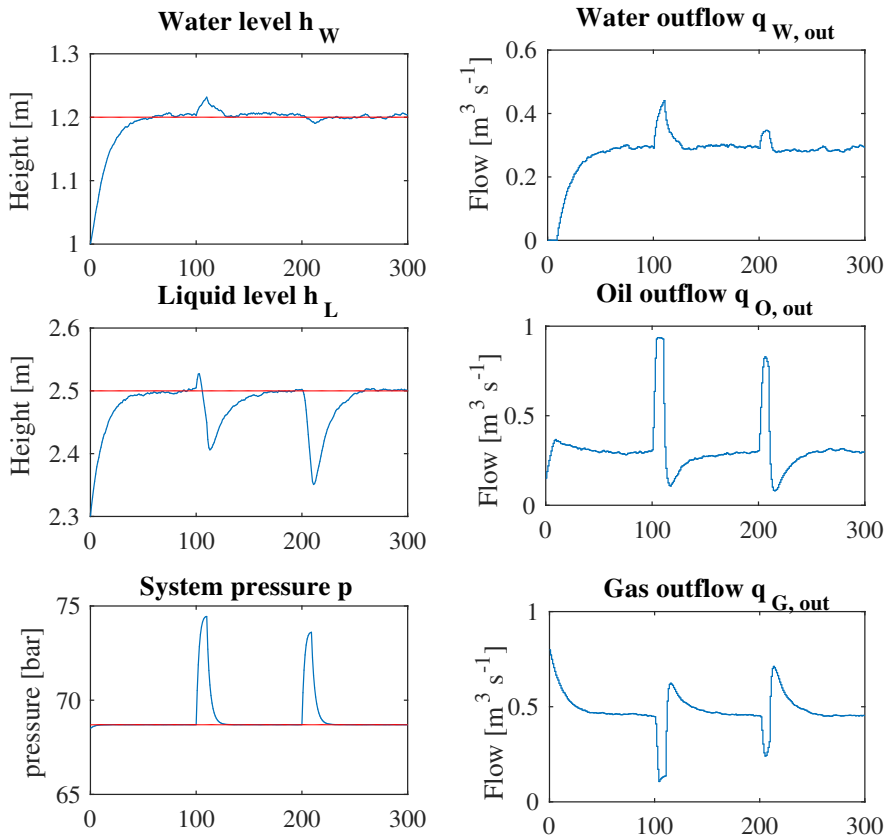


Figure 5.4: Disturbance rejection when there is sudden increase in inflow liquid $q_{L,in}$ and inflow gas $q_{G,in}$. Left: h_W , h_L and p , and the dotted red line indicates the reference values. Right: Manipulated variables $q_{W,out}$, $q_{O,out}$ and $q_{G,out}$

level and the liquid level increases, thereby increasing the pressure of the system. The controller tries to flush the accumulated liquid out of the system to restore the levels to their reference values.

Similarly, a pulse disturbance in the gas inflow rate is given to the system. At $t = 200$ s, a sudden increase in gas inflow rate is applied, therefore increasing from the steady state value $q_{G,in} = 0.456 \text{ m}^3 \text{ s}^{-1}$ to $0.775 \text{ m}^3 \text{ s}^{-1}$ and then at $t = 210$ s back to the initial value as shown in the Figure 5.3. It can be observed

that the disturbance in $q_{G,in}$ results in abrupt increase in the system pressure. The controller starts flushing more gas out of the system to bring back the pressure to normal level. It can be seen that other manipulated variables change appropriately to bring back water and liquid levels back to the reference values.

CASE IV: Maximize the volume of oil entering its homogeneous phase (V_{OoW})

In this case, the objective is to maximize the total volume of oil entering its native (mother) phase, and obtain the optimal value of the water level while satisfying the constraints. It is desired to control the water level in the range given by (5.2d). Therefore, the objective function does not include tracking for the water level since maximizing V_{OoW} and tracking of the water level to the desired setpoint cannot be done simultaneously. Since the goal is to minimize the objective function while maximizing V_{OoW} , this term is preceded with a minus sign in the objective function. Additionally, because it is very important to maximize V_{OoW} , a higher weight was assigned to this term in comparison to the weights on state variables and the control moves. The NMPC in this case is formulated as follows.

$$\begin{aligned}
 \min_{q_{W,out}, q_{O,out}, q_{G,out}} \quad & \sum_{k=0}^{N-1} 100(h_L(k) - h_L^{SP})^2 + 100(p(k) - p^{SP})^2 \\
 & + \sum_{k=0}^{N-1} 100(q_{W,out}(k) - q_{W,out}(k-1))^2 \\
 & + \sum_{k=0}^{N-1} 100(q_{O,out}(k) - q_{O,out}(k-1))^2 \\
 & + \sum_{k=0}^{N-1} 100(q_{G,out}(k) - q_{G,out}(k-1))^2 \\
 & - \sum_{k=0}^{N-1} 1000(V_{OoW}(k))^2 \tag{5.2a}
 \end{aligned}$$

$$\text{Subject to } (3.21), \tag{5.2b}$$

$$(3.29), \tag{5.2c}$$

$$0.9 \leq h_W \leq 2.2, \tag{5.2d}$$

$$2.2 \leq h_L \leq 3.3, \tag{5.2e}$$

$$50 \leq p \leq 100, \tag{5.2f}$$

$$-0.05 \leq q_{W,out}(k) - q_{W,out}(k-1) \leq 0.05, \tag{5.2g}$$

$$-0.05 \leq q_{O,out}(k) - q_{O,out}(k-1) \leq 0.05, \tag{5.2h}$$

$$-0.05 \leq q_{G,out}(k) - q_{G,out}(k-1) \leq 0.05 \tag{5.2i}$$

$$h_L(0), p(0), q_{W,out}(-1), q_{O,out}(-1), q_{G,out}(-1) \text{ given} \tag{5.2j}$$

In Figure 5.5, the two plots on the top show the state variables h_L , p and the plots on the bottom show the manipulated variables $q_{O,out}$, $q_{G,out}$. It is seen from the plots of the controlled variables, that the NMPC is able to keep them at their desired reference values, (see Table 5.1). It can be observed from the oil outflow

plot, that there are a few drops in the outlet flow rate thus violating the constraints 5.2h. The reason is that, while solving the above optimization problem, IPOPT gives a message "Maximum number of iterations reached" at those points which means that the solver could not find the optimal control input at those time instants (see Appendix D, Figure D.3). Similar observation is also made in Figure 5.6.

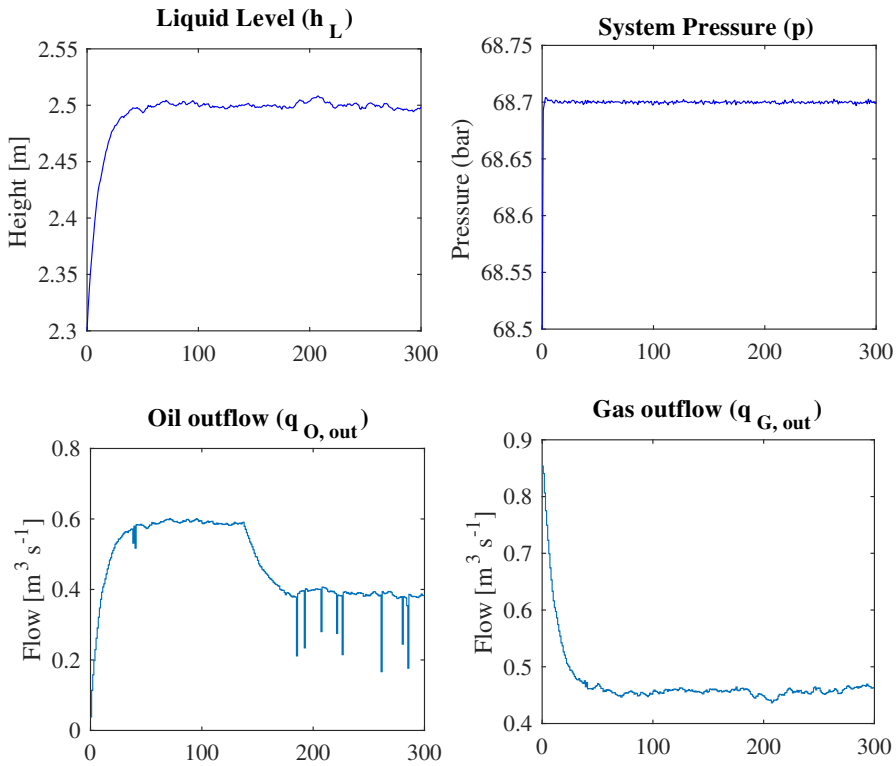


Figure 5.5: Controlled variables h_L and p on top with their respective manipulated variables $q_{O, out}$ and $q_{G, out}$ on the bottom

In Figure 5.6, the plot on top shows the water level h_W , which starts at an initial value of 1 m, and stabilizes at 2 m. The red dotted lines indicate the upper and lower bounds given by (5.2d). It is observed from the Figure that the controlled

variable h_W is kept within its bounds and stabilizes at an optimal value. It can be seen that the optimal value for the level of water in order to maximize the amount of oil leaving the water phase and entering its continuous phase V_{OoW} is at 2 m.

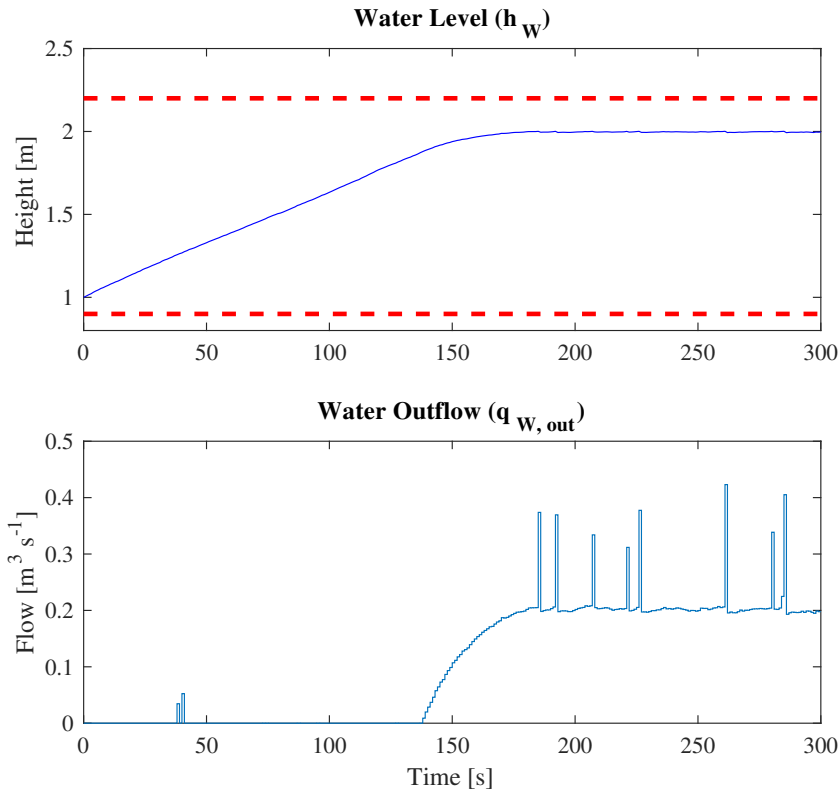


Figure 5.6: Top: Controlled variable h_W indicated in (blue) and the bounds on h_W in dotted red. Below: Manipulated variable $q_{W,out}$

In order to investigate the optimal water level, the total volume of oil at the water outlet, efficiency of oil removal from the water phase as well as the horizontal and the vertical residence times were analyzed. The lower bound on the water level is chosen to be 0.9 m, while the upper bound is set to 2.2 m. In order to investigate the optimal water level, four different water levels are considered within the given

bounds.

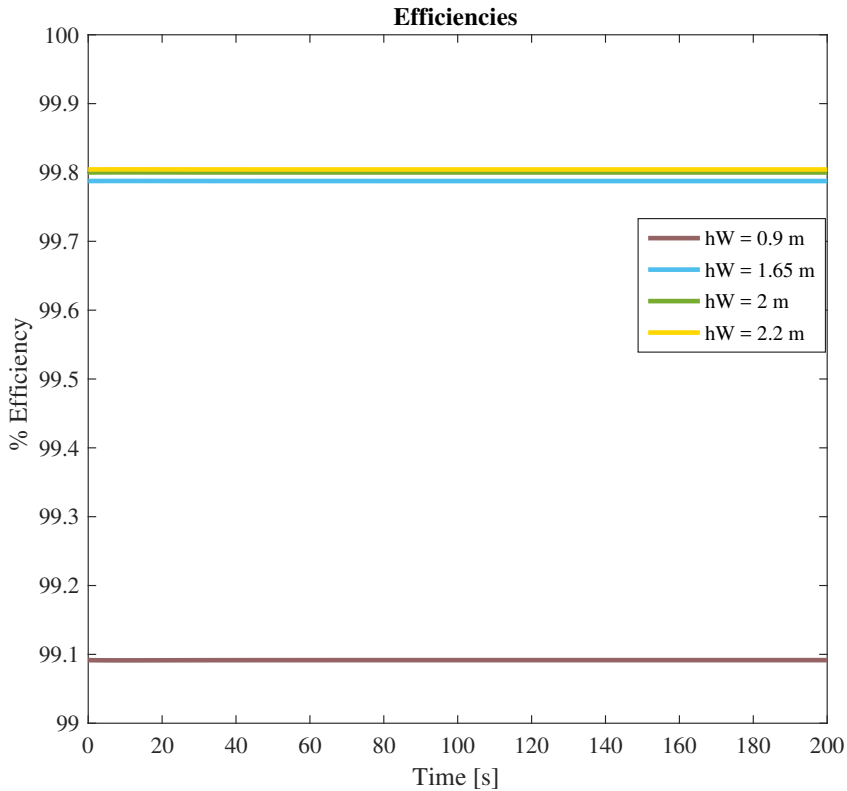


Figure 5.7: Efficiencies of oil removal from the water phase for different water levels

From Figure 5.7, it can be observed that at lower water levels the efficiency of oil removed from the water phase is smaller, whilst it increases with an increase in the water level. It can be noticed that the efficiency of oil removal for a water level greater than 2 m is almost constant at around 99.8%.

This can also be observed from the oil droplet distribution at the water outlet, which is demonstrated in the following Figures 5.8 - 5.11.

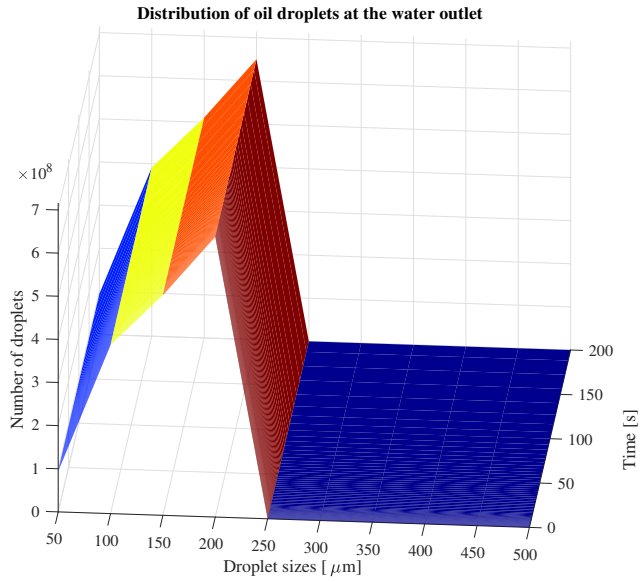


Figure 5.8: Distribution of oil droplets at the water outlet for water level = 0.9 m

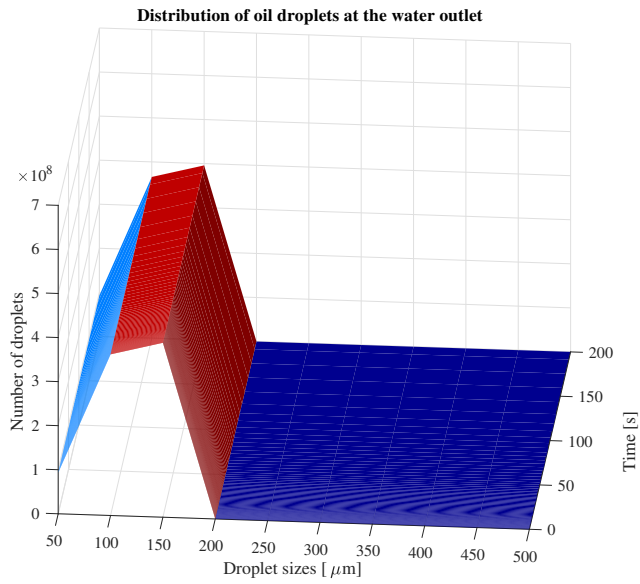


Figure 5.9: Distribution of oil droplets at the water outlet for water level = 1.65 m

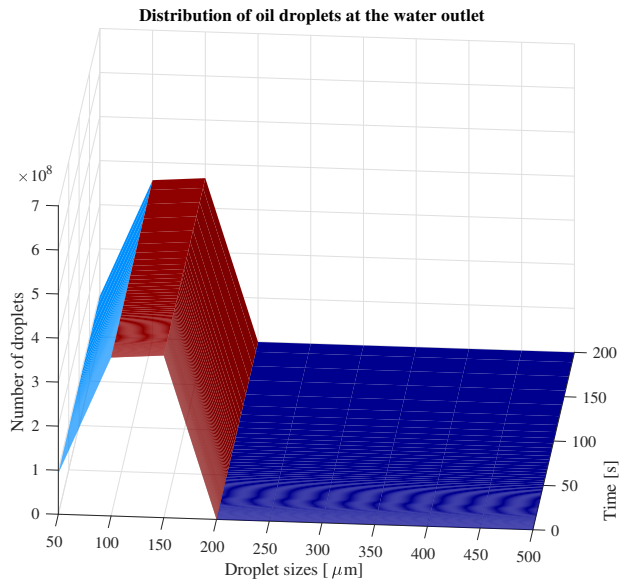


Figure 5.10: Distribution of oil droplets at the water outlet for water level = 2 m

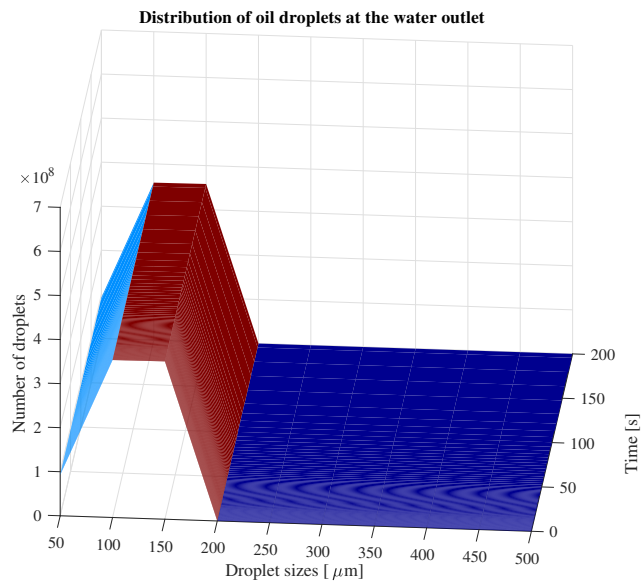


Figure 5.11: Distribution of oil droplets at the water outlet for water level = 2.2 m

Figures 5.8 - 5.11 show the amount of oil droplets present at the water outlet for the corresponding droplet size and time for different water levels, respectively. From Figure 5.8, it can be seen that for a water level of 0.9 m, the cut-off oil droplet size (minimum size above which the dispersed droplets present at the water outlet is zero) is observed to be 250 μm . This means that all oil droplets larger than 250 μm leave out of the water phase into their mother phase. With the increase in water level, the cut-off oil droplet size decreases (200 μm), meaning that more particle classes leave the water phase and enter their mother phase. Therefore, also the total volume of oil droplets leaving the water phase (V_{OoW}) will be higher. This is due to the fact that if the water level is smaller, the horizontal velocity of the droplets is higher. Hence, the droplets need less time to reach the outlet in comparison to the time it takes to reach their oil-water interphase (see Appendix C).

It can be observed from Figures 5.10 and 5.11 that for levels of water above 2 m the oil distribution at the water outlet remains almost nearly the same.

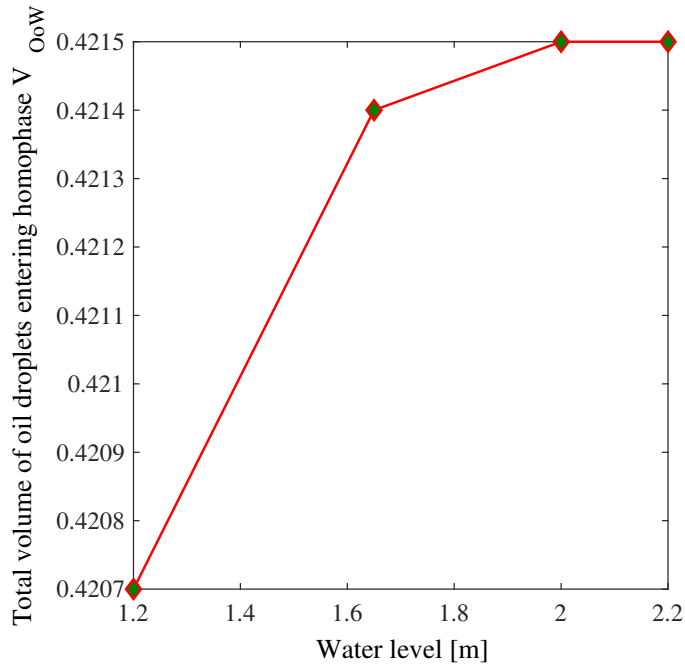


Figure 5.12: Total volume of oil droplets entering their continuous phase V_{OoW} for the respective water levels.

From Figure 5.12 it can be seen that with the increase in water level, the total volume of oil droplets leaving the water phase and entering the oil phase is numerically almost the same for water levels above 2 m. This explains the optimal level of the water obtained.

Conclusion

In this thesis, a Nonlinear Model Predictive Control (NMPC) strategy is proposed to control the levels and the system pressure of a three phase gravity separator. The NMPC optimization problem was formulated for the three phase system and solved using the Direct collocation approach (Simultaneous method) using CasADi within the MATLAB programming environment. Nonlinear dynamic models were derived for the water level, liquid level and the pressure of the gas in the system based on inflow and outflow dynamics. The equations were implemented in MATLAB as separate script files and were solved using ode15s solver. The models were formulated as a semi-explicit DAE system and the discontinuous algebraic equations were embedded into the optimizing controller as continuous algebraic equations using arctan functions.

The simulations were conducted for several cases in order to analyze the performance of the NMPC controller on the system behavior. Firstly, the system behavior was studied to analyze the performance of the controller to sensitivity in measurement noise. It was observed that the controller was sensitive to the large

magnitude in measurement noise. This could lead to the wear and tear of the control valves very easily. Therefore for the successive cases, the system performance was studied for a low magnitude noise added to the measurements. Secondly, the controller was studied for the tracking of the setpoint using NMPC scheme. It was seen that the states converged to their given setpoints while satisfying the constraints. However, at the time instants, where the step change in water level was introduced, constraint violation of the rate of change in manipulated variables was observed. Thirdly, the controller was studied for the case, where sudden disturbances in liquid and gas inflows to the gravity separator system were introduced. It was observed that the controller was able to reject the disturbances and regulate the output at the desired setpoint without any offset. Finally, a case was investigated to maximize the total volume of oil entering its native phase, and thereby to obtain the optimal water level. It was noticed that the optimal level of water was at 2 m and within the given bounds. This was because, at water level greater than 2 m, the efficiency of oil removal from the water phase was almost constant. Also, the distribution of oil droplets at the water outlet remained almost the same. Additionally, the numerical values from the simulation show that the total volume of oil entering its continuous phase at a water level greater than 2 m is almost constant.

From the simulations, it is observed that the controller is effectively able to track the setpoints, reject disturbances for a low magnitude of measurement noise since the controller is sensitive to high magnitude measurement noise thereby reacting positively to changes in the parameters. Results also agree with the simulations, and it can be concluded that the Nonlinear Model Predictive Controller is a good control method for oil-water-gas separators.

Future Work

The fluid dynamics inside the separators involve a lot of complexity, and hence it is necessary to study the phenomena like droplet breakage and coalescence. In actuality, there exists an initial distribution of droplet size at the inlet of the separator, and the size distribution changes inside the separator due to nonuniform flow patterns and factors such as coalescence and droplet breakage. This evolution of the droplet size distribution can be estimated by the use of Population Balance Equations (PBE) and incorporating these models can further improve the exactness of estimation.

For the absence of droplet breakage and coalescence, only Stokes law is sufficient to determine the terminal settling velocity of the droplets which has been considered for the purpose of modeling. However, while considering the coalescence and the break up of droplets which exists commonly in the separators, the velocity is hindered and the Stokes law for settling will need to take into account the correction factor. Introducing PBEs within the CasADi environment seems a bit challenging and also the simulations might take a lot of computation time.

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Appendices

Appendix A

Simulation Parameters

Table A.1: Simulation Parameters (Backi and Skogestad, 2017; Laleh et al., 2013).

Parameter	Description	Value
M_G	Molar mass of the gas	$0.01604 \text{ kg mol}^{-1}$
R	Universal gas constant	$8.314 \text{ kg m}^2 \text{ s}^{-2} \text{ mol}^{-1} \text{ K}^{-1}$
T	Temperature	328.5 K
g	Gravitational acceleration	9.8 m s^{-2}
$q_{G,in}$	Volumetric gas inflow	$0.456 \text{ m}^3 \text{ s}^{-1}$
$q_{L,in}$	Volumetric liquid inflow	$0.59 \text{ m}^3 \text{ s}^{-1}$
μ_O	Viscosity of oil	$0.001 \text{ kg (m s)}^{-1}$
μ_W	Viscosity of water	$0.0005 \text{ kg (m s)}^{-1}$
ρ_G	Density of gas	49.7 kg m^{-3}
ρ_O	Density of oil	831.5 kg m^{-3}
ρ_W	Density of water	1030 kg m^{-3}

Quadrature coefficients

The aim is to evaluate the integral $I = \int_a^b f(x) dx$. Assuming that the values of the function $f(x)$ is known at $n + 1$ points; $x_0, \dots, x_n \in [a, b]$. With the function values $f(x_0), \dots, f(x_n)$, the Lagrange interpolating polynomial is written in the form

$$L_n(x) = \sum_{i=0}^n f(x_i)P_i(x), \tag{B.1}$$

where

$$P_i(x) = \prod_{j=0, j \neq i}^n \frac{x - x_j}{x_i - x_j}, \quad 0 \leq i \leq n \tag{B.2}$$

The integral $I = \int_a^b f(x) dx$ is then approximated by the integral of $L_n(x)$:

$$\int_a^b f(x) dx \approx \int_a^b L_n(x) dx = \sum_{i=0}^n f(x_i) \int_a^b P_i(x) dx = \sum_{i=0}^n \omega_i f(x_i) \quad (\text{B.3})$$

This way the integral is converted to a summation

$$\int_a^b f(x) dx = \sum_{i=0}^n \omega_i f(x_i)$$

Therefore the quadrature coefficients ω_i in Equation B.3 is expressed by the equation given below

$$\omega_i = \int_a^b P_i(x) dx. \quad (\text{B.4})$$

Residence time of droplet classes

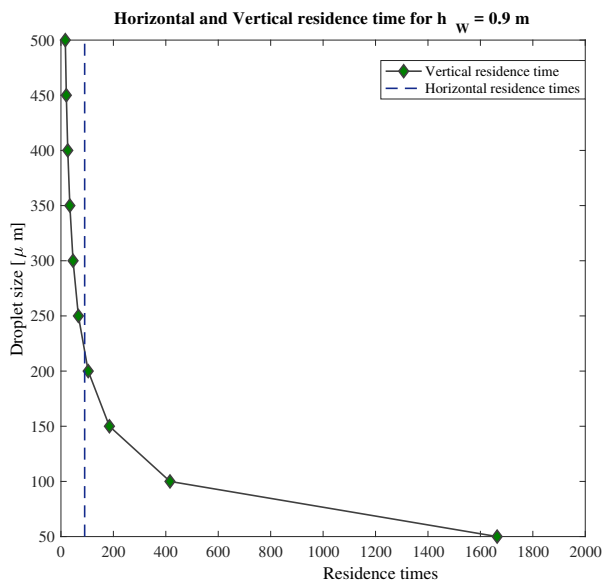


Figure C.1: Horizontal and vertical residence times of oil droplets in the water phase for water level = 0.9 m. The horizontal residence time = 90.45 s.

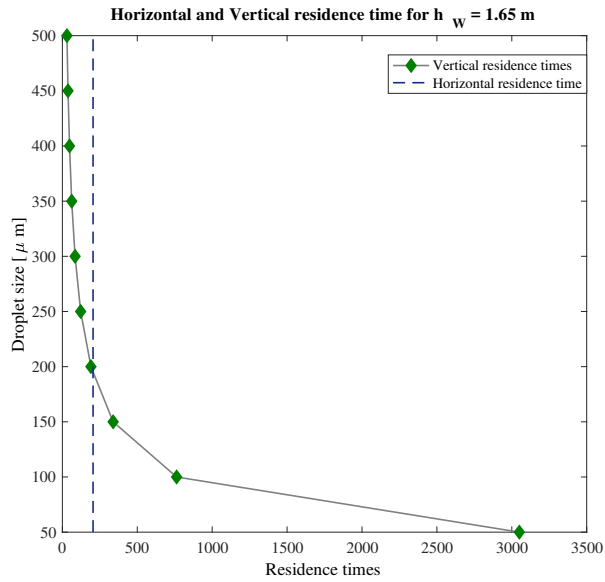


Figure C.2: Horizontal and vertical residence times of oil droplets in the water phase for water level = 1.65 m. The horizontal residence time = 204.715 s.

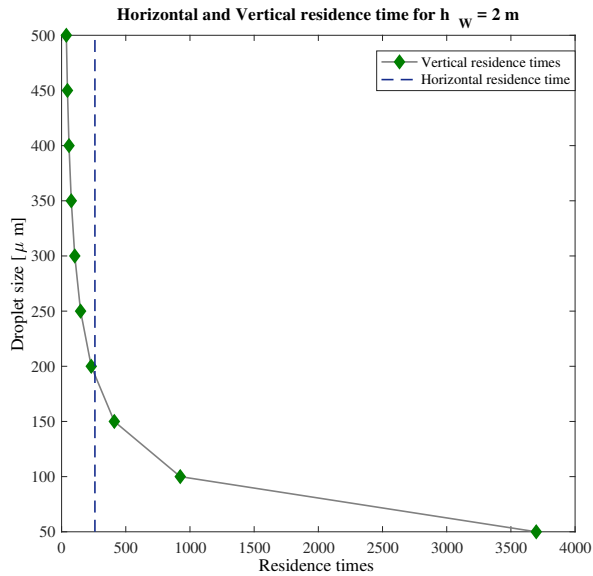


Figure C.3: Horizontal and vertical residence times of oil droplets in the water phase for water level = 2 m. The horizontal residence time = 259.58 s.

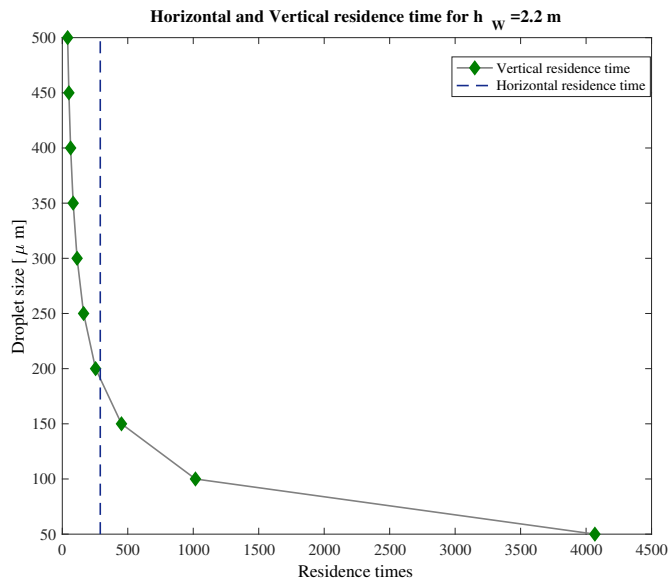


Figure C.4: Horizontal and vertical residence times of oil droplets in the water phase for water level = 2.2 m. The horizontal residence time = 289.96 s.

The above figures indicate the horizontal (blue dotted line) and vertical residence times (green diamond marked) of droplets in the water phase. It is understood that, if the horizontal residence time is greater than the vertical residence times, the droplets will leave to their continuous phase.

It can be observed from Figure C.1 that all droplets above 250 μm leave to their continuous phase for a water level of 0.9 m. With the increase in water level, the cut-off droplet size is 200 μm as can be seen from Figures C.2 - C.4 meaning all droplets above the cut-off droplet size enter their mother phase.

Appendix D

Rate of change constraints

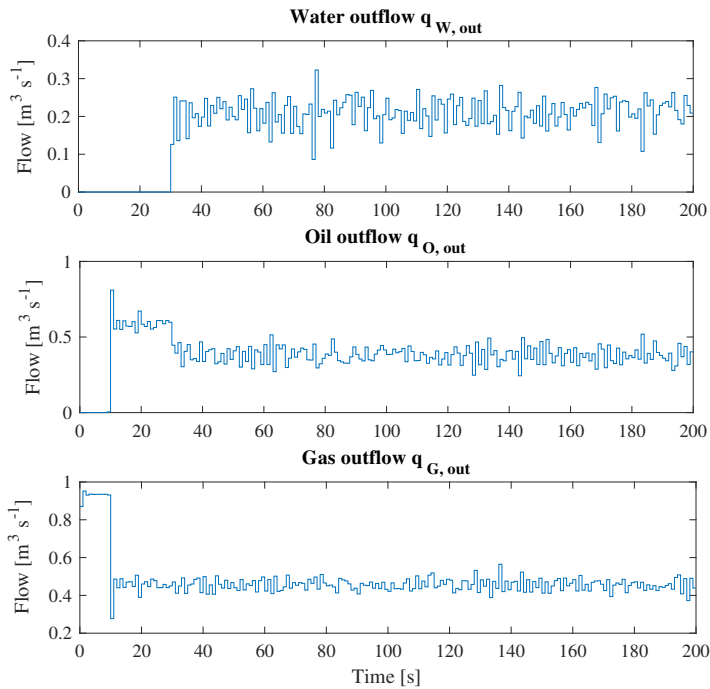


Figure D.1: The control inputs without the inclusion of rate of change constraints

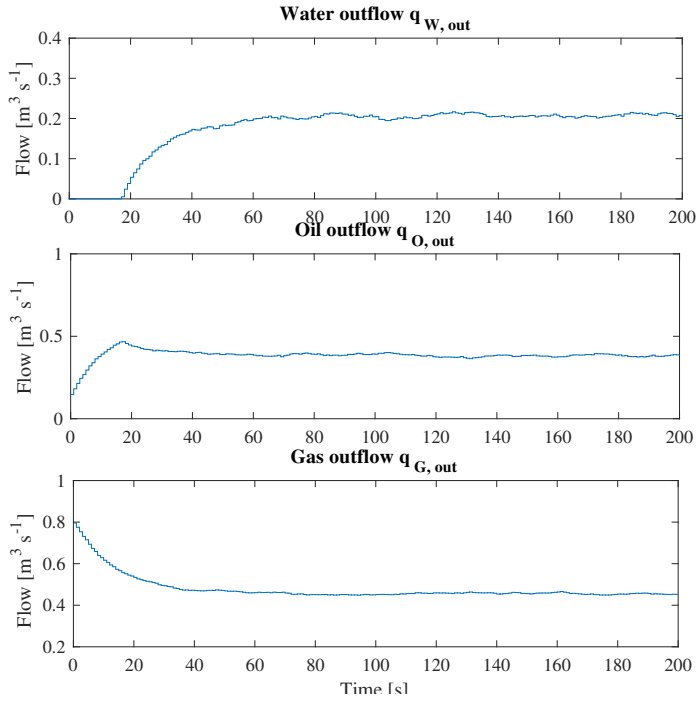


Figure D.2: The control inputs with the inclusion of rate of change constraints

Without the implementation of the rate of change constraint $\Delta u^T R \Delta u$ the plots for the control inputs is as shown in Figure D.1, and with the implementation of the above mentioned constraint, the plot as shown in Figure D.2 is obtained. The inclusion of this constraint in the optimization problem penalizes the control moves ($\Delta u_t = u_t - u_{t-1}$), thereby reduces the wear and tear of valves.

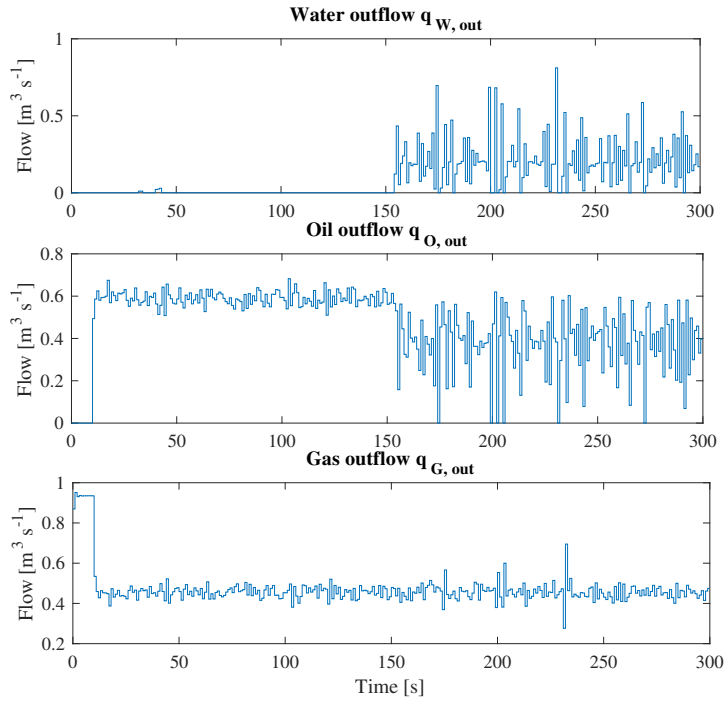


Figure D.3: The control inputs without the inclusion of rate of change constraints

Without the implementation of the rate of change constraint $\Delta u^T R \Delta u$ the plots for the control inputs is as shown in Figure D.3.

Appendix E

MATLAB CODE

E.1 velocity_inflows.m

```
1 function [vel_oil , vel_water , qWin] = velocity_inflows (qLin , hL , hW ,  
    alpha , phi_wo , phi_ow , radius )  
2  
3 abs_wo = alpha*phi_wo;           % Fraction of water going into  
    the continuous oil phase (subscript wo -> water to oil)  
4 abs_ww = alpha*(1-phi_wo);       % Fraction of water going into  
    the continuous water phase (subscript ww -> water to water)  
5 abs_ow = (1-alpha)*phi_ow;       % Fraction of oil going into  
    the continuous water phase (subscript ow -> oil to water)  
6 abs_oo = (1-alpha)*(1-phi_ow);   % Fraction of oil going into  
    the continuous oil phase (subscript oo -> oil to oil)  
7  
8 flow_fraction_w = abs_ww + abs_ow; % Split factor describing the  
    amount of oil and water going into the water phase  
9 flow_fraction_o = abs_oo + abs_wo; % Split factor describing the  
    amount of oil and water going into the oil phase
```

```

10
11 AL = radius^2/2*(2*acos((radius-hL)/radius)-sin(2*acos((radius-hL)
    /radius))); % Cross sectional area of liquid
12 AW = radius^2/2*(2*acos((radius-hW)/radius)-sin(2*acos((radius-hW)
    /radius))); % Cross sectional area of water
13 AO = AL-AW; % Cross
    sectional area of oil
14
15 vel_oil = flow_fraction_o*qLin/AO; % Velocity of the
    continuous oil phase
16 vel_water = flow_fraction_w*qLin/AW; % Velocity of the
    continuous water phase
17
18 qWin = qLin*flow_fraction_w; % Fraction of liquid inflow
    going into to water phase

```

E.2 gravity_ode.m

```

1 function dydt = gravity_ode(t,y, parameters_plant)
2
3 % ASSIGN EMPTY VECTOR
4 dydt = zeros(3,1);
5
6 % DEFINE THE PASSING PARAMETERS
7 qLin= parameters_plant(1);
8 alpha = parameters_plant(2);
9 phi_wo = parameters_plant(3) ;
10 phi_ow = parameters_plant(4);
11 radius = parameters_plant(5);
12 L = parameters_plant(6) ;
13 R = parameters_plant(7);
14 Temp = parameters_plant(8) ;

```

```

15 density_gas = parameters_plant(9);
16 MG = parameters_plant(10);
17 qGin = parameters_plant(11);
18 qWout = parameters_plant(12);
19 qOout = parameters_plant(13);
20 qGout = parameters_plant(14);
21 hW = parameters_plant(15);
22 hL = parameters_plant(16);
23 p = parameters_plant(17);
24
25 % CALLING THE FUNCTIONS
26 [vel_oil , vel_water , qWin] = velocity_inflows (qLin , hL , hW , alpha ,
    phi_wo , phi_ow , radius); % horizontal velocity , inflows
    calculation model
27 [~, VOOw_vector, ~] = oildroplets_in_water(vel_water , hW , L);
    % oildroplets_in_Water calculation
    model
28 [~, VWoO_vector, ~] = waterdroplets_in_oil(vel_oil , hL , hW , L);
    % waterdroplets_in_oil calculation model
29
30
31 % DEFINING THE VARIABLES
32 time_oil = L/vel_oil; % used in dhW/dt model
33 time_water = L/vel_water; % used in dhW/dt model
34 constant = R*Temp*density_gas/MG; % used in dp/dt model
35
36 % DIFFERENTIAL EQUATIONS
37
38 %dhW/dt
39 dydt(1) = (qWin- qWout + sum(VWoO_vector)/time_oil- sum(
    VOOw_vector)/time_water)/(2*L*sqrt(2*radius*y(1) - y(1)^2));
40
41 %dhL/dt

```

```

42 dydt(2) = (qLin - qOout-qWout)/(2*L*sqrt(2*radius*y(2) - y(2)^2));
43
44 %dp/dt
45 dydt(3) = 1e-5* 1/(pi*radius^2*L-radius^2/2*(2*acos((radius-y(2))/
    radius)-sin(2*acos((radius-y(2))/radius))*L) * (constant*(
    qGin-qGout) + 1e5*y(3)*(qLin-qOout-qWout));

```

E.3 oildroplets_in_water.m

```

1 function [noil_new, Voil_new, lpp_oil_new] = oildroplets_in_water(
    v_horizontal, h_water, n, L)
2
3 % v_horizontal is the horizontal velocity of oil droplets =
    horizontal
4 % velocity of the water phase
5
6 % h_water is the height of the water.
7 % n is the vector of initial oil particles size distribution.
8 % L is the length of the Gravity Separator
9 % d is the vector of diameter of all the oil particles.
10 % s is the vertical velocity of the droplets.
11 % V is the volume of oil droplets
12
13 d = zeros(1,10)';
14 V = zeros(1,10)';
15 s = zeros(1,10)';
16 lpp_oil_new = zeros(1,10)';
17 noil_new = zeros(1,10)';
18 Voil_new = zeros(1,10)';
19 density_oil = 831.5;
20 density_water = 1030;
21 dyn_visc_water = 0.5e-3;

```

```

22
23 for i = 1:10
24 d(i) = 50*10^(-6)*i;
25 V(i) = 4/3*pi*(d(i)/2)^3;
26 s(i) = abs(((d(i))^2)*9.81*(density_oil-density_water)/(18*
      dyn_visc_water));
27
28 if (L/v_horizontal)>= ( h_water/s(i))
29
30 noil_new(i) = 0;           % number of oil droplets present in the
      water phase
31 Voil_new(i) = n(i)* V(i); % volume of oil droplets leaving the
      water phase
32 lpp_oil_new(i)= h_water; % position of the oil particles
33 else
34 lpp_oil_new(i)= (L/v_horizontal)*s(i);
35 noil_new(i) = (1-(lpp_oil_new(i)/h_water))* n(i);
36 Voil_new(i) = n(i) * V(i)*(lpp_oil_new(i)/h_water);
37 end
38 end

```

E.4 waterdroplets_in_oil.m

```

1 function [nwater_new , Vwater_new , lpp_water_new] =
      waterdroplets_in_oil(v_horizontal , h_liquid , h_water , n , L)
2
3 % v_horizontal is the horizontal velocity of water droplets =
      horizontal
4 % velocity of the bulk oil phase
5
6 % h_water is the height of the water.
7 % n is the vector of initial water particles size distribution.

```

```

8 % L is the length of the Gravity Separator
9 % d is the vector of diameter of all the water particles .
10 % s is the vertical velocity of the droplets .
11 % V is the volume of water droplets
12
13 h_oil = h_liquid - h_water;
14 d = zeros(1,10)';
15 V = zeros(1,10)';
16 s = zeros(1,10)';
17 lpp_water_new = zeros(1,10)';
18 nwater_new = zeros(1,10)';
19 Vwater_new = zeros(1,10)';
20 density_oil = 831.5;
21 density_water = 1030;
22 dyn_visc_oil = 1e-3;
23
24 for i = 1:10
25 d(i) = 50*10^(-6)*i;
26 V(i) = 4/3*pi*(d(i)/2)^3;
27 s(i) = abs(((d(i))^2)*9.81*(density_oil-density_water)/(18*
    dyn_visc_oil));
28
29 if (L/v_horizontal)>= (h_oil/s(i))
30     nwater_new(i) = 0;           % number of water droplets
    present
31                                     % in the oil phase
32     Vwater_new(i) = n(i)* V(i); % volume of water droplets
    leaving
33                                     % the oil phase
34     lpp_water_new(i)= h_water; % position of the water
    particles
35 else
36     lpp_water_new(i)= (L/v_horizontal)*s(i);

```

```

37     nwater_new(i) = (1-(lpp_water_new(i)/h_oil))* n(i);
38     Vwater_new(i) = n(i) * V(i)*(lpp_water_new(i)/h_oil);
39     end
40 end

```

E.5 NMPC.m

```

1  clear all
2
3  addpath('F:\NTNU')
4  import casadi.*
5  % Running using only oil droplets in water algebraic equations ,
   here
6  % optimal solution is found
7
8  % TIME HORIZON
9  T = 20;
10
11 % DECLARE VARIABLES
12 qLin= 0.59 ;           % Inlet liquid flow rate
13 alpha = 0.135;       % Water cut
14 phi_wo = 0.3;        % Fraction of incoming water entering
   the oil phase
15 phi_ow = 0.3;        % Fraction of incoming oil entering the
   water phase
16 radius = 1.65;       % Radius of the gravity separator
17 L = 10 ;             % Length of gravity separator
18 R = 8.314;           % Universal gas constant
19 Temp = 328.5 ;      % Temperature
20 density_gas = 49.7;  % Density of gas
21 MG = 0.01604;       % Molecular weight of gas
22 qGin = 0.456;       % Inlet gas flowrate

```

```

23
24 phi_ww = 1-phi_wo ;           % fraction of water entering the water
    phase
25 phi_oo = 1-phi_ow ;           % fraction of oil entering the oil phase
26 V_sep = 3.14*radius*radius*L; % Total separator volume
27
28 density_oil = 831.5;
29 density_water = 1030;
30 dyn_visc_water = 0.5e-3;
31 dyn_visc_oil =1e-3;
32
33 % INITIAL VALUES FOR THE INTEGRATORS OF THE DYNAMIC STATES
34 hWinit = 1;                    % Initial value for the water level [m]
35 hLinit = 2.3;                  % Initial value for the liquid level [m]
36 pinit = 68.5;                 % Initial pressure [bar]
37
38 % PARTICLE DIAMETERS
39 d1 = 50*10^(-6);
40 d2 = 100*10^(-6);
41 d3 = 150*10^(-6);
42 d4 = 200*10^(-6);
43 d5 = 250*10^(-6);
44 d6 = 300*10^(-6);
45 d7 = 350*10^(-6);
46 d8 = 400*10^(-6);
47 d9 = 450*10^(-6);
48 d10= 500*10^(-6);
49
50 % VOLUME OF EACH PARTICLE CLASS
51 V1 = 4/3*pi*(d1/2)^3;
52 V2 = 4/3*pi*(d2/2)^3;
53 V3 = 4/3*pi*(d3/2)^3;
54 V4 = 4/3*pi*(d4/2)^3;

```

```

55 V5 = 4/3*pi*(d5/2)^3;
56 V6 = 4/3*pi*(d6/2)^3;
57 V7 = 4/3*pi*(d7/2)^3;
58 V8 = 4/3*pi*(d8/2)^3;
59 V9 = 4/3*pi*(d9/2)^3;
60 V10 = 4/3*pi*(d10/2)^3;
61
62 V = [V1 V2 V3 V4 V5 V6 V7 V8 V9 V10];
63
64 % TOTAL PARTICLE DISTRIBUTION
65 n = [1e8 5e8 1e9 5e9 1e10 1e10 5e9 1e9 5e8 1e8];
66
67 % TOTAL INITIAL VOLUE DISTRIBUTION OF ALL PARTICLES
68 init_V = sum(n .* V);
69
70 % SETTLING VELOCITY OF OIL PARTICLE CLASS IN THE WATER PHASE
71 so_1 = abs(((d1)^2)*9.81*(density_oil-density_water)/(18*
        dyn_visc_water));
72 so_2 = abs(((d2)^2)*9.81*(density_oil-density_water)/(18*
        dyn_visc_water));
73 so_3 = abs(((d3)^2)*9.81*(density_oil-density_water)/(18*
        dyn_visc_water));
74 so_4 = abs(((d4)^2)*9.81*(density_oil-density_water)/(18*
        dyn_visc_water));
75 so_5 = abs(((d5)^2)*9.81*(density_oil-density_water)/(18*
        dyn_visc_water));
76 so_6 = abs(((d6)^2)*9.81*(density_oil-density_water)/(18*
        dyn_visc_water));
77 so_7 = abs(((d7)^2)*9.81*(density_oil-density_water)/(18*
        dyn_visc_water));
78 so_8 = abs(((d8)^2)*9.81*(density_oil-density_water)/(18*
        dyn_visc_water));

```

```

79 so_9 = abs(((d9)^2)*9.81*(density_oil-density_water)/(18*
      dyn_visc_water));
80 so_10 = abs(((d10)^2)*9.81*(density_oil-density_water)/(18*
      dyn_visc_water));
81
82 % DECLARE MODEL VARIABLES
83 % State variables
84 x1 = MX.sym('x1');           % Water level
85 x2 = MX.sym('x2');           % Liquid level
86 x3 = MX.sym('x3');           % System Pressure
87
88 % Control variables
89 u1 = MX.sym('u1');           % Outlet Water flowrate
90 u2 = MX.sym('u2');           % Outlet Oil flowrate
91 u3 = MX.sym('u3');           % Outlet gas flowrate
92
93 %Algebraic variables (particle position of oil in the water phase)
94 lpp1 = MX.sym('lpp1');
95 lpp2 = MX.sym('lpp2');
96 lpp3 = MX.sym('lpp3');
97 lpp4 = MX.sym('lpp4');
98 lpp5 = MX.sym('lpp5');
99 lpp6 = MX.sym('lpp6');
100 lpp7 = MX.sym('lpp7');
101 lpp8 = MX.sym('lpp8');
102 lpp9 = MX.sym('lpp9');
103 lpp10 = MX.sym('lpp10');
104
105 % CONCATENATE VARIABLES
106 x = [x1;x2;x3];             % state variables
107 u = [u1; u2; u3];           % control variables
108 lpp = [lpp1; lpp2; lpp3; lpp4; lpp5; lpp6; lpp7; lpp8; lpp9; lpp10
      ]; % particle position alg. variables

```

```

109
110 % DECLARE VARIABLES
111 flow_fraction_w = (alpha*(1-phi_wo)) + ((1-alpha)*phi_ow);
112 AW = (radius^2*acos((radius-x1)/radius)) - ((radius - x1)*sqrt(2*
      radius*x1 - x1*x1));
113 vel_water = (flow_fraction_w*qLin)/AW;
114 time_water = L/ vel_water ;
115
116 A.L = (radius^2*acos((radius-x2)/radius)) - ((radius - x2)*sqrt(2*
      radius*x2 - x2*x2));
117 AO = A.L-AW;
118 flow_fraction_o = ((1-alpha)*(1-phi_ow)) + (alpha*phi_wo);
119 vel_oil = (flow_fraction_o*qLin)/AO;
120 time_oil = L/vel_oil;
121
122 % DEFINE ARGUMENT FUNCTION(residence time difference of oil
      droplets)
123 arg1 = (L/vel_water)-(x1/so_1);
124 arg2 = (L/vel_water)-(x1/so_2);
125 arg3 = (L/vel_water)-(x1/so_3);
126 arg4 = (L/vel_water)-(x1/so_4);
127 arg5 = (L/vel_water)-(x1/so_5);
128 arg6 = (L/vel_water)-(x1/so_6);
129 arg7 = (L/vel_water)-(x1/so_7);
130 arg8 = (L/vel_water)-(x1/so_8);
131 arg9 = (L/vel_water)-(x1/so_9);
132 arg10 = (L/vel_water)-(x1/so_10);
133
134 % USE ARCTAN FUNCTION TO EXPRESS CONTINUOUS SYSTEM
135 arctan_func1 = (atan((arg1).*1000*pi)+pi/2) ./ pi;
136 arctan_func2 = (atan((arg2).*1000*pi)+pi/2) ./ pi;
137 arctan_func3 = (atan((arg3).*1000*pi)+pi/2) ./ pi;
138 arctan_func4 = (atan((arg4).*1000*pi)+pi/2) ./ pi;

```

```

139 arctan_func5 = (atan((arg5).*1000*pi)+pi/2) ./ pi;
140 arctan_func6 = (atan((arg6).*1000*pi)+pi/2) ./ pi;
141 arctan_func7 = (atan((arg7).*1000*pi)+pi/2) ./ pi;
142 arctan_func8 = (atan((arg8).*1000*pi)+pi/2) ./ pi;
143 arctan_func9 = (atan((arg9).*1000*pi)+pi/2) ./ pi;
144 arctan_func10 = (atan((arg10).*1000*pi)+pi/2) ./ pi;
145
146 % DEFINING ALGEBRAIC EQUATIONS
147 f_z1 = lpp1 - ((arctan_func1)*x1 + (1-arctan_func1)*(L*so_1 ./
    vel_water));
148 f_z2 = lpp2 - ((arctan_func2)*x1 + (1-arctan_func2)*(L*so_2 ./
    vel_water));
149 f_z3 = lpp3 - ((arctan_func3)*x1 + (1-arctan_func3)*(L*so_3 ./
    vel_water));
150 f_z4 = lpp4 - ((arctan_func4)*x1 + (1-arctan_func4)*(L*so_4 ./
    vel_water));
151 f_z5 = lpp5 - ((arctan_func5)*x1 + (1-arctan_func5)*(L*so_5 ./
    vel_water));
152 f_z6 = lpp6 - ((arctan_func6)*x1 + (1-arctan_func6)*(L*so_6 ./
    vel_water));
153 f_z7 = lpp7 - ((arctan_func7)*x1 + (1-arctan_func7)*(L*so_7 ./
    vel_water));
154 f_z8 = lpp8 - ((arctan_func8)*x1 + (1-arctan_func8)*(L*so_8 ./
    vel_water));
155 f_z9 = lpp9 - ((arctan_func9)*x1 + (1-arctan_func9)*(L*so_9 ./
    vel_water));
156 f_z10 = lpp10 - ((arctan_func10)*x1 + (1-arctan_func10)*(L*so_10 ./
    vel_water));
157
158 % VOLUME OF OIL DROPLETS ENTERING THEIR NATIVE PHASE
159 VOoW1 = (lpp1 ./ x1)*n(1)* V1;
160 VOoW2 = (lpp2 ./ x1)*n(2)* V2;
161 VOoW3 = (lpp3 ./ x1)*n(3)* V3;

```

```

162 VOoW4 = (lpp4 ./ x1)*n(4)* V4;
163 VOoW5 = (lpp5 ./ x1)*n(5)* V5;
164 VOoW6 = (lpp6 ./ x1)*n(6)* V6;
165 VOoW7 = (lpp7 ./ x1)*n(7)* V7;
166 VOoW8 = (lpp8 ./ x1)*n(8)* V8;
167 VOoW9 = (lpp9 ./ x1)*n(9)* V9;
168 VOoW10 =(lpp10 ./ x1)*n(10)* V10;
169
170 VOoW = (VOoW1+ VOoW2 + VOoW3 + VOoW4 + VOoW5 + VOoW6 + VOoW7 +
        VOoW8 + VOoW9 + VOoW10);
171
172 % MPC DIFFERENTIAL EQUATIONS
173
174 xdot1 = (qLin*(alpha*phi_ww+(1-alpha)*phi_ow)-u1 -(VOoW/time_water
        ))/(2*L*sqrt(2*radius*x1 - x1^2));
175 xdot2 = (qLin - u2 -u1)/(2*L*sqrt(2*radius*x2 - x2^2));
176     VG = V_sep - A.L*L;
177 xdot3 = (1/VG)*((((R*Temp*density_gas)/(MG))*(qGin - u3))+(x3
        *10^5)*(qLin-u1-u2))*10^-5;
178
179 % CONCATENATE EQUATIONS TO FORM SEMI EXPLICIT DAE SYSTEM
180 diff = [xdot1; xdot2; xdot3]; % Differential eqns
181 alg = [f_z1; f_z2; f_z3; f_z4; f_z5; f_z6; f_z7; f_z8; f_z9; f_z10
        ];% Particle position alg expression
182
183 % MPC OBJECTIVE TERM
184
185 L1 = (x1 - 1.2)^2 + (x2 - 2.5 )^2 + (x3 - 68.7)^2; % CASE 1
186 %L1 = 100* (x2 - 2.5)^2 + 100*(x3 - 68.7)^2 - 1000*(VOoW)^2; %
        CASE 3
187
188 % CONTINUOUS TIME DYNAMICS

```

```

189 f = Function('f',{x,lpp,u},{diff,alg,L1},{ 'x','z','p'},{ 'xdot','zj
      ','qj'});
190
191 %% -----DIRECT COLLOCATION SCHEME-----%%
192
193 % DEGREE OF INTERPOLATING POLYNOMIAL
194 d = 3;
195
196 % GET COLLOCATION POINTS
197 tau_root = [0 collocation_points(d, 'radau')];
198
199 % COEFFICIENTS OF THE COLLOCATION EQUATION
200 C = zeros(d+1,d+1);
201
202 % COEFFICIENTS OF THE CONTINUITY EQUATION
203 D = zeros(d+1, 1);
204
205 % COEFFICIENTS OF THE QUADRATURE FUNCTION
206 B = zeros(d+1, 1);
207
208 % CONSTRUCT POLYNOMIAL BASIS
209 for j=1:d+1
210     % CONSTRUCT LAGRANGE POLYNOMIALS TO GET THE POLYNOMIAL BASIS AT
           THE COLLOCATION POINT
211     coeff = 1;
212     for r=1:d+1
213         if r ~= j
214             coeff = conv(coeff, [1, -tau_root(r)]);
215             coeff = coeff / (tau_root(j)-tau_root(r));
216         end
217     end
218

```

```

219 % EVALUATE POLYNOMIAL AT THE FINAL TIME TO GET THE COEFFICIENTS
      OF CONTINUITY EQUATION
220 D(j) = polyval(coeff, 1.0);
221
222 % EVALUATE TIME DERIVATIVE OF THE POLYNOMIAL AT ALL COLLOCATION
      POINTS TO GET COEFFICIENTS OF CONTINUITY EQUATION
223 pder = polyder(coeff);
224 for r=1:d+1
225     C(j,r) = polyval(pder, tau_root(r));
226 end
227
228 % EVALUATE INTEGRAL OF THE POLYNOMIAL TO GET COEFFICIENTS OF THE
      QUADRATURE FUNCTION
229 pint = polyint(coeff);
230 B(j) = polyval(pint, 1.0);
231 end
232
233 % CONTROL DISCRETIZATION
234 N = 10; % Number of control intervals
235 h = T/N; % Size of each control interval
236
237 % START WITH AN EMPTY NLP
238 w = {};
239 w0 = [];
240 lbw = [];
241 ubw = [];
242 J = 0;
243 g={};
244 lbg = [];
245 ubg = [];
246
247 % "LIFT" INITIAL CONDITIONS
248 X0 = MX.sym('X0', 3);

```

```

249 Z0 = MX.sym('ZO', 10);
250 w = {w{:}}, X0, Z0};
251 lbw = [lbw; 1; 2.3; 68.5; zeros(10,1)]; % Assign initial
      conditions to lower bound
252 ubw = [ubw; 1; 2.3; 68.5; zeros(10,1)]; % Assign initial
      conditions to upper bound
253 w0 = [w0; 1; 2.3; 68.5; zeros(10,1)]; % Initial conditions
      value
254
255
256 %% FORMULATE THE NLP (DIRECT COLLOCATION)
257 Xk = X0;
258 for k = 0: N-1
259     % NEW NLP VARIABLE FOR EACH CONTROL
260     Uk = MX.sym(['U_' num2str(k)], 3);
261     w = {w{:}}, Uk};
262     lbw = [lbw; 0;0;0];
263     ubw = [ubw; 2 ; 2; 5];
264     w0 = [w0; 0; 0; 0];
265
266     % STATE AT COLLOCATION POINTS
267     Xkj = {}; Zkj = {};
268     for j=1:d
269         Xkj{j} = MX.sym(['X_' num2str(k) '_' num2str(j)], 3);
270         Zkj{j} = MX.sym(['Z_' num2str(k) '_' num2str(j)], 10);
271         w = {w{:}}, Xkj{j}, Zkj{j}};
272         lbw = [lbw; 0.9; 1.9; 50; zeros(10,1)];
273         ubw = [ubw; 2.2; 2*radius; 100; 2*ones(10,1)];
274         w0 = [w0; 1; 2.3; 68.5; zeros(10,1)];
275     end
276
277     % LOOP OVER COLLOCATION POINTS
278     Xk_end = D(1)*Xk; %where D is the continuity eqn coeff

```

```

279     for j=1:d
280         % EXPRESSION FOR STATE DERIVATIVE AT COLLOCATION POINT(
                polynomial approx)
281         xp = C(1,j+1)*Xk; % C is the derivative at the collocation
                points
282         for r=1:d
283             xp = xp + C(r+1,j+1)*Xkj{r};
284         end
285         % APPEND COLLOCATION EQUATIONS
286         [fj ,zj ,qj] = f(Xkj{j},Zkj{j},Uk);          % fj = dxdt & qj =
                objective function value , zj = solution of alg
                expression
287         g = {g{:}, h*fj - xp , zj};                % slope constraints
288         lbg = [lbg; zeros(3,1); zeros(10,1)];
289         ubg = [ubg; zeros(3,1); zeros(10,1)];
290
291         % ADD CONTRIBUTION TO THE END STATE
292         Xk_end = Xk_end + D(j+1)*Xkj{j};
293
294         % ADD CONTRIBUTION TO THE QUADRATURE FUNCTION
295         J = J + B(j+1)*qj*h;
296
297     end
298     % NEW NLP VARIABLE FOR STATE AT THE END OF THE EQUATION
299     Xk = MX.sym(['X_' num2str(k+1)], 3); % this value is updated
300     w = {w{:}, Xk};
301     lbw = [lbw; 0.9 ; 1.9; 50];
302     ubw = [ubw; 2.2; 2*radius; 100];
303     w0 = [w0 ; 1; 2.3; 68.5];
304
305     % ADD EQUALITY CONSTRAINT
306     g = {g{:}, Xk_end-Xk};                % shooting gap constraint
307     lbg = [lbg; 0; 0 ;0];

```

```

308     ubg = [ubg; 0; 0 ;0];
309
310     % ADD CONTROL MOVES AND PENALTY TERM IN THE OBJECTIVE FUNCTION
311     if k>0
312         % ADD EQUALITY CONSTRAINT
313         g = {g{:}, Uk - Uk_prev};
314         lbg = [lbg; -0.05; -0.05 ; -0.05];
315         ubg = [ubg; 0.05; 0.05 ; 0.05];
316         % INCLUDE THE PENALTY TERM IN OBJECTIVE FUNCTION WITH WEIGHT
           FACTOR
317         J = J + ((Uk(1) - Uk_prev(1)).^2) + ((Uk(2) - Uk_prev(2)).^2)
           + ((Uk(3) - Uk_prev(3)).^2);
318     end
319
320     Uk_prev = Uk;
321 end
322 %%
323
324 % CREATE AN NLP SOLVER
325 prob = struct('f', J, 'x', vertcat(w{:}), 'g', vertcat(g{:}));
326 options.ipopt.print_level = 0; % options does not allow to print
           at every iteration
327 options.print_time = false;
328 solver = nlpsol('solver', 'ipopt', prob, options);
329
330 % TOTAL SIMULATION TIME
331 tsim =300;
332
333 % DEFINE EMPTY MATRICES
334 qWout_vec = zeros(1,tsim);
335 qOout_vec = zeros(1,tsim);
336 qGout_vec = zeros(1,tsim);
337

```

```

338 position_oilparticles_vec = zeros(tsim,10);
339 volume_oilparticles_vec = zeros(tsim,10);
340
341 % PASSING INITIAL VALUES TO PLANT MODEL
342 hW = hWinit;
343 hL = hLinit;
344 p = pinit;
345
346 for sim = 1: tsim
347     % SOLVE THE NLP
348     sol = solver('x0', w0, 'lbx', lbw, 'ubx', ubw, 'lbg', lbg, 'ubg
        ', ubg);
349     w_opt = full(sol.x);
350     sim           % helps to know the current
        iteration
351
352     % SOLUTION
353     x1_opt = w_opt(56:6+13*d:end);
354     x2_opt = w_opt(57:6+13*d:end);
355     x3_opt = w_opt(58:6+13*d:end);
356
357     u1_opt = w_opt(14:6+13*d:end);
358     u2_opt = w_opt(15:6+13*d:end);
359     u3_opt = w_opt(16:6+13*d:end);
360
361     % FIRST VALUE OF THE OPTIMIZER SENT TO PLANT MODEL
362     qWout= u1_opt(1);
363     qOout= u2_opt(1);
364     qGout= u3_opt(1);
365
366     % SET EMPTY MATRICES TO STORE THE OPTIMUM VALUES
367     qWout_vec(sim) = qWout;
368     qOout_vec(sim) = qOout;

```

```

369     qGout_vec(sim) = qGout;
370
371     % CONCATENATE THE OPTIMUM INPUTS
372     control_inputs = [qWout;qOout;qGout];
373     control_inputs_all(:,sim) = [control_inputs];
374
375     % SIMULATE PLANT MODEL
376     parameters_plant = [qLin, alpha, phi_wo, phi_ow, radius, L, R,
        Temp, density_gas, MG, qGin, qWout, qOout, qGout, hW, hL,
        p];
377     [t,y]=ode15s(@gravity_ode,[sim-1 sim],[hW, hL, p],[],
        parameters_plant);
378
379     % EXTRACT EACH FINAL ELEMENT OF THE STATES
380     hW = y(end,1);
381     hL = y(end,2);
382     p = y(end,3);
383
384     % ADDING NOISE TO THE STATES
385     hW = hW + 1e-3*randn(1);
386     hL = hL + 1e-3*randn(1);
387     p = p + 1e-2*randn(1);
388
389     [vel_oil,vel_water,qWin] = velocity_inflows(qLin,hL,hW,alpha,
        phi_wo,phi_ow,radius);
390     [~,VOoW_vector,lpp_oil_new] = oildroplets_in_water(
        vel_water, hW, L);
391
392     position_oilparticles_vec(sim,:) = [lpp_oil_new]';
393     volume_oilparticles_vec(sim,:) = [VOoW_vector]';
394
395     % STATES PERTURBED BY NOISE
396     states = [hW; hL; p];

```

```

397     states_all(:,sim) = [states];
398
399 %% SET NEW INITIAL VALUES FOR THE NEXT ITERATION
400 lbw = [];
401 ubw = [];
402 w0 = [];
403
404 % "LIFT" INITIAL CONDITIONS
405 lbw = [lbw; states; lpp_oil_new'];
406 ubw = [ubw; states; lpp_oil_new'];
407 w0 = [w0; states; lpp_oil_new'];
408
409     for k= 1: N
410         lbw = [lbw; 0; 0; 0];
411         ubw = [ubw; 2 ; 2; 5];
412         w0 = [w0; control_inputs];
413
414         % STATE AT COLLOCATION POINTS
415         for j=1:d
416             lbw = [lbw; 0.9; 1.9; 50; zeros(10,1)];
417             ubw = [ubw; 2.2; 2*radius; 100; 2*ones(10,1)];
418             w0 = [w0; states; lpp_oil_new'];
419         end
420
421         % NEW NLP VARIABLE FOR STATES AT THE END OF THE INTERVAL
422         lbw = [lbw; 0.9 ; 1.9; 50];
423         ubw = [ubw; 2.2 ; 2*radius; 100];
424         w0 = [w0 ; states];
425
426     end
427
428 end
429

```

```
430 tgrid = 0:1:tsim;
431 level_W1 = [1, states_all(1,:)];
432 level_L1 = [2.3, states_all(2,:)];
433 P1 = [68.5, states_all(3,:)];
434
435 qWout1 = [control_inputs_all(1,:), nan];
436 qOout1 = [control_inputs_all(2,:), nan];
437 qGout1 = [control_inputs_all(3,:), nan];
438
439 % PLOTTING
440 figure;
441 subplot(321)
442 plot(tgrid, level_W1)
443 subplot(322)
444 stairs(tgrid, qWout1)
445 subplot(323)
446 plot(tgrid, level_L1)
447 subplot(324)
448 stairs(tgrid, qOout1)
449 subplot(325)
450 plot(tgrid, P1)
451 subplot(326)
452 stairs(tgrid, qGout1)
```