

# Non-linear Analysis of Low-Frequency Combustion Instabilities in Liquid Rocket Engines

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## Abstract

In this paper we propose an analysis of low frequency combustion instabilities, taking advantage of the software EcosimPro. We implement a specific module based on the double time lag model and investigate the coupling of combustion chamber and feed line oscillations by using a complete set of non-linear equations. We identify the characteristic time lags following two approaches: (i) a constant time lag approach and (ii) a variable time lag approach based on correlations available in open literature. To prove the module capabilities we reproduce an experimental set up and generate a stability map, comparing our results with literature data from both experiments and a linear double time lag model. The stability boundaries obtained with the chugging module are in good agreement with those obtained in open literature, and the first characteristic frequency of the engine is well predicted. Furthermore, the model proves his capability in reconstructing the reversal in the slope of the stability boundary at low fuel injector pressure drops and in detecting the high frequency content typically observed when in presence of multi-mode oscillations. However, in our calculations the higher frequency does not dominate the instabilities, that is in the unstable regime the model diverges with a frequency equal to the first characteristic frequency. In the last part of the paper we use the variable time lag approach to investigate a portion of the aforementioned stability map. Thanks to the semi-empirical correlations we are able to improve the prediction of the first characteristic frequency, whereas the stability boundary does not change significantly and remains comparable with the one predicted by the constant double time lag approach.

## Nomenclature

$A$	cross section area	$\rho$	density
$Dp$	injector pressure drop	$\xi$	pressure drop coefficient
$E$	total energy	<i>Subscripts</i>	
$H$	total enthalpy	<i>atom</i>	atomisation
$p$	pressure	<i>c</i>	chamber
$R$	gas constant	<i>crit</i>	critical conditions
$T$	temperature	<i>fu</i>	fuel
$t$	time	$g$	gas
$u$	velocity	$l$	liquid
$V$	volume	$ox$	oxidizer
$x$	axial coordinate	$th$	throat
<i>Greek Symbols</i>		$vap$	vaporisation

## 1. Introduction

In this paper we propose an analysis of low frequency combustion instabilities, frequently referred to as chugging, where the typical components of a Liquid Rocket Engine (LRE) are modelled with an unsteady non-linear approach. Low frequency combustion instabilities in liquid rocket engines are characterised by oscillations in the mass flow rate of propellants that result in pulsations of the chamber pressure. The response of the combustion chamber to the upstream oscillations is not immediate. The delay is due to the different characteristic times associated with the processes of injection, atomisation, vaporisation, mixing and combustion, taking place in quick succession inside the combustion chamber. Moreover, the mass flow rate is not immediately responsive to pressure oscillations, because of the propellant inertia in the feed lines. Hence, both feed lines and combustion chamber physics must be taken into account. Low frequency instabilities have been usually investigated with the help of linearised models [1, 2, 3, 4, 5], whereas only a few attempts have been made to detail the phenomena with non-linear models [6, 7]. In the present study we propose an unsteady non-linear analysis of chugging instabilities, with particular attention to the appropriate modelling of delays induced by atomisation, vaporisation and mixing processes.

The analysis of chug instabilities requires to model the subsystems of a LRE to investigate the mutual influence of each component. Such a comprehensive simulation inherently needs simplified sub-models able to give a satisfactory level of reliability along with a reasonable computational time. In this context an object oriented simulation tool is a valid option to handle the complexity of a propulsion system. EcosimPro [8] is a simulation platform focused on the analysis of complete systems that spans different engineering fields. Thanks to the object oriented philosophy, single components can be connected to each other in order to analyse systems at different complexity levels. Propulsion systems can be investigated by the European Space Propulsion System Simulation (ESPSS) library [9], in which the typical components of a liquid rocket engine (tanks, turbo machinery, feed lines, valves, gas generator/pre-burner, combustion chamber, etc.) are modelled to study both steady state and transient phases.

To detect low frequency instabilities it is important to properly model the characteristic times associated with both liquid and gases injected into the combustion chamber. Chug instabilities arise when the total time lag is such that oscillations in the rate of propellant injection are coupled with the corresponding oscillations of the chamber pressure. Different studies have been performed during the past decades to investigate which time lag is the most important for the generation of low frequency instabilities. Most of the proposed chug models focused on steady linear analyses based on the linearisation of the continuity equation in the combustion chamber. In the attempt of explaining damages caused by chugging in LRE, Gunder and Friant [3] introduced the time lag concept. They linearised the energy conservation equation and solved the relative delay differential equation. The time lag approach was then further extended by Summerfield [2], that retrieved a stability analysis from the linearisation of the continuity equation introducing the influence of the feed lines. He investigated the effects of feed lines, injector pressure drops and combustion chamber length on chug instabilities. A comprehensive theoretical approach based on the concept of sensitive and in-sensitive time lag was then proposed by Crocco and Cheng [5, 10, 11]. Their approach is based on the concentrated combustion model and applies to both mono-propellant and bi-propellant cases. Later, Wenzel and Szuch [4, 12, 13, 14] developed a model that incorporates a double time lag approach, hence extending the potentialities of the original Crocco's model to cases in which the two propellants have different time lags. They applied the model to analyse a system ignoring the influence of the feed lines and considering only the injector impedances. They used a useful engineering approach to express the solutions, highlighting the importance of injector pressure drops on instabilities. Recently, Casiano [1] extended the Wenzel and Szuch's model by formulating the problem in a parametric form and adding the influence of the feed lines. Beside these linear approaches a few attempts have been made in the field of non-linear analysis, mainly because the computational cost of such approaches was too high for the hardware resources. Webber's [6] work was the first attempt of performing an unsteady analysis based on non-linear equation for both the feed lines and the combustion chamber. He focused his attention towards the modelling of the liquid phase in the combustion chamber, showing the effects of droplet sizes on the amplitude of the low frequency oscillations. Bartrand [7] proposed an extension of this model, including the effects of finite rate chemistry and showed that, in the Space Shuttle Main Engine pre-burner case, the assumption of finite combustion velocity affects the magnitude of the damping rather than the frequency of oscillations.

The analyses conducted in this paper take advantage of the ESPSS feed lines modelling approach, whereas the combustion chamber component is suitably improved by means of a chug dedicated module. It is well known that the shift, or time lag, between disturbances coming from injectors and pressure response is the sum of five time lags [15], each related to a specific process taking place in the combustion chamber. The five processes are, in order: injector response, atomisation, vaporisation, mixing, combustion. It is commonly accepted that the most important delays are those related with atomisation, vaporisation and mixing [16, 7], whereas spatial droplet distribution produced by injection has been shown to have a great effect on high frequency instabilities but has a little or no effect on chugging [13]. Moreover, combustion processes can be considered fast as far as excessively slow reactions are not involved. On the

contrary, the vaporisation rate is of great importance for the evaluation of chugging and is strictly related with the atomisation process (droplet size and surface area). Therefore, we focus our attention on atomisation, vaporisation and mixing phenomena. Two different approaches are foreseen: a constant time lag approach and a variable time lag approach. In the constant approach both atomisation and mixing time lag remains unchanged throughout the entire simulation. Instead, for the variable approach both atomisation [17] and vaporisation [17, 18] will be modelled with the aim of semi-empirical correlations while mixing times will be retrieved from an empirical curve [12]. In this way the characteristic times depend on the main combustion chamber parameters, thus changing during the simulation. The chug module with the two approaches will be validated against experimental and numerical data available in open literature [19] and the differences between the two methods will be discussed.

## 2. Numerical Method and Approach

The LRE system is here modelled with EcosimPro, taking advantage of the components included in the European Space Propulsion System Simulation (ESPSS) library, for which a complete description can be found in different works [9, 20]. In the following the feed lines and injector components are briefly introduced, whereas more details are provided for the combustion chamber component. The modelling approach for the gas phase in the combustion chamber is particularly important for the scope of the present paper, because it is the frame in which the vaporisation, atomization and mixing models are included.

### 2.1 System Components

#### 2.1.1 Feed Line and Injector Components

The tube component is modelled with a quasi-one-dimensional, unsteady approach for two phase flow. The two phase flow evolution is described by the Homogeneous Equilibrium Model [21] (HEM), which is based on two assumptions: (i) thermodynamic equilibrium between the two phases, (ii) same velocity for both liquid and vapour. Hence, two mass conservation equations are foreseen, one for the total mass and one for the vapour phase, completed by a momentum and an energy equation. This approach has the advantage of being governed by a set of equations that is unconditionally hyperbolic, hence allowing the use of widespread techniques for the calculation of the interface fluxes.

Injectors and injector cavities are located between the feed lines and the combustion chamber. Injector cavities are modelled by means of a two-phase set of equations for a non adiabatic variable volume. Two mass conservation equations and one energy equation describe the evolution of the two-phase mixture. The velocity in the cavity is calculated as an average velocity retrieved from the mass flow rate of the incoming fluid (calculated by the tube component) and the fluid going to the combustion chamber (calculated by the injector component), hence no momentum equation is solved. Finally, the injector component is based on a momentum equation and calculates the mass flow across the injector in both turbulent and laminar regimes. The momentum equation accounts for concentrated pressure losses and for the inertia associated with the connected components. In this way both mass and energy fluxes are provided to the combustion chamber component.

#### 2.1.2 Combustion Chamber

The thrust chamber component consists of two sub-components. The first one is an unsteady component that models the combustion chamber up to the throat. In this part of the combustion chamber the flow is considered either in chemical equilibrium or in a “time delayed” equilibrium state, that is a simple approach mimicking finite rate models. In the present analyses we will consider the component as based on the chemical equilibrium assumption. Instead, the divergent part is a quasi-steady component that mimics the supersonic expansion by means of analytical correlations for a flow in either frozen or equilibrium conditions.

The first part of the combustion chamber is based on a quasi-one-dimensional formulation of the governing equations. The component is non adiabatic, hence allowing for thermal connections with the external components, i.e. cooling channels. The unsteady conservation equations for the gas mixture are:

**Mass**

$$A \frac{\partial \rho}{\partial t} + \frac{\partial(\rho u A)}{\partial x} = \frac{A \dot{m}_{vap}}{V} \quad (1)$$

**Momentum**

$$A \frac{\partial(\rho u)}{\partial t} + \frac{\partial[(\rho u^2 + p)A]}{\partial x} = -\frac{1}{2} \frac{d\xi}{dx} \rho u |u| A + p \left( \frac{dA}{dx} \right) \quad (2)$$

## Energy

$$A \frac{\partial(\rho E)}{\partial t} + \frac{\partial(\rho u H A)}{\partial x} = \frac{d\dot{q}_w}{dx} + \frac{A\dot{q}_{unst}}{V} \quad (3)$$

The first source term on the right hand side of the momentum equation models the friction effects, while the second is related to the effects of a variable area. The first source term in the energy equation  $\dot{q}_w$  is the heat exchanged with the combustion chamber walls while the second term  $\dot{q}_{unst}$  is the unsteady heat release introduced in the present analysis to mimic the roughness of the combustion processes. Finally, the vaporised mass source term in the continuity equation  $\dot{m}_{vap}$  is expressed according to the vaporisation models described in the next section.

The nozzle component is a quasi-steady component able to describe the evolution of a flow in a duct with variable cross section. The nozzle takes the upstream thermodynamic state from the last computational cell of the combustion chamber and characterises the throat and the diverging section conditions. The evolution of the thermodynamic state throughout the nozzle is given for either isentropic or non-isentropic conditions using algebraic correlations.

### 2.1.3 Time lag approach

Once injected into the combustion chamber, the liquid subsequently undergoes atomisation, vaporisation, mixing and combustion. The time lags of each process must hence be characterised in an appropriate way in order to catch the coupling between the feed lines and the combustion chamber. The general numerical procedure is the following: once injected in the combustion chamber the mass associated with the liquid droplet is stored in a memory array and conserved for a time equal to the total time lag (vaporisation, atomisation and mixing) associated with the droplet. After the ascribed time lag the incoming mass is injected as a vaporised source term in the continuity equation (1). Two different cases are foreseen. A first one in which the time lags are user-defined input parameters and remain constant throughout the simulation. The second case is instead based on a time delayed approach in which the vaporisation and atomisation processes are described by time lags based on chamber conditions and liquid phase properties. The mixing times are retrieved from an empirical curve [12], while the combustion is considered as sufficiently fast such that its time lag can be neglected. In the present study we decided to focus our attention towards the oxygen-hydrogen combination, hence the correlations applied to the characteristic times hold only for these propellants. In particular, the hydrogen is injected in a gaseous phase, therefore the delay characterising this propellant is only the mixing time lag. The oxidizer is instead injected in a liquid phase, hence it undergoes also vaporisation and atomisation. The time lags and the related parameters are calculated as follows. For the atomisation processes we apply the correlation [17]:

$$\tau_{atom} = 6 \cdot 10^{-4} \cdot \left(\frac{\rho_g}{\rho_l}\right)^{-0.32} We_g^{0.03} Re_l^{0.55} \frac{D_l}{u_l} \quad (4)$$

$$Re_l = \frac{u_l D_l \rho_l}{\mu_l} \quad We_g = 2\rho_g \frac{(u_g - u_l)^2 D_l}{\sigma_l} \quad (5)$$

where  $D_l$  is the liquid post inner diameter,  $u_l$  is the liquid velocity at injection,  $\sigma_l$  the liquid surface tension,  $Re_l$  the Reynolds number for the injected droplet and  $We_g$  the Weber's number for the combustion chamber gases.

Vaporisation characteristic times for droplets when in presence of convection [22, 23, 17, 24] are usually retrieved as a correction of those obtained for droplets immersed in a quiescent gaseous environment [25]. We here make use of the correlation [17]:

$$\tau_{vap} = \frac{\tau_{vap,Re=0}}{1 + 1.5\alpha} \quad \alpha = 1 - \frac{0.3p_c}{100} \quad (6)$$

$$\tau_{vap,Re=0} = \frac{D_0^2}{k} \quad k = 10^{-6} \left[ 1.01 \frac{1}{1 + MR} + 1.16 \cdot 10^{-3} (T_\infty - T_{cr})^{0.93} \right]^{0.86} \quad (7)$$

where  $p_c$  is the chamber pressure,  $MR$  the mixture ratio,  $T_\infty$  the temperature of the combustion chamber gases,  $T_{cr}$  the liquid critical temperature and  $D_0$  the initial droplet diameter. The initial droplet diameter  $D_0$  is here calculated from a correlation specifically developed for coaxial injectors and liquid oxygen [18]:

$$D_0 = C_{inj} K 127 \left(\frac{V_r}{160}\right)^{-0.93} \left(\frac{D_l}{2}\right)^{2.25} \left(\frac{D_g}{3}\right)^{-2.65} \left(\frac{R_{post}}{D_l}\right)^{-0.26} \quad (8)$$

being  $V_r$  the relative velocity between injected liquid and gases in the combustion chamber,  $D_g$  the annular gap outer diameter,  $R_{post}$  the recess length,  $C_{inj}$  a parameter taking into account the injector design ranging from 0.4 to 1.2 and  $K$  a correction factor to be applied when the experimental simulants differ from the actual propellants [26].

Table 1: Main engine parameters

$D_c$	chamber diameter [cm]	5.08
$D_{th}$	throat diameter [cm]	1.24
$L^*$	characteristic length [m]	2.31
$L_c$	chamber length [cm]	10.4
$MR$	mixture ratio [-]	5
$p_c$	chamber pressure [bar]	44.8
$\dot{m}_{ox}$	mass flow rate [kg/s]	0.249

Table 2: EcosimPro component descriptions of figure 1

Component	Description
LOX/GH2	Propellant definitions
BC_LOX/LH2	Boundary conditions for propellant lines
MOV	Main Oxidizer Valve
MFV	Main Fuel Valve
Heater_1	Heat source
Loss	Heat losses time law
Insulation_1	Thermal insulation (zero heat flux)
SZC_CHUG	Combustion chamber

### 3. Test Case

In order to prove our chugging module capabilities we refer to an experimental test case specifically developed to validate a linear double time lag model [19]. The experimental apparatus is a gaseous-hydrogen liquid-oxygen engine. The injector cavities and injector plate were specifically designed to ensure the desired pressure drops and a constant pressure in the injector cavities. The main engine parameters are described in Table 1, while the EcosimPro schematic and its component descriptions are proposed respectively in Fig.1 and in Table 2.

Following the approach proposed in the reference work [19], we define a stability map for the engine by performing simulations with variable injector pressure drops. Changing both fuel and oxidizer pressure drops it is possible to create a map in which stable and unstable regions can be identified. The combustion inside the chamber is a rough process, hence spontaneous oscillations in the pressure values are usually experienced. In order to mimic these oscillations and trigger the instabilities we decided to introduce an unsteady term in the energy equation. The purpose is here to reproduce the unsteady heat release associated with combustion, phenomena commonly identified as the triggering mechanism for combustion instabilities. The unsteady heat release term is modelled by means of a broadband signal. The broadband reported in eq.(9) precisely selects the frequencies of interest and, in the hypothesis of small perturbations, enables the analysis of the chamber response at the selected frequencies

$$\dot{q}_{unst} = \dot{q}_0 \cdot \left[ 1 + \epsilon \sum_{i=1}^N \sin(2\pi(f_0 + i\Delta f)t) \right] \quad (9)$$

being  $f_0$ ,  $\Delta f$ ,  $N$ ,  $\epsilon$  respectively the first excited frequency, the sampling frequency, the number of samples and the magnitude of the perturbation. In this way frequencies between  $f_0$  and  $f_0 + N\Delta f$  are excited every  $\Delta f$ . The steady value  $\dot{q}_0$  is selected to mimic the losses measured in the combustion facility in order to match the experimental combustion efficiency  $\eta_{c^*} = 0.75$ . The final value is kept constant throughout all the simulations and equal to  $\dot{q}_0 = 1.933$  MW. This gives a low value of the combustion chamber temperature  $T_c = 2038$  K, as expected from the low combustion efficiency.

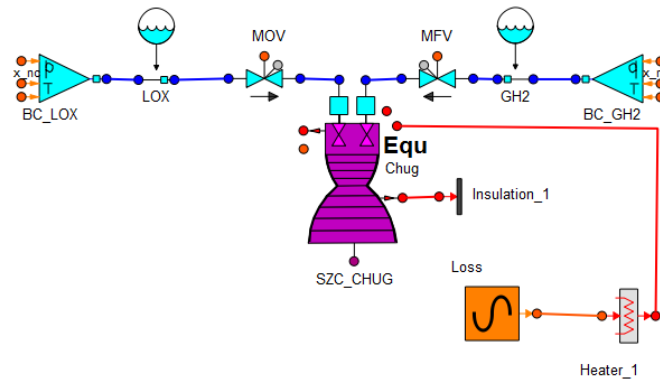


Figure 1: System schematic for chug analysis

Pressure boundary conditions on both feed lines are changed throughout the simulations to ensure constant mass flow rates at steady state for different injectors pressure drops. It is worth saying that EcosimPro has its own component to ensure a constant mass flow rate in the feed line. This component could have been replaced in the schematic in Fig.1 to both main oxidizer and main fuel valves, thus imposing the desired mass flow rates. However, it has been verified that the use of this component adds an artificial damping preventing the system from falling in any instability regime, hence failing in detecting chugging instabilities.

## 4. Results

### 4.1 Constant Time Lag

In the case of constant time lag the characteristic times are imposed as user input data. The values here used are inferred from the reference work [19]. In particular, the atomisation time is in this case neglected and the vaporisation lag is retrieved from the length needed to vaporise the 50 percent of the injected liquid mass  $l_{50}$  [27]. Once the length is known the vaporisation time is calculated as the ratio of the vaporisation length and the injection velocity of the liquid,  $\tau_{vap} = l_{50}/v_{inj}$ . Calculations for the examined engine configuration give a value of  $\tau_{vap} = 4.4$  msec. The mixing and reaction times are then retrieved from the neutral stability condition. Starting from the experimentally observed chugging frequency at high injector pressure drops of 68 Hz, the phase requirement for stability gives an overall time delay of  $\tau_{tot} = 6.7$  msec. Hence, the mixing time lag is obtained as a difference of the previously calculated delays  $\tau_{mix} = 2.23$  msec.

The results obtained with the chugging module are here compared to those coming from both experimental evidences and Szuch's double time lag model. The tests are performed by fixing one of the two injector pressure drops, varying the other and measuring the response in terms of chamber pressure. The Root Mean Square (RMS) of the oscillations of the measured signal is then extracted and presented in terms of percentage values with respect to the mean chamber pressure, while the frequency content is analysed by means of the signal Power Spectral Density (PSD). In this way we are able to retrieve both amplitude of the response and characteristic frequencies of the system. It is worth noticing that the amplitude of the chamber response is strictly related to the amplitude of the input perturbation, that in turn affects the definition of the stability region. In the experimental campaign here referenced [19] a 10 percent RMS criterion was used to identify stable and unstable operating regimes. In particular, all the configurations that give a pressure signal with a RMS exceeding the 10 percent of the mean chamber pressure value are considered unstable, while all the others can be considered stable. The same consideration applies to the results coming from the simulation performed by Szuch and Wenzel for the double time lag model, where they imposed a white noise signal on the mean combustion products values to mimic the roughness of the combustion processes. The white noise input perturbation affects the entire spectrum, but we are just interested in the lower part of the spectrum, since chug instabilities are associated with low frequencies. This consideration justifies our choice of a broadband input. In the subsequent analyses, the broadband signal in eq.(9) is characterised by the values  $f_0 = 0$  Hz,  $\Delta f = 5$  Hz,  $N = 20$  and  $\epsilon = 0.01$ , with reference to eq.(9). In Fig.2 we compare results obtained with the chug module for both single and double time lag cases against those presented by Szuch and Wenzel for both analog and experimental cases. In the single time lag case only the injected liquid droplets are included in the chug module and are introduced with a time lag equal to the sum of vaporisation, atomisation and mixing delays into the continuity equation. In the double time lag case a delay equal to the mixing time lag is also imposed to the injected gaseous fuel. Results are obtained imposing a constant fuel injector pressure drop of  $Dp_f/p_c = 0.5$ , varying the oxidizer injector pressure drops and measuring the pressure response in the chamber. Both of our methods (single and double time lag) are able to catch the increase in the RMS pressure values associated with the reduction of pressure drops. This is a well known behaviour since increasing the injector pressure drops has a stabilising effect on the overall system and decouples the oscillations in the feed lines of liquid oxygen from those in the combustion chamber, thus preventing the system from falling in the unstable regime. The predominant frequency resulting from the PSD in the single time lag case (55 Hz) differs from the one of the double time lag case (60 Hz). The result of this prediction has to be discussed considering that the broadband sampling frequency  $\Delta f$  introduces an uncertainty of  $\pm 5$  Hz. The difference between the two approaches is related to the fact that in the double time lag also the fuel feed line couples dynamically with the oscillations in the combustion chamber, thus changing the phase relation between lines and combustion chamber. However, the small registered shift indicates that the chug frequency is mostly determined by the liquid oxygen, since the greater time lag is associated to this propellant and it dominates the phase relation between oscillations. Furthermore, in the single time lag case the RMS values are always smaller than those of the double time lag approach and a plateau is observed between  $Dp_f/p_c = 0.6$  and  $Dp_f/p_c = 0.7$ . Furthermore, both time lag approaches give RMS values of the chamber pressure that are lower than those registered in the reference test cases. This is related with the amplitude  $\epsilon$  of the broadband in eq.(9).

In Fig.3 we show the influence of the amplitude of the broadband signal on the chamber pressure. Increasing the

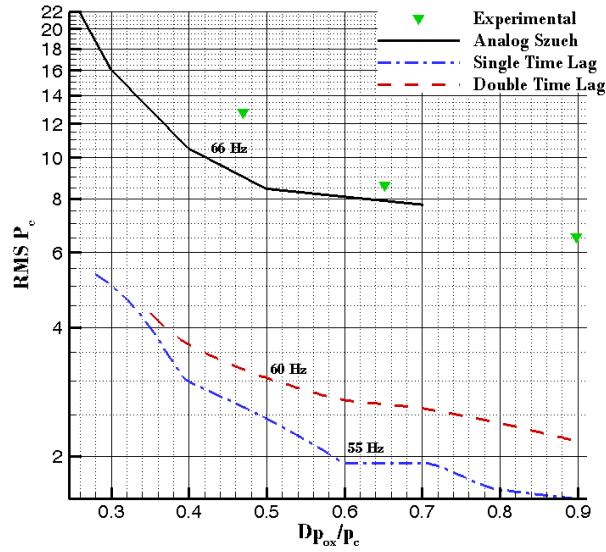


Figure 2: Double vs. single time lag,  $Dp_{fu}/p_c = 0.5$ : percentage chamber pressure RMS vs oxidizer injector pressure drops

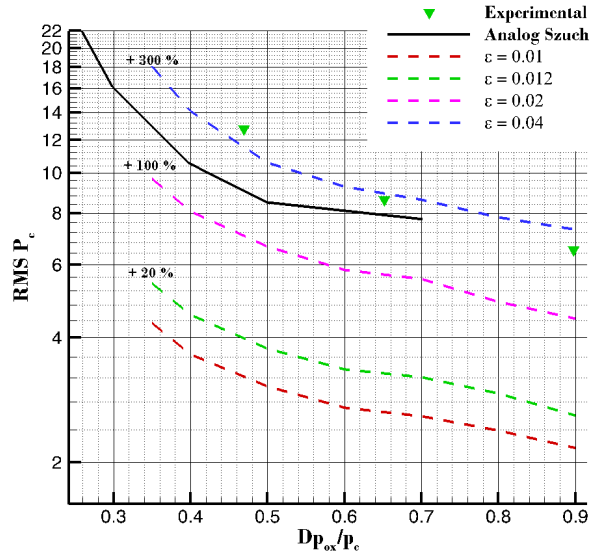


Figure 3: Double time lag,  $Dp_{fu}/p_c = 0.5$ : percentage chamber pressure RMS vs oxidizer injector pressure drops at different input amplitudes

amplitude of the input perturbation increases the chamber pressure response and the values reported in the reference paper are approached. It must be noted that in our double time lag model the inherent stability limit is placed before the one reported in the reference work, and this is true for all the selected level of amplitudes.

In both single and double time lag approaches the system response to the broadband can be divided in three zones characterised by a different behaviour in terms of both frequency content and signal shape. In Fig.4 and Fig.5 we show respectively the time evolution of chamber pressure and the PSD of the same variable in the case the double time lag model. Similar trends have been noted in the single time lag case. At high oxidizer pressure drops the chamber pressure signal is proportional to the input signal, with a frequency content characterised by a small peak at 60 Hz. When the oxidizer injector pressure drop is reduced, the 60 Hz frequency becomes predominant and the pressure signal assumes a characteristic damped behaviour. At  $Dp_{fu}/p_c = 0.35$  the system is still able to damp the perturbation and to prevent

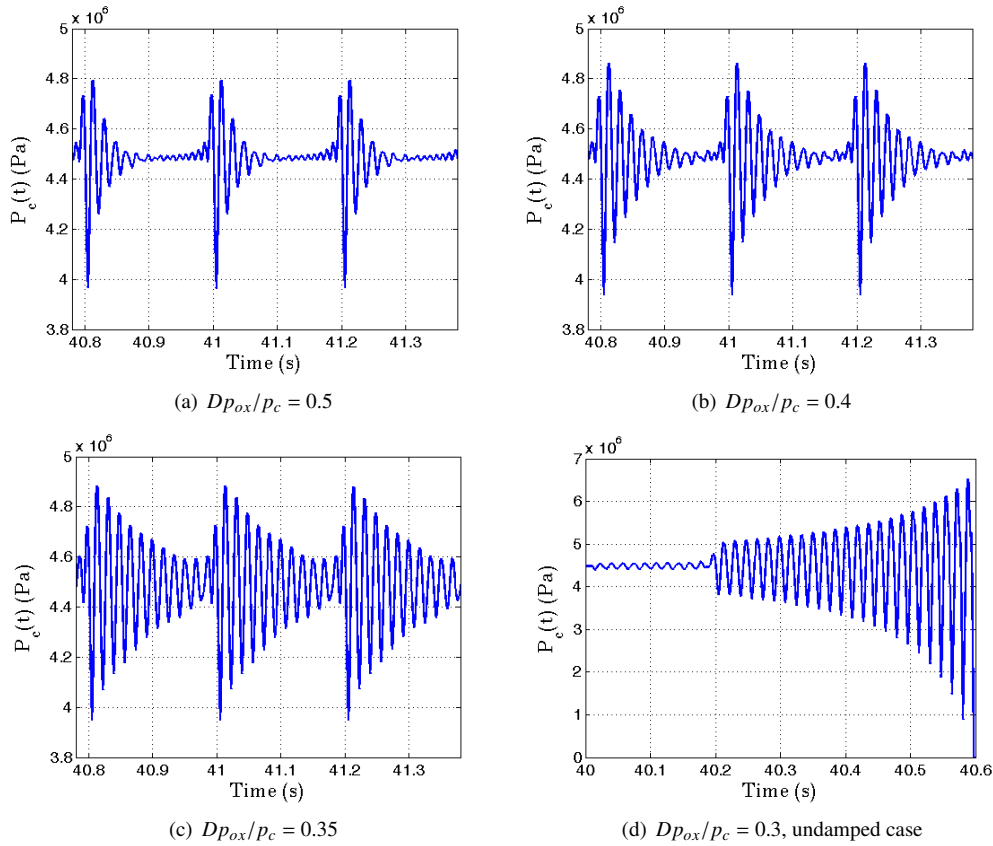


Figure 4: Double time lag,  $Dp_{fu}/p_c = 0.5$ : time evolution of the chamber pressure response at different oxidizer injector pressure drops

the oscillations from growing. No limit cycle is observed and the chamber pressure oscillations are damped and tend to disappear if not sustained with the input broadband signal. Below the inherent stability limit,  $Dp_{ox}/p_c < 0.35$  the system is not able to damp the oscillations and the pressure response grows to high values causing the simulation to fail, as shown in Fig.4(d). In this case the system oscillates at the system characteristic frequency of 60 Hz and the oscillations are self sustained even when the input is deactivated. Furthermore, the system oscillates at the characteristic frequency regardless the frequency content of the input, that is: the system diverges with a 60 Hz frequency even if a monochromatic input of 20 Hz is imposed for a short time and then deactivated.

In Fig.6 we propose the stability map obtained for the selected test case with the double time lag model. The results are compared with the boundaries identified by Szuch et al. It is worth noticing that the 10% RMS criterion is a practical criterion [19] to define a stability boundary. Since this criterion is directly related with the amplitude of the input perturbation we choose to use a value of  $\epsilon = 0.01$  throughout the simulations and we make use of also a 50% RMS criterion to identify the stability boundary. The boundaries identified with both the 10% and the 50% RMS criterion are shown with white dashed lines, whereas the reference results are shown in black. When the boundary is identified with the 50% criterion the reversal in slope at lower  $Dp_f/p_c$  is well matched, and, differently from the analog referenced computer program, no unstable regimes are identified at high oxidizer injector pressure drops when the fuel injector pressure drop is lower than 0.15. The authors of the reference paper [19] justified this discrepancy assuming a back-flow in the fuel injector element at lower pressure drops, however this back-flow has never been registered in our simulations. Furthermore, the reversal in slope is due to the intersection of two boundaries: the low frequency boundary and the high frequency boundary. At lower fuel injector pressure drops characteristic system frequencies ranging from 147 to 210 Hz were observed both experimentally and numerically [19]. In order to investigate the presence of this second frequency we extend the broadband input up to 200 Hz, measuring the resulting chamber pressure signal. With this input we are able to register a higher frequency content of 175 Hz that is well within the limits of the aforementioned high frequency range, giving a ratio of 2.9 between higher and lower frequency that well approximates the 2.7 ratio noted during multi-mode oscillations [28]. However, when the system enters its unstable regime at fuel injector pressure drops lower than 0.15 the pressure diverges with a frequency of 60 Hz, thus meaning that in our analysis the 60 Hz frequency is the characteristic frequency of the system even at lower fuel injector pressure

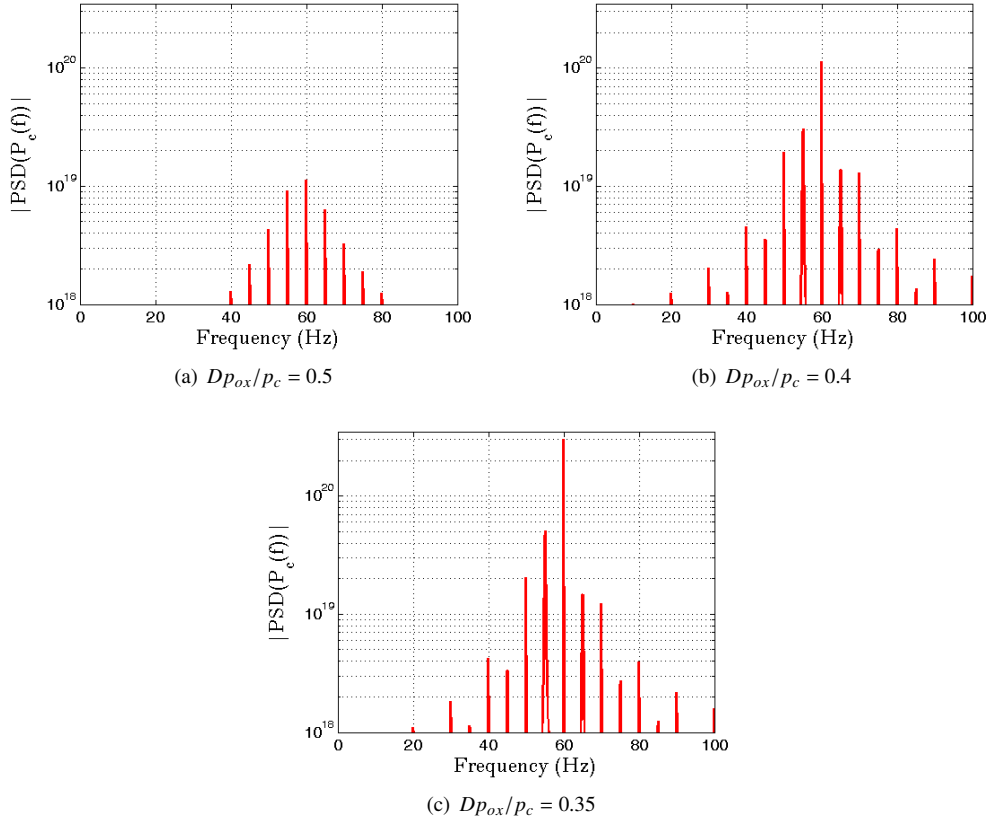


Figure 5: Double time lag,  $Dp_{fu}/p_c = 0.5$ : Power Spectral Density at different oxidizer injector pressure drops

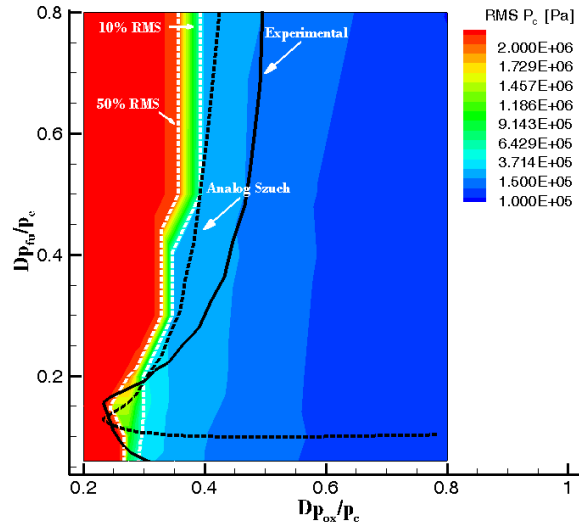


Figure 6: Stability map. Chamber pressure RMS dependency on oxidizer and fuel injector pressure drops

drops. Furthermore, the PSD shows that the power associated to the 60 Hz component is always greater than the one at 175 Hz, explaining the predominance of the first frequency throughout the entire stability map.

For the constant time lag model we finally investigate the influence of different input amplitudes for various fuel pressure drops. The scope is here to investigate how the stability boundaries are influenced by the level of the input perturbations. In particular, we present the results obtained when the amplitude  $\epsilon$  in eq.(9) is increased of 20% and

100%. The results in terms of the RMS of chamber pressure oscillations are then normalised with those obtained with the reference case  $\epsilon = 0.01$  and then divided by the respective percentage increment. The aim of these normalisations is to investigate whether the response of the system is proportional to the input increment. Results shown in Fig.7 highlight that for different fuel injector pressure drops the system response has a maximum at higher oxidizer injector pressure drops. Thus evidences that, when the amplitude of the input perturbation is increased, the boundary of the stability regions experience a greater shift if placed in the high pressure drop region and a smaller shift if placed in the low pressure drop region. Furthermore, increasing the input amplitude does not affect the characteristic frequency of the system that remains equal to 60 Hz.

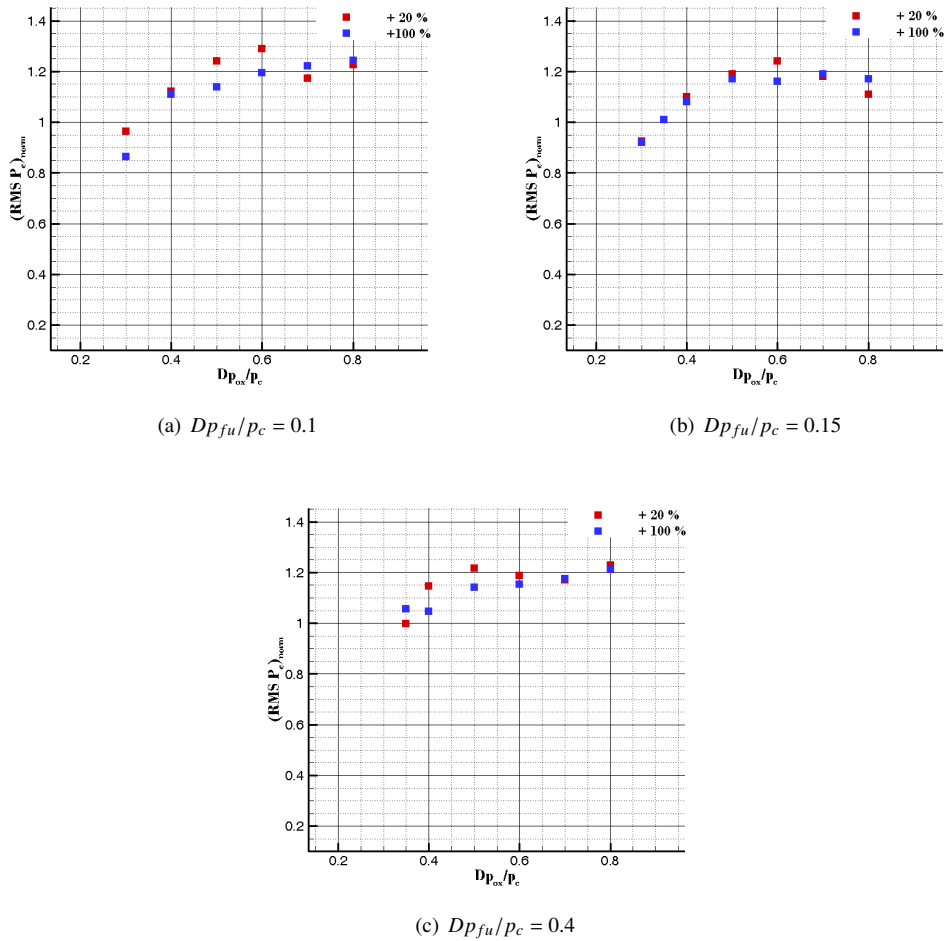


Figure 7: Double time lag: normalised RMS of the chamber pressure signal for different input amplitudes at various fuel injector pressure drops

## 4.2 Variable Time Lag

The time lags are typically not known a priori, hence when a new engine configuration is analysed it is useful to retrieve the needed time delays by means of dedicated correlations. The benefit of introducing correlations for the time lags is dual: on one hand the user is free from any a priori calculation of the time lags, and on the other the time lags become time varying functions of the combustion chamber variables. In this frame we introduced semi-empirical correlations to calculate atomisation and vaporisation time lags, considering the combustion time lag as infinitely fast. As mentioned in Section 2.1.3 the time lags for atomisation and vaporisation processes are retrieved respectively from eq.(6) and eq.(4), while mixing time lag is evaluated from an empirical curve [12]. A double time lag approach is still used, hence the total time delay for the liquid oxygen is retrieved as the sum of vaporisation, atomisation and mixing time lags while only the mixing time lag characterises the gaseous hydrogen delay. Some considerations must be done on the correlation for the initial droplet diameter eq.(8). Due to the lack of informations on the experimental set up, the recess

of the oxidizer post  $R_{post}$  is an unknown data that greatly influences the value of the initial droplet diameter and in turn the vaporisation time. Furthermore, the injector coefficient  $C_{inj}$  can range from 0.4 to 1.2, thus changing significantly the vaporisation delay. We know, from experimental evidences [25], that the initial droplet diameters of liquid oxygen at high pressure range from  $60 \mu\text{m}$  to  $160 \mu\text{m}$ . Hence, stated that the recess length is a construction parameter of a few millimetres, we decided to use a combination of values that, while respecting all of the aforementioned constraints, ensures a vaporisation time lag at steady state equal to  $\tau_{vap} = 4.4 \text{ msec}$ . In this way we retrieved a time varying time lag matching with the steady state values suggested from the experimental evidences [19]. The reference chosen working point is characterised by the values  $R_{post} = 0.003 \text{ m}$ ,  $C_{inj} = 0.4$  and  $K = 0.121$ .

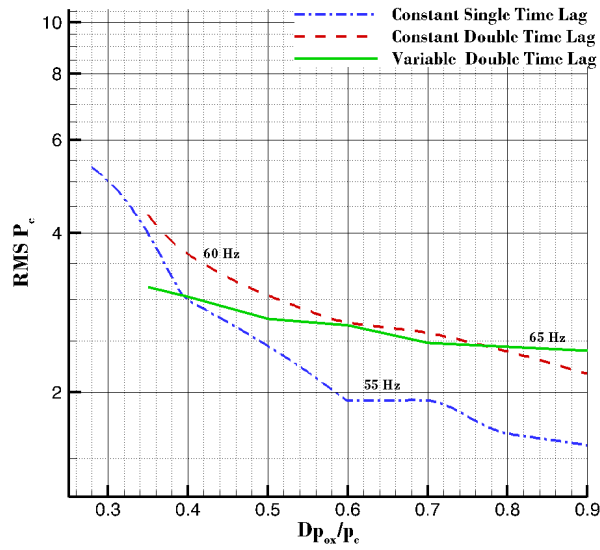


Figure 8: Double time lag,  $Dp_{fu}/p_c = 0.5$ : chamber pressure RMS vs oxidizer injector pressure drops

The time dependent approach is then tested on the engine configuration presented in Section 3. A single fuel injector pressure drop is here investigated at different oxidizer injector pressure drops, and results are compared in Fig.8 against those obtained for the constant time lag approaches. The variable time lag approach gives a pressure response in line with the fixed time lag approaches. The amplitudes of the oscillations are less influenced by the reduction of the the oxidizer injector pressure drop and a smoother growth is observed while passing from high to low injector pressure drops. Finally, the time dependent time lags give a characteristic frequency of 65 Hz approaching the experimental measured value of 68 Hz, with an uncertainty of 5 Hz related to the broadband sampling frequency. Furthermore, the tendency of the double time lag approach in anticipating the insurgence of the inherent stability limit is here confirmed. In fact, in both double time lag approaches the limit moves towards higher oxidizer pressure drops and for pressure drops lower than  $Dp_{ox} = 0.23$  the system is not able to damp the imposed perturbation, causing the chamber pressure to diverge with a frequency equal to the characteristic frequency of 65 Hz.

## 5. Conclusion

A double time lag model to investigate low frequency combustion instabilities has been implemented in the system analysis tool EcosimPro and tested against experimental results. Two different approaches have been followed: a constant time lag and a variable time lag approach based on semi-empirical correlations. The constant time lag approach demonstrates its effectiveness in reproducing the unstable behaviour associated with the reduction of the injector pressure drops, in particular no stable region has been identified for an oxidizer injector pressure drop lower than 0.35. The chug module is able to reproduce the reversal in slope of the stability boundary at low injector pressure drops, thus giving a result close to the one found in the reference experiment. However, it must be pointed out that the location of the stability boundary depends on both the criterion used to identify a regime as unstable and on the amplitude of the input perturbation. We demonstrated that the boundaries can be easily shifted by changing the amplitude of the input perturbation and that this shift slightly depends on the selected injector pressure drop. Hence, in the computer program the boundary of the stability limits can be shifted in accordance with the level of the input perturbation. Despite its

simplicity, the constant time lag approach is able to predict a characteristic frequency close to the one measured experimentally. The registered high frequency content is characterised by a ratio of higher to lower frequency close to 2.7, as reported in open literature. It must be noted that in our case the high frequency does not dominate the dynamic response at low fuel injector pressure drops and the instability is still related to the first characteristic frequency. The inability in catching the suppression of the first mode could be related with a lack of data for parameters that usually influence the damping effect and the associated system impedances, i.e. injector cavity volumes and length of the feed lines. These values have been selected to be consistent with the downstream injector geometry and mass flow rates, but a fine tuning with the real data is unavoidable to improve the results of the model. The chugging module has been finally tested with semi-empirical correlations to retrieve characteristic delays that, depending on the main flow variables, are time dependent functions of the main combustion chamber parameters. With the implemented correlations the model retrieves a lower characteristic frequency that better approximates the experimental results, while the predicted RMS values of the chamber pressure remain comparable with those obtained with the constant double time lag model.

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