

UNIVERSITAT POLITÈCNICA DE VALÈNCIA  
DEPARTAMENTO DE MÁQUINAS Y MOTORES TÉRMICOS

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UNIVERSITAT  
POLITÈCNICA  
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MODELLING AND OBSERVATION OF EXHAUST  
GAS CONCENTRATIONS FOR DIESEL ENGINE  
CONTROL

PHD DISSERTATION  
Presented by  
David Blanco-Rodríguez  
Advised by  
Dr. Carlos Guardiola

Valencia, September 2013



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MODELLING AND OBSERVATION OF EXHAUST GAS  
CONCENTRATIONS FOR DIESEL ENGINE CONTROL

Presented by: David Blanco-Rodriguez  
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*A mis abuelos Manolo y Feli,  
a Mikel, Ama y Aita.*

*Shine on you crazy diamond.*



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**Resumen.** La presente Tesis Doctoral estudia la observación en tiempo real de la concentración en el colector de escape de los óxidos de nitrógeno ( $\text{NO}_x$ ) y del dosado ( $\lambda^{-1}$ ) en motores diesel sobrealimentados. Para ello se fusionan dos fuentes de información diferentes:

- Sensores capaces de proporcionar una medida de dichas variables, y
- modelos orientados a control que estiman estas variables a partir de otras medidas del motor.

El trabajo parte de la evaluación de la precisión de los sensores, realizada mediante la comparación de su medida con la proporcionada por equipos analíticos de alta precisión, que son usados como estándares de calibración estática. También se desarrollan en la Tesis métodos para la calibración de la dinámica del sensor; dichos métodos permiten identificar un modelo de comportamiento del sensor y revelar su velocidad de respuesta y retraso asociados. En general, estos sensores demuestran ser precisos pero relativamente lentos.

Por otra parte, se proponen modelos para la estimación de  $\text{NO}_x$  y  $\lambda^{-1}$ . Estos métodos, basados en relaciones físicas, tablas de parámetros y una serie de correcciones, emplean las medidas proporcionadas por otros sensores con el fin de proporcionar una estimación de las variables de interés. Los modelos permiten una estimación muy rápida capaz de reproducir las características dinámicas, pero resultan afectados por efectos de deriva que comprometen su precisión.

Con el fin de aprovechar las características dinámicas del modelo y mantener la precisión en estado estacionario del sensor, se proponen técnicas de fusión de la información basadas en la aplicación de filtros de Kalman (KF). En primer lugar, se diseña un KF capaz de combinar ambas fuentes de información y corregir en tiempo real el sesgo entre las dos señales. Posteriormente, se estudia la adaptación en tiempo real de los parámetros del modelo con el fin de corregir de forma automática los problemas de deriva asociados al uso de modelos.

Todos los métodos y procedimientos desarrollados a lo largo de la presente Tesis Doctoral se han aplicado de forma experimental a la estimación de  $\text{NO}_x$  y  $\lambda^{-1}$ . De forma adicional, la Tesis Doctoral desarrolla aspectos relativos a la transferencia de estos métodos a los motores diesel de serie.

**Resum.** La present Tesis Doctoral estudia l'observació en temps real de òxids de nitrogen ( $\text{NO}_x$ ) i del dosat ( $\lambda^{-1}$ ) a motors Diesel sobrealimentats. Per tal d'assolir suficient precisió i velocitat es combinen dues fonts de informació:

- Sensors que proporcionen una mesura de les variables, i
- models orientats a control basats en variables més directes del motor.

El treball desenvolupat parteix d'estudis per avaluar la precisió dels sensors. Aquests estudis tracten de comparar la mesura dels sensors amb la proporcionada per equips analítics d'alta precisió, equips utilitzats com estàndards de calibratge estàtic. A partir del calibratge estàtic es desenvolupen diversos mètodes per el calibratge dinàmic dels sensors, aquests mètodes permeten identificar un model de comportament del sensor, simulant la seva velocitat de resposta.

En general, els sensors demostren ser precisos però relativament lents. Per millorar els temps de resposta de les mesures, a una segona part de la tesis es proposen models orientats a control per la estimació de  $\text{NO}_x$  i  $\lambda^{-1}$ . Aquests mètodes, basats en relacions físiques, taules de paràmetres i una sèrie de correccions, utilitzen mesures d'altres sensors més directes per tal de proporcionar una estimació de les variables d'interès. La estimació obtinguda als models, a pesar de ser molt ràpida, sol estar afectada per efectes de deriva del model, raó per la que la precisió pot veures greument afectada.

Finalment, per tal d'aprofitar les característiques dinàmiques del model i mantenir la precisió en estat estacionari del motor, es proposen noves tècniques de fusió de la informació basades en l'aplicació de filtres de Kalman (KF). En primer lloc, es dissenya un KF capaç de combinar models i sensors, per corregir en temps real la desviació de les dos senyals. I posteriorment, amb la finalitat d'eliminar els problemes de deriva associats al model, s'estudia la adaptació en temps real dels paràmetres del model.

Tots els mètodes i procediments desenvolupats a la present Tesis Doctoral, han sigut aplicats de forma experimental a la estimació de  $\text{NO}_x$  i  $\lambda^{-1}$ . De forma addicional, la Tesis Doctoral desenvolupa aspectes relatius a la possible implementació dels mètodes a motors diesel de sèrie.



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**Abstract.** The dissertation covers the problem of the online estimation of the diesel engine exhaust concentrations of the nitrogen oxides ( $\text{NO}_x$ ) and the fuel-to-air ratio or richness ( $\lambda^{-1}$ ). Two information sources are utilised:

- On-board sensors for measuring  $\text{NO}_x$  and  $\lambda^{-1}$ , and
- control oriented models in order to predict  $\text{NO}_x$  and  $\lambda^{-1}$ .

The evaluation of the static accuracy of these sensors is made by comparing the outputs with a gas analyser, while the dynamics are identified on-board by performing step-like transitions on  $\text{NO}_x$  and  $\lambda^{-1}$  after modifying ECU actuation variables. Different methods for identifying the dynamic output of the sensors are developed in this work; these methods allow to identify the response time and delay of the sensors if a sufficient data set is available. In general, these sensors are accurate but present slow responses.

Afterwards, control oriented models for estimating  $\text{NO}_x$  and  $\lambda^{-1}$  are proposed. Regarding the  $\lambda^{-1}$  prediction, the computation is based on the relative fuel-to-air ratio, where the fuel comes from an ECU model and the air mass flow is measured by a sensor. For the case of  $\text{NO}_x$ , a set-point relative model based on look-up tables is fitted for representing nominal engine emissions with an exponential correction based on the intake oxygen variation. Different correction factors for modeling other effects such as the thermal loading of the engine are also proposed. The model is able to predict  $\text{NO}_x$  fast with a low error and a simple structure.

Despite of using models or sensors, the model drift and sensor dynamic deficiencies affect the final estimation. In order to solve these problems, data fusion strategies are proposed by combining the steady-state accuracy of the sensor and the fast estimation of the models by means of applying Kalman filters (KF). In a first approach, a drift correction model tracks the bias between the model and the sensor but keeping the fast response of the model. In a second approach, the updating of look-up tables by using observers is coped with different versions based on the extended Kalman filter (EKF). Particularly, a simplified KF (SKF) allows to observe the parameters with a low computational effort.

Finally, the methods and algorithms developed in this work are combined and applied to the estimation of  $\text{NO}_x$  and  $\lambda^{-1}$ . Additionally, the dissertation covers aspects relative to the implementation of the methods in series diesel engines.



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Mi vida siempre ha estado unida de una manera u otra a correr. Parece una coincidencia, pero hoy hace apenas 4 años y unos 2 meses que comencé mi última aventura, como si de un paralelismo con la maratón (casi 42 kms) se tratase. Ha sido sin duda una carrera de fondo, en la que he disfrutado de los primeros kilómetros, los primeros años, he sufrido y he tenido que trabajar mucho durante bastante tiempo (sufriendo lo que en el argot atlético se conoce como "El Muro"), pero a la postre y ahora que veo la recta final, ya sólo me queda esprintar y entrar sonriendo por todo lo vivido. Ahora todas esas noches en vela, todos los fines de semana sin apenas descansar, y todos esos encierros en mi despacho sonando Pink Floyd (Another Brick on the Wall?), me parecen una anécdota, una experiencia que siempre había querido vivir. En esta parte del documento no quisiera olvidarme de todas "las liebres" que durante estos años me han ayudado a seguir adelante.

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*Sopelana, a 3 de Septiembre de 2013*

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# Nomenclature

## Acronyms

AR	Autoregressive
ARMAX	Autoregressive moving-average with exogenous inputs
ARX	Autoregressive moving-average
AT	After-treatment
BTDC	Before the top dead centre
CAD	Crank angle degree
CAE	Computer aided engineering
CAN	Controller area network
CARB	California air resources board
CAS	Crank angle solved
CFD	Computational fluid dynamics
CI	Compression ignition
CO	Carbon monoxide
CO <sub>2</sub>	Carbon dioxide
COM	Control oriented model
CR	Common-rail
DDM	Data driven model
DENO <sub>x</sub>	After-treatment system for reducing NO <sub>x</sub>
DF	Data fusion
DI	Direct injection
DICI	Direct injection compression ignition
DOC	Diesel oxidation catalyst
DPF	Diesel particulate filter
E&FM	Emptying and filling models
ECU	Electronic control unit
EGO	Exhaust gas oxygen
EGR	Exhaust gas recirculation
EKF	Extended Kalman filter
EMS	Engine management strategy
EPA	Environmental protect agency
ET	Energising time

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ETK	Ethernet based ECU connection
FDI	Failure detection and isolation
FTP	Federal test procedure
GDP	Gross domestic product
GPS	Global positioning system
HC	Hydrocarbons
HCCI	Homogeneous charge compression ignition
HC-SCR	Hydrocarbons selective catalyst reduction
HD	Heavy-duty
HIL	Hardware in the loop
HPEGR	High pressure exhaust gas recirculation
HW	Hammertein-Wiener
IC	Internal combustion
ID	Injection duration
KF	Kalman filter
LD	Light duty
LEV	Low emission vehicle
LNT	Lean NO <sub>x</sub> trap
LPEGR	Low pressure exhaust gas recirculation
LPV	Linear parameter varying
LS	Least squares
LTC	Low temperature combustion
MI	Main Injection
MIL	Malfunction indicator lamps
MPC	Model predictive control
MV	Mean value
MVEM	Mean value engine model
NEDC	New European driving cycle
NH <sub>3</sub>	Ammonia
NN	Neural network
NO	Nitrogen monoxide
NO <sub>2</sub>	Nitrogen dioxide
NO <sub>x</sub>	Nitrogen oxides (NO + NO <sub>2</sub> )
O <sub>2</sub>	Oxygen
OBD	On-board diagnostics
OEM	Original equipment manufacturer
PEMS	Portable emissions measurement system
PM	Particulate matter
PWA	Piece-wise affine
RT	Real time
SC	Supercharger
SCCI	Stratified charge compression ignition
SCR	Selective catalyst reduction
SI	Spark ignition

SIL	Software in the loop
SKF	Simplified Kalman filter
SOE	Start of energising
SOI	Start of injection
SSKF	Steady-state Kalman filter
TC	Turbocharger
TCCI	Turbocharger compression ignition
TCD	Turbo compounding
TDC	Top dead centre
TWC	Three way catalyst
UEGO	Universal exhaust gas oxygen
VGC	Variable geometry compressor
VGT	Variable geometry turbine
VVT	Variable valve timing
WAM	Wave action model
WG	Waste-gate
ZrO2	Zirconium oxide (zirconia)

**Engine positions**

<i>adm0</i>	Admission downstream of the first air filter installed in the test bench
<i>adm1</i>	Admission downstream of the series air filter
<i>boost</i>	Downstream of the engine intercooler
<i>egr</i>	EGR runner upstream of the intake junction
<i>int</i>	Intake runner
<i>exh</i>	Exhaust runner
<i>inl</i>	Turbine inlet
<i>out</i>	Turbine outlet
<i>wic</i>	Water (coolant) in the engine intercooler
<i>cyl</i>	In-cylinder
<i>crank</i>	Engine crankshaft
<i>CR</i>	Common-rail

**Engine variables**

<i>u<sub>wg</sub></i>	Waste-gate valve position command 0-open, 100-closed	%
<i>u<sub>egr</sub></i>	EGR valve position command 0-open, 100-closed	%
<i>u<sub>lp</sub></i>	Low pressure EGR valve position command	%
<i>u<sub>hp</sub></i>	High pressure EGR valve position command	%
<i>u<sub>prail</sub></i>	Rail pressure command	bar
<i>u<sub>soi</sub></i>	Start of injection of the main injection command	°BTDC
<i>u<sub>id</sub></i>	Total injection duration command	μs
<i>u<sub>α</sub></i>	Pedal position command	%
<i>u<sub>vgt</sub></i>	VGT position command	%
<i>u<sub>it</sub></i>	Intake throttle valve command	%
<i>u<sub>bp</sub></i>	Backpressure valve command	%

$\dot{m}$	Mass flows	kg/h
$m$	Mass per stroke	mg/str
$p$	Pressures	bar
$T$	Temperatures	K (°C)
$\dot{m}_a$	Air mass flow	kg/h
$\dot{m}_a^*$	$\dot{m}_a$ set-point reference for controller	kg/h
$\dot{m}_f$	Injected fuel mass flow	kg/h
$m_f$	Injected fuel mass	mg/str
$t_{mi}$	Main injection duration	$\mu$ s
$t_{id}$	Total injection duration	$\mu$ s
$p_{boost}$	Boost pressure	bar
$p_{boost}^*$	$p_{boost}$ set-point reference for controller	bar
$T_{cool}$	Engine coolant temperature	K (°C)
$T_{exh}$	Exhaust temperature	K (°C)
$T_{int}$	Intake temperature	bar
$p_{exh}$	Exhaust pressure	bar
$p_{int}$	Intake pressure	bar
$T_{wic}$	Temperature of the water of the intercooler	K (°C)
$T_{egr}$	EGR temperature	K (°C)
$n$	Engine speed	rpm
$p_{rail}$	Rail pressure	bar
$M_e$	Engine torque	Nm
$p_{cyl}$	In-cylinder pressure	bar
$I_p$	Current output from the UEGO sensor	mA
$\eta_v$	Volumetric efficiency	-
$EGR$	EGR rate	%
$z$	Sensor measurements	-
$x$	State-vector	-
$u$	Input	-
$\hat{x}$	State observation	-
$x_r$	Actual signal	-
$x_f$	Filtered signal	-
	in order to compare with sensor outputs	-
$\Theta$	Adaptive look-up table	-
$x_\Theta$	Output from interpolating the look-up table $\Theta$	-
$K$	Kalman gain	-
$K_{\infty}$	Steady-state Kalman gain	-
$K^{SS}$	Steady-state Kalman gain for the SSKF	-
$\mathbf{P}$	Covariance matrix	-
$\sigma_w^2$	Process variance	-
$\sigma_v^2$	Output variance	-

$\tau$	Sensor delay	s (or ms)
$T_{\text{NO}_x}$	Response time of the sensor	s (or ms)
$a$	Discrete response time parameter	-
$T_s$	Sample time	ms
$t$	time	s
$k$	Discrete instant	-
$s$	S-transform variable	-
$z$	Z-transform variable	-
$a_{\text{NO}_x}$	Discrete response time of the $\text{NO}_x$ output	-
$\tau_{\text{NO}_x}$	Delay of the $\text{NO}_x$ output	$\mu\text{s}$
$z_{\text{NO}_x}$	$\text{NO}_x$ output from the $\text{NO}_x$ sensor	ppm
$x_{\text{NO}_x}$	Output from the $\text{NO}_x$ model	g/h (or ppm)
$y_{\text{NO}_x}$	Filtered and delayed output from the $\text{NO}_x$ model	g/h (or ppm)
$x_{\text{NO}_x f}$	Filtered and delayed output from the $\text{NO}_x$ model in the state-space model	g/h (or ppm)
$\hat{x}_{\text{NO}_x}$	Actual $\text{NO}_x$ observation	g/h (or ppm)
$\hat{x}_{\text{NO}_x f}$	Filtered $\text{NO}_x$ observation	g/h (or ppm)
$\theta_{\text{NO}_x}$	Bias on the $\text{NO}_x$ output	g/h (or ppm)
$\hat{\theta}_{\text{NO}_x}$	Observed bias on the $\text{NO}_x$ output	g/h (or ppm)
$\Theta_{\text{NO}_x}$	Adaptive look-up table for modelling $\text{NO}_x$	g/h (or ppm)
$a_{\lambda^{-1}}$	Discrete response time of the $\lambda^{-1}$ output	-
$\tau_{\lambda^{-1}}$	Delay of the $\lambda^{-1}$ output	$\mu\text{s}$
$z_{\lambda^{-1}}$	$\lambda^{-1}$ output from the $\text{NO}_x$ sensor	-
$x_{\lambda^{-1}}$	$\lambda^{-1}$ model from the fuel-to-air ratio calculation	-
$y_{\lambda^{-1}}$	Filtered and delayed output from the $\lambda^{-1}$ model	-
$x_{\lambda^{-1} f}$	Filtered and delayed output from the $\lambda^{-1}$ model in the state-space model	-
$\hat{x}_{\lambda^{-1}}$	Actual $\lambda^{-1}$ observation	-
$\hat{x}_{\lambda^{-1} f}$	Filtered $\lambda^{-1}$ observation	-
$\theta_{\lambda^{-1}}$	Bias on the $\lambda^{-1}$ output	-
$\hat{\theta}_{\lambda^{-1}}$	Observed bias on the $\lambda^{-1}$ output	-
$\Theta_{\lambda^{-1}}$	Adaptive look-up table for modelling $\lambda^{-1}$	-





# Chapter 1

## Introduction

### Contents

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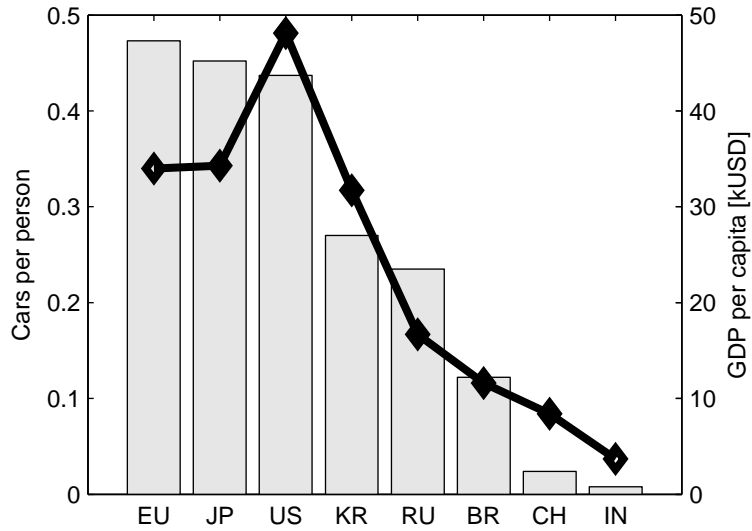
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### 1.1 Background

The evolution of transportation systems is closely linked with the degree of development of modern societies. As shown in Figure 1.1, developed countries present passenger car densities around 0.5 cars per inhabitant [1,2], while developing countries, such as China or India around 0.03 [3–5], although the economic development is expected to make this average grow. The transportation fleet is responsible for about 40% of the global fuel consumption and the expected growth will rise the number.

**The emissions regulation in diesel engines.** The world attention about environmental protection has resulted in new strict laws which establish the requirements for pollutants emissions, and therefore define priorities in the technology development [6]. Traditionally, two types of internal combustion



**Figure 1.1.** Cars per person (bars) and gross domestic product (GDP) per capita (line) in some selected countries, adapted from [5].

(IC) engines have dominated the transportation market: the spark ignition (SI) or gasoline and the compression ignition (CI) or diesel engines. Particularly in diesel engines, main pollutants are:

- Nitrogen oxides ( $\text{NO}_x$ ) are benefited from high temperatures and a lean combustion.  $\text{NO}_x$  at the engine exhaust contains NO with a proportion between 70 and 90% while the rest is mainly  $\text{NO}_2$  [7].
- Particulate matter (PM) is a complex aggregate formed by soot, hydrocarbons (HC) resulting from fuel and lubrication and other minor products. The proportion of these components varies with the engine and operating point conditions. PM emissions have an opposite trend with  $\text{NO}_x$ , and if fuel is burnt appropriately, PM concentration decreases while  $\text{NO}_x$  increases.
- Carbon dioxide ( $\text{CO}_2$ ) formation is proportional to the fuel consumption. Traditionally,  $\text{CO}_2$  emissions were not regulated, however they have been driven by the fuel economy demanded by the users.  $\text{CO}_2$  is today a main concern for manufacturers not only due to the fuel saving requirements

but due to the new regulations on CO<sub>2</sub><sup>1</sup>. The passenger cars are particularly responsible for around the 12% of the CO<sub>2</sub> emissions caused by humans [9].

- Carbon monoxide (CO) is a sub-product of the combustion (more important in SI than in CI engines) and a dangerous colourless, odourless and tasteless gas. The usual is that CO oxides at the atmosphere for forming CO<sub>2</sub>.
- Hydrocarbons (HC) are a product of an incomplete combustion of the injected fuel due to rich conditions and low temperatures that can locally be achieved inside the cylinder. Partially oxidised HC and non-burnt HC are usually included inside this group. The emissions of CO and HC in diesel engines are of less importance than NO<sub>x</sub>, PM and CO<sub>2</sub>.

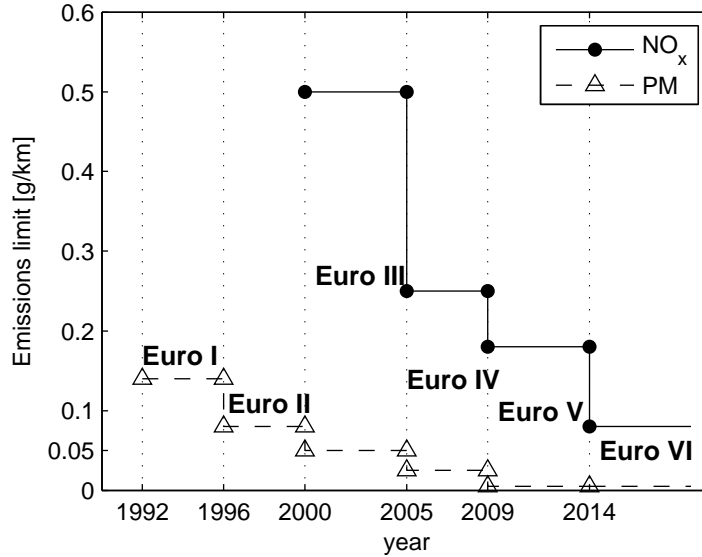
Figure 1.2 shows emissions limits for NO<sub>x</sub> and PM for the different versions of European legislation standards [8] for diesel passenger cars and light-duty (LD) vehicles. Specifically in mobile sources, diesel engines must reduce NO<sub>x</sub> emissions by 20% with regards to EURO V and 50% with regards to EURO VI if comparing with the previous EURO IV standard.

Driven by these legislations and the market demands, the diesel engine has substantially changed from the original one designed by Rudolf Diesel. The modern diesel engine comprises of variable geometry turbocharging [10] for increasing the intake air mass flow, common-rail systems for allowing multiple and controlled injections [11], an exhaust gas recirculation (EGR) [12] for reducing NO<sub>x</sub> and different after-treatment (AT) systems [13] such as the diesel particulate filter (DPF) for trapping soot, the diesel oxidation catalyst (DOC) for eliminating HC and the selective catalyst reduction (SCR) or the lean NO<sub>x</sub> trap (LNT) for removing NO<sub>x</sub>. These advances make the diesel engine a sophisticated system that requires an engine control unit (ECU) for an appropriate management and a growing complexity on calibration and implementation.

The SI engine has traditionally solved the emissions problem easier than the CI engine by installing a three way catalyst (TWC), which simultaneously reduces NO<sub>x</sub> and oxidises CO and HC. Nevertheless, the fuel efficiency is lower in SI than in CI engines. In the last years, and with the idea of taking advantage of the best properties of CI and SI engines, manufacturers and researchers have put a big effort in developing a more efficient and environmentally sustainable engine. These concepts usually lie on low temperature combustions

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<sup>1</sup>CO<sub>2</sub> emissions are regulated in Europe since the year 2009 by the European directive 443/2009 [8].



**Figure 1.2.**  $NO_x$  and PM emission limits for diesel passenger cars and light-duty vehicles (M1) from Euro I to Euro VI standards.  $NO_x$  limits were first imposed in Euro III. Euro VI is to be applied in 2014.

(LTC) [14]. Alternatively, hybrid-electric and pure electric power sources are also becoming an option for the present and future vehicle.

Whatever the engine selected (IC, electric or hybrid), the appropriate management of all subsystems merits a special focus on the ways of retrieving information about the states (i.e. variables that determine the behaviour of the engine). The work presented in this dissertation is focused on the estimation of the exhaust gas concentrations of diesel engines, whose motivation is further presented in the next section.

## 1.2 The need of information in diesel engines

The advancements in the diesel engines must be in the direction of guarantying low emissions while achieving high efficiency, performance and reliability. These purposes require not only efforts on the engine design and performance, but also in the development of reliable measurement systems to get information about the processes, which is needed for the implementation of control strategies, paying a special attention to the transient operation. In such sense,

standards also specify on-board diagnostics procedures (OBD) that, based on models and measured data, are capable of detecting faults on engine or vehicle system. Driver must be notified when emissions are out of the limits with the malfunction indicator lamps (MIL) and the ECU must correct this fault or recommend what to do for a correct and safe actuation.

The diesel engine itself may be split in two main systems: the fuel and the air paths. The fuel path is controlled by the measured rail pressure and the estimated injection fuel rate. Globally, the engine is controlled for delivering a given referenced torque on the wheels. However, no variable is measured downstream the combustion chamber and that means that engine management is not based on final measurements, and emissions limits or performance are tracked by feed-forward controllers with closed loop control on intermediate variables.

Regarding the emissions control in LD diesel engines without SCR and LNT, the common layout is installing no physical nor virtual sensor for tracking pipe-out emissions; the engine is calibrated offline before going to market. During the calibration procedure, the emissions are measured by laboratory sensors with limited dynamic capabilities. Furthermore, standard procedures are often based on steady-state measurements and indeed, dynamic homologation cycles, such as the new European driving cycle (NEDC) (see Chapter 3), are not representative of real-life driving cycles [15]. Finally, hardware unit-to-unit dispersion (engine, sensors, etc.), system ageing and the effects of external variables (temperature, pressure, humidity, etc.) make this procedure does not ensure that the engine keeps working as when it was calibrated.

Narrow band lambda sensors for measuring lean or rich mixture have been installed in SI engines from 1968 (for controlling the TWC) and wide band lambda sensor for providing full resolution oxygen concentration can be found in some diesel engines from the end of the 90s. Anyway, the exhaust oxygen measurement in diesel engines is not standard and is not usually utilised for controlling the air path (except for detecting injector drifts).

Furthermore, it is less usual finding on-board sensors for measuring engine-out emissions in LD diesel engines. Different reasons can be stated:

- Exhaust gas concentration sensors (mainly  $\text{NO}_x$  and soot) with a sufficient small size and acceptable accuracy for on-board measurements have not been available until the last decade. More recently, on-engine  $\text{NO}_x$  can be measured with  $\text{NO}_x$  sensors based on the  $\text{ZrO}_2$  planar technology [16, 17]. On-board soot sensors are also starting to be marketed.

- The high cost of these on-board sensors with respect to the traditional ones used in automotive until EURO V standards (the cost can be several times higher according to [18]).
- Even though the steady-state accuracy of these sensors is acceptable, the dynamic performance is still problematic due to the slow and filtered responses in the order of 1 s, which prevent from using them with real-time purposes.

Anyway, these sensors are being gradually introduced in the market, enhanced by the development of after-treatment devices<sup>2</sup>. Admittedly, some luxury cars have already installed NO<sub>x</sub> sensors some years before. However, the problems derived of using the raw output signal in real-time functions must be addressed.

Models may be used as an alternative to sensors, since they can reproduce fast transients and present offline prediction capabilities [19,20]. Nevertheless, models suffer from drift depending on time and engine operating conditions, and uncertainties linked with the model structure or non-considered inputs. In commercial engine ECUs, the prevailing approach is to use look-up tables to model nonlinear and operating point dependencies because of the simple programming and easy comprehension. An intensive test campaign should be made for filling the tables with a big number of steady-states for offline calibration and a limited number of dynamic tests for validation. Anyhow, the drift problem cannot be eliminated unless sensors with acceptable accuracy are used.

Two main options can be considered for cancelling drift: modelling the bias as an additive or multiplicative factor to the model, or recalibrating the model. For both adaptive options, observers can be designed. For the former, the bias can be tracked by designing a proper drift correction observer, while for the latter, the model parameters can be updated for minimising the errors.

For all those techniques and cases, the Kalman filter (KF) [21] is a standard tool for observing variables. The multiple possibilities for designing observers make feasible to estimate drift, correct variables and indeed calibrate and update tables and models, all in an online basis required for diesel diagnostics or control. Anyway, computational limitations should be issued, and optimized algorithms and simplifications are crucial, jointly with the robustness of the algorithms.

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<sup>2</sup>In HD engines it is usual finding NO<sub>x</sub> sensors for a correct urea dosing in SCR systems.

### 1.3 Scope of the work

The dissertation covers the problem of the online estimation of diesel engine exhaust concentrations of  $\text{NO}_x$  and the fuel-to-air ratio or richness ( $\lambda^{-1}$ ). Two information sources are utilised:

- on-board sensors for measuring  $\text{NO}_x$  and  $\lambda^{-1}$ , and
- control oriented models (COM) in order to predict  $\text{NO}_x$  and  $\lambda^{-1}$ .

The evaluation of the static accuracy of these sensors is made by comparing the outputs with a gas analyser, while the dynamics are identified on-board by performing step-like transitions on  $\text{NO}_x$  and  $\lambda^{-1}$  after modifying ECU actuation variables. The former are produced by provoking sharp changes in the start of injection (SOI) while the latter by sharp changes in the injected fuel. With a sufficient data set, the response time and delay of the sensor are identified.

Afterwards, control oriented models for estimating  $\text{NO}_x$  and  $\lambda^{-1}$  are proposed. Regarding the  $\lambda^{-1}$  prediction, the computation is based on the relative fuel-to-air ratio, where fuel comes from an ECU model and air mass flow is measured by a sensor. For the case of  $\text{NO}_x$ , the model is set-point relative and based on look-up tables fitted for representing nominal engine emissions with an exponential correction based on the intake oxygen variation. Different correction factors for modelling other effects such as the thermal loading of the engine are proposed. The model is able to predict  $\text{NO}_x$  with a low error and a simple structure.

Regardless of using models or sensors, model drift and sensor dynamic deficiencies affect the final estimation. In order to solve this problem, data fusion strategies are proposed by combining the steady-state accuracy of the sensor and the fast estimation of the models. In a first approach, a drift correction model tracks the bias between the model and the sensor but keeping the fast response of the model. A KF is used for observing the bias, acting as a high-pass filter for the model and a low-pass filter for the sensor. In a second approach, the updating of look-up tables by using observers is coped with different versions based on the extended Kalman filter (EKF). Particularly, a simplified KF (SKF) allows to observe the parameters with a low computational effort.

Finally, the methods and algorithms developed in this work are combined and applied to the estimation of  $\text{NO}_x$  and  $\lambda^{-1}$  by drift correction algorithms and by developing adaptive look-up tables. The methods developed in this

work are applied to  $\text{NO}_x$  and  $\lambda^{-1}$ , however these could be employed for other variables such as soot, or engine parameters such as the volumetric efficiency.

## 1.4 Objectives

The present work provides solutions and ideas in the field of adaptive estimation and drift correction of signals and models. Particularly, the following are the main objectives of the present work:

- The introduction of a novel method for the online characterisation of  $\text{NO}_x$  sensors (Chapter 3).
- The development of a simple ECU-oriented model for diesel  $\text{NO}_x$  prediction (Chapter 4).
- The design of a simplified version of the EKF with similar accuracy as the standard KF, but that requires a much lower memory resources and calculation time, for updating look-up tables and models (Chapter 5).
- The development of a full methodology covering design, tuning and on-engine implementation of observers for the estimation of  $\text{NO}_x$  and  $\lambda^{-1}$  (Chapters 5 and 6).
- The design of a set of observers for updating an ECU-oriented model based on physical and virtual sensors and applied to the  $\text{NO}_x$  prediction (Chapter 6).

### 1.4.1 Methodology

In the following, the work methodology is sketched up and split up among the chapters.

Chapter 2 emphasises the need of information for a proper control of the diesel engine and reviews sensors and models for estimating gas concentrations in diesel engines, while also presents different methods for fast estimation of variables. These methods are focused on the drift correction and updating algorithms for look-up tables.

Chapter 3 describes the system set-up including the engine characteristics, which in the present work has been a 2.2 litre turbocharged DIC I engine, and the different engine tests performed on that engine. Furthermore, a novel procedure for the online characterisation of  $\text{NO}_x$  sensors based on SOI variations



is depicted. Characterisation of  $\lambda^{-1}$  output from  $\text{NO}_x$  sensors is also presented. This is a key point for the application of the algorithms developed in subsequent chapters, as the knowledge on the sensor response greatly affects the results.

Chapter 4 presents the design and validation results of ECU-oriented models for  $\lambda^{-1}$  and  $\text{NO}_x$  prediction. The first is just based on the calculation of the fuel-to-air ratio by the injected fuel mass flow and the air mass flow signals from the ECU while the second is based on a nominal set-point relative fitting of the  $\text{NO}_x$  with a series of corrections for accounting with variations on  $\lambda^{-1}$ , temperatures and other signals. The  $\text{NO}_x$  model combines look-up tables with physical-based equations and is designed for being implemented on commercial ECUs.

Chapter 5 presents two data fusion strategies for the fast estimation of variables when a steady-state accurate sensor and a fast model are presented. The first is based on drift correction algorithms in order to track the bias between a model and a sensor. In order to observe the bias, a steady-state KF is designed and the observer tuning is also discussed.

The second strategy comes from that look-up tables are often used in automotive control systems and there is a need for systematic methods that can estimate or update them on-line. According to that, learning algorithms for the online updating of look-up tables are proposed. Based on the EKF for observing look-up tables parameters [22], two different versions that keep some of the properties of the original KF but optimise the computation are designed.

The methods presented in the Chapter 5 can be applied to different engine variables. In this work, Chapter 6 applies them to the fast estimation of  $\lambda^{-1}$  and  $\text{NO}_x$ . In order to do that, different algorithms based on the dynamic estimation methods of Chapter 5 are proposed.

Afterwards, a full methodology for the fast estimation of  $\lambda^{-1}$  is presented. This part covers the methodology, the implementation of robustness conditions for coping with signal uncertainties and proposes an adaptive look-up table for modelling operating point dependency of the bias. Finally, experimental results are presented.

The second part of the chapter is centred on the fast estimation of  $\text{NO}_x$  by designing different algorithms. First, the evaluation of the methods for updating look-up tables designed in the Chapter 5 is validated by using real engine data for the  $\text{NO}_x$  estimation. Furthermore, single 2D adaptive look-up tables scheduled by injection and speed are proposed for modelling slow

varying cycles. Second, different set of observers are proposed in order to update parameters on the NO<sub>x</sub> model developed in the Chapter 4 and for minimising the errors in the model inputs when sensor signals are available. To conclude, the estimation of the actual NO<sub>x</sub>, solving the problems of filtering and delay of sensors, is proposed.

Finally, Chapter 7 presents the conclusions and the future works derived from this dissertation.

## 1.A Publications

The publications by the author that are directly linked with the dissertation are presented next, where the Chapters related with the papers are shown in brackets.

### Journal papers

- Galindo J, Serrano JR, Guardiola C, Blanco-Rodriguez D and Cuadrado IG. An on-Engine Method for Dynamic Characterisation of NO<sub>x</sub> Concentration Sensors. *Experimental Thermal and Fluid Science*, Vol. 35 n° 3: 470-476, 2011. [Chapter 3]
- Guardiola C, Pla B, Blanco-Rodriguez D and Cabrera P. A Learning Algorithm Concept for Updating Look-up Tables for Automotive Applications. *Mathematical and Computer Modelling*, Vol. 57 n° 7-8: 1979-1989, 2013. [Chapter 5]
- Guardiola C, Pla B, Blanco-Rodriguez D, Mazer A and Hayat O. A Bias Correction Method for Fast Fuel-to-Air Ratio Estimation in Diesel Engines. *Proceedings of the Institution of Mechanical Engineers, Part D: Journal of Automobile Engineering*. Vol. 227 n° 8: 1099-1111, 2013. [Chapters 3, 4, 5 and 6]
- Guardiola C, Pla B, Blanco-Rodriguez D and Eriksson L. A Computationally Efficient Kalman Filter based Estimator for Updating Look-up Tables Applied to NO<sub>x</sub> Estimation in Diesel Engines. *Control Engineering Practice*. Vol. 21 n° 11: 1455-1468, 2013. [Chapters 5 and 6]
- Guardiola C, Pla B, Blanco-Rodriguez D and Calendini PO. ECU Oriented Models for NO<sub>x</sub> Prediction. Part 1: A Mean Value Engine Model for NO<sub>x</sub> Prediction. Submitted to *Proceedings of the Institution of Mechanical Engineers, Part D: Journal of Automobile Engineering*. [Chapters 4]
- Guardiola C, Climent H, Pla B and Blanco-Rodriguez D. ECU Oriented Models for NO<sub>x</sub> Prediction. Part 2: Adaptive Estimation by Using a NO<sub>x</sub> Sensor. Submitted to *Proceedings of the Institution of Mechanical Engineers, Part D: Journal of Automobile Engineering*. [Chapters 5 and 6]

### Conference papers

- Desantes JM, Luján JM, Guardiola C and Blanco-Rodriguez D. Development of NO<sub>x</sub> Fast Estimate Using NO<sub>x</sub> Sensors. *EAEC 2011 Congress*, 2011. [Chapter 6]

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# Chapter 2

## Exhaust gas concentrations estimation in diesel engines

### Contents

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### 2.1 Introduction

Driven by the advancements in diesel engines and especially due to the stringent regulations in the recent years that have enhanced the introduction of the after-treatment (AT) systems, the control and diagnostic logics in the modern ECU are accomplishing a renewal. The typical configuration of a short route EGR and a variable geometry turbine (VGT) shall be extended to include dual EGR loops, and different systems for eliminating particles,

hydrocarbons and  $\text{NO}_x$  in the exhaust. This modernisation must be run jointly with the improvements in the ways of retrieving information from the states of the engine, particularly from engine-out concentrations. This chapter gives an overview on the actual state on technologies related with diesel engine and the different sources of information from the exhaust gas concentration on diesel engines: sensors, models and adaptive estimators. A specific discussion about the  $\lambda$  and  $\text{NO}_x$  estimation in CI engines is made due to their relevance in the present dissertation.

## 2.2 Diesel engine subsystems

There exists different possible diesel engine layouts, but they all share at some point the following subsystems [1]: the fuel path, the air path, the after-treatment and the control system.

### 2.2.1 The fuel path system

The common-rail (CR) system and the direct injection (DI) is standard in current diesel engines [2]. The main advantages of the CR against other systems, such as the distributor pump and the unit pump injector, are the multiple injections and the control flexibility. A pressurised deposit monitored by a rail pressure sensor is capable of maintaining the pressure highly constant during the injection.

In the ECU, the total injected fuel mass ( $m_f$ ) is modelled by a look-up table function of the injection duration ( $t_{id}$ ) and the rail pressure ( $p_{rail}$ ). The SOI actuation ( $u_{soi}$ ) is determined by a calibrated look-up table as function of speed ( $n$ ) and the desired  $m_f$  and is measured in degrees with respect to the top dead centre (TDC). In the case of sharp load steps and in order to avoid the excessive emissions of HC and soot, the smoke limiter function limits  $m_f$  until the air path responds and supplies the air mass flow ( $\dot{m}_a$ ) required for the combustion, as depicted later in Figure 2.2.

$$m_f = f(t_{id}, p_{rail}) \quad (2.1a)$$

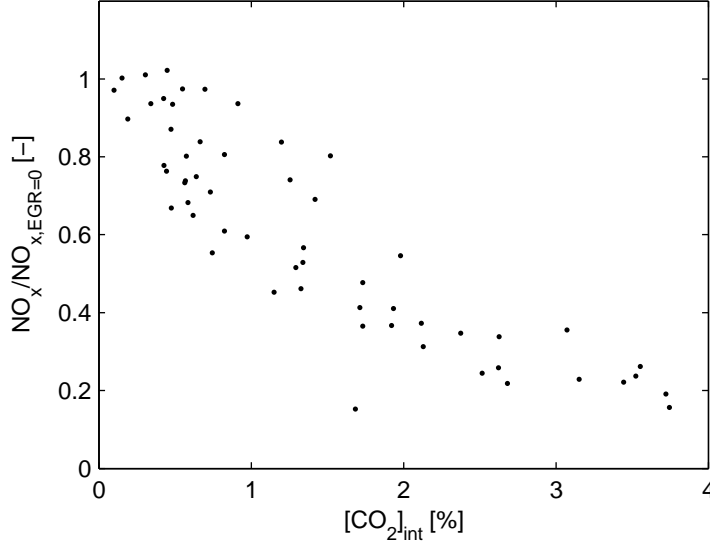
$$u_{soi} = f(n, m_f) \quad (2.1b)$$

### 2.2.2 The air path system

In a commercial CI engine, the following subsystems can usually be found in the air path: the air filter (in order to clean the entering air, although causing a minimum pressure drop in the line), a (single or double stage) turbocharger (TC) including an intercooler for cooling the compressed air (thus increasing the density), the intake and exhaust manifolds, the EGR loops (high and low pressure loops) and the pipe-out line (the after-treatment devices are considered as an independent subsystem because of the different possibilities existing today, even though they are really connected to the pipe-out line).

A single TC is the common layout in commercial engines, but some mount also dual stage TC [3]. The main function is increasing the intake air mass flow by increasing the density with a compressor usually powered by a turbine shaft located at the engine exhaust [4]. One of the main drawbacks of the system is the turbolag of the TC which is associated with the inertial response of the turbo shaft to accelerations, especially at low speed and low loads, when the exhaust energy is quite low [5]. In order to get the maximum efficiency of the compressor and turbine, two types of control are usual: waste-gate (WG) valves and a variable geometry turbine (VGT), also known as variable nozzle turbine (VNT). The WG configuration is based on the use of a fixed geometry in the turbine with a valve that acts bypassing the flow minimising the effective exhaust flow through the turbine. This solution has practically been replaced by the use of the VGT, which is based on movable nozzles that vary the turbine work depending on the operating point. The VGT allows to increment the engine performance reducing the specific fuel consumption and also adds a major flexibility to the control, admittedly that increments the related complexity and calibration effort. In a dual stage TC, there exists a wide variety of possibilities of combining WG, VGT and fixed geometries in the high pressure (HP) and low pressure (LP) turbines [6].

The exhaust gas recirculation (EGR) is the most widely extended  $\text{NO}_x$  reduction system [7] and is used in CI engines since the middle of the 90s. The basic idea is recirculating a portion of the exhaust gas to the intake. The principal effect of the EGR system is reducing the effective  $\text{O}_2$  (increasing effective  $[\text{CO}_2]_{int}$ ) at the intake and then diminishing the peak temperature in the cylinder and thus reducing  $\text{NO}_x$  as it is shown in Figure 2.1.



**Figure 2.1.** Nominal  $NO_x$  emissions depending on  $[CO_2]_{int}$  for nominal tests in the CI engine explained in Chapter 3. Values are normalised with the maximum  $NO_x$  value for each operating point, which corresponds to  $u_{egr} = 100$  (fully closed) for each pair of  $[n, m_f]$ .

The EGR rate in steady-state operation can be defined as the quotient between the total EGR mass flow  $\dot{m}_{egr}$  and the intake mass flow ( $\dot{m}_{int}$ )

$$\dot{m}_{int} = \dot{m}_a + \dot{m}_{egr} \quad (2.2a)$$

$$EGR = \frac{\dot{m}_{egr}}{\dot{m}_{int}} \quad (2.2b)$$

The work by Ladommatos *et al.* [8–12] is a good reference for understanding the effects of the EGR in the combustion and emissions. In addition to the internal EGR (inert gas fraction that stays in the cylinder after the combustion), the external EGR can comprise high pressure EGR (HPEGR) and low pressure EGR (LPEGR) loops. The HPEGR is based on the extraction of a portion of the exhaust gas upstream of the turbine and driven normally to the intercooler output (hot EGR). The LPEGR extracts a portion of the exhaust gas at some point downstream of the turbine (it is usual to locate it downstream of the diesel particulate filter for avoiding damage in the compressor) and is guided at the compressor inlet (after the air filter). Besides the installation of valves for controlling the flow, coolers are installed for increasing the flow density.



The HPEGR, namely EGR for simplification, is the standard in CI engines due to the simple layout and control: the pressure differences between exhaust and intake facilitate that gas goes to the intake with some exceptions, which can occur for high EGR rates, problem that can be solved e.g. by introducing a throttle valve [13]. Even though the LPEGR did not enter into massive production, there exists today a certain interest in combining both loops for further reduction of emissions that can push up to relax the AT efficiencies [14, 15] and thus the cost of complex systems.

**EGR/VGT control.** The air path is managed independently of the fuel injection system. The effects of the injection settings are considered in the air path as instantaneous inputs for a multiple input multiple output (MIMO) control problem. More concretely, the fuel injection quantity ( $m_f$ ) and the engine speed ( $n$ ) are used as scheduling variables for defining the set-points for the EGR and VGT actuation. Then, for a common layout of HPEGR and VGT (or WG), the inputs are the EGR valve position ( $u_{egr}$ ) and the turbine actuation ( $u_{vgt}$  or  $u_{wg}$ ), besides other actuators such as valves for bypassing the flow over the EGR cooler ( $u_{bp,egr}$ ) if necessary. If the TC is dual stage, it is usual that only one of the turbines can be actuated. For other layouts, different control strategies can be designed, e.g. with a twin turbo sequential parallel configuration, an extra variable is necessary in order to coordinate the switching [6]. Varnier [3] gives a complete review on boosting technologies and its associated control.

The EGR/VGT control strategy basically consists on two maps for determining the set-points for the air mass flow ( $\dot{m}_a$ ) and the boost pressure ( $p_{boost}$ ), while  $u_{egr}$  and  $u_{vgt}$  are commanded for reaching these set-points [16]. The coupling of the EGR/VGT system is avoided by defining control regions, with and without EGR. When engine is working in the EGR area,  $u_{egr}$  is controlled in closed loop for reaching the required air mass flow set-point scheduled by  $m_f$  and  $n$ . The VGT is commanded in open loop by a feed-forward controller that imposes a given  $u_{vgt}$  depending also on  $m_f$  and  $n$ . This structure avoids the problem of the EGR/VGT coupling that can be exemplified as follows: when the EGR valve is opened, a closing on turbine nozzles increments the exhaust backpressure and the turbine regime, this effect produces a higher recirculation of EGR reducing the effective air mass flow at the intake (the EGR mass flow replaces the fresh air). However, when the EGR valve is fully closed, a closing of the turbine nozzles produces an increase of the intake air mass flow. This occurs for medium-high engine loads and  $u_{vgt}$  is controlled in closed loop, while the EGR valve keeps closed. The control objective is then reaching the

intake air mass flow set-point. Figure 2.2 schematises the air path and the smoke limiter controller.

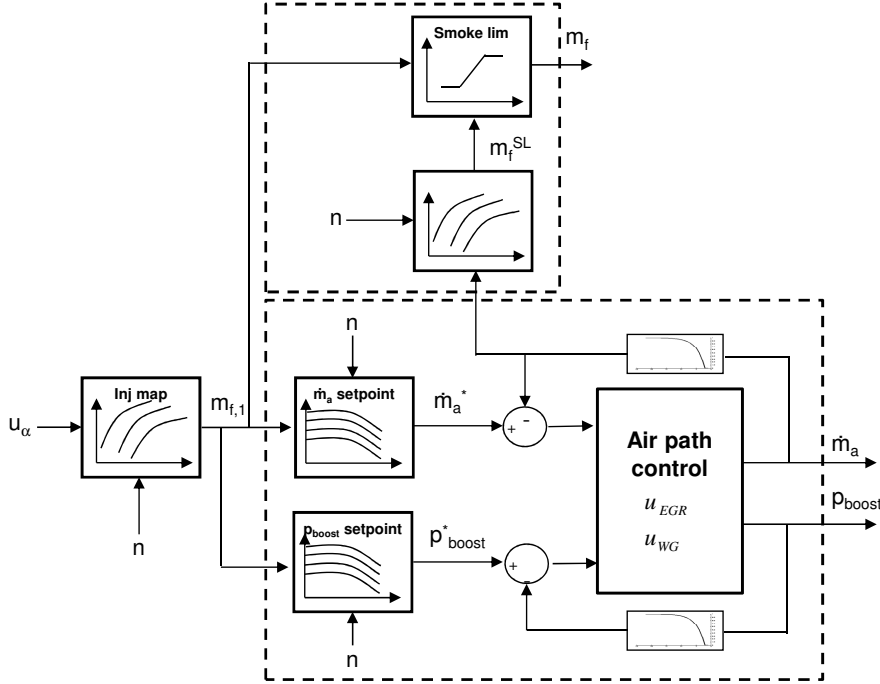


Figure 2.2. Smoke limiter and air path control blocks.

Furthermore, some corrections or maps switching are applied depending on combustion modes or for cold starting strategies (the references for injection and air path are modified). The engine coolant temperature ( $T_{cool}$ ) is also used for corrections and for defining the cold starting strategies.

An alternative to the commercial air path control is using virtual and physical sensors or including adaptive strategies by feeding back the controllers with direct measurements of exhaust conditions, e.g. exhaust gas concentrations whose estimation is further discussed in the Section 2.3. Then, model predictive control (MPC) [17, 18] or other strategies [19] can be exploited for such control.

**Dual EGR/VGT control.** When using a dual EGR loop, in addition to the control variables presented for the single loop, the following inputs are necessary: LPEGR valve ( $u_{lp}$ ), intake throttle valve ( $u_{it}$ ) and/or the back-pressure valve positions ( $u_{bp}$ ). The total EGR rate is calculated as in (2.2)

but considering mass flows of the low and high pressure loops respectively  $\dot{m}_{lp}$  and  $\dot{m}_{hp}$

$$\dot{m}_{egr} = \dot{m}_{lp} + \dot{m}_{hp} \quad (2.3)$$

Depending on the engine conditions and control needs, the EGR split ( $\dot{m}_{lp}/\dot{m}_{hp}$ ) is selected, i.e. the set-points for  $\dot{m}_{lp}$  and  $\dot{m}_{hp}$ . The control is similar to that of the single EGR/VGT control but designing an appropriate algorithm for defining the EGR split [20]. A common solution in steady-state is controlling  $u_{lp}$  and if necessary opening  $u_{hp}$  for reaching the  $\dot{m}_a$  set-point (EGR is not directly measured). Indeed, the LP loop itself might be sufficient for getting the required  $\dot{m}_a$  set-point. For transient operation, the discussion must evaluate the characteristic response times of the systems: the HPEGR line is faster than the LPEGR line due to the higher pressure differences and the shorter length. This produces that different trade-offs between  $u_{hp}$  and  $u_{lp}$  may be stated during transients. An intrinsic benefit of using a dual loop EGR is that for the EGR engine area, the total  $\dot{m}_{egr}$  quantity may be increased leading to further reductions on  $\text{NO}_x$ , which can alleviate the needs of AT or could move the  $\text{deNO}_x$ <sup>1</sup> trade-off to select a lean  $\text{NO}_x$  trap (LNT) rather than a selective catalyst reduction (SCR) system. Model based and optimal approaches to the control problem of dual EGR loops can be found in [15, 21, 22].

**Control of the air and fuel paths.** Some authors have proposed a joint control of the air and fuel paths for optimising the engine transient operation. For that, fast acting variables linked to the fuel path ( $u_{soi}$  or  $p_{rail}$ ) are able to track  $\text{NO}_x$  or soot while  $u_{egr}$  and  $u_{vgt}$  are arranged to improve the engine transient response while keeping the engine operation (ensuring the torque on wheels) without exceeding emission limits. This can be made by feeding back the controllers with real time measurements of exhaust gas concentrations (typically  $\text{NO}_x$  and soot, or exhaust oxygen) [17, 23]. Table 2.1 sums up the characteristic response times for step variations in the inputs, showing how the associated fuel path responses are quite faster than the air path ones. Furthermore, the response times of the TC are in general higher than those related to the EGR loop due to the turbolag problem.

Some authors have proposed a transient operation optimization of the engine management strategy (EMS), see e.g. [24]. However, manufacturers are still reluctant to change the actual implementation in the fuel and air path

<sup>1</sup>The term  $\text{deNO}_x$  refers in general to an AT system for reducing  $\text{NO}_x$ , see Section 2.2.3.

**Table 2.1.** Approximate characteristic response times for the main actuation variables in a CI engine.

Command	Response time
$u_{p_{rail}}, u_{soi}, u_{id}$	100-600 $\mu$ s
$u_{hp}$ ( $u_{egr}$ )	0.2-0.5 s
$u_{lp}$	0.2-1 s
$u_{vgt}, u_{wg}$	0.5-1.5 s

system: the EMS is confined to fixed calibrations and optimised for steady-state operation with a validation in transient operation without feedback on exhaust engine states.

### 2.2.3 The after-treatment systems

The new stringent regulations have forced to get an impressive development in diesel engines, and particularly with the after-treatment devices (AT); Johnson [25] and Twigg [26] present complete reviews on emissions control and available AT technologies. The standard configuration for AT in diesel engines for EURO VI will comprise

- a diesel oxidation catalyst (DOC) for three main functions: oxidising CO and HC, heating the exhaust gas and oxidising NO to NO<sub>2</sub>, which is useful in the DPF regeneration and also facilitates the SCR performance;
- a diesel particulate filter (DPF) for eliminating soot; and
- a deNO<sub>x</sub> system for reducing NO<sub>x</sub> that it is usually a selective catalyst reduction (SCR) or a lean NO<sub>x</sub> trap (LNT).

The DPF and LNT action is based on substrate materials that store the pollutants through mechanisms or chemical principles. When the system is full, a regeneration by means of post-injections and thus increasing exhaust heat is necessary. The DOC system is built from ceramic and catalytic noble metals, such as platinum or rhodium, for enhancing the oxidation of CO and HC as well as NO to NO<sub>2</sub>. SCR systems eliminate NO<sub>x</sub> by dosing urea and water (principle component is ammonia) and reacting with NO<sub>x</sub> for obtaining N<sub>2</sub> and water, where a dedicated control management is necessary for optimising its dosing [27, 28]. The SCR requires an specific installation for the urea storage that increases the cost and weight, and for this reason, the SCR

is more extended in HD than in LD. Major concerns in SCR are the light-off related with the minimum temperature for guarantying the efficiency of the system (critical when the engine is cold or it is working in low-load operation); and the ammonia slip, due to the excessive injection of urea, which does not react with  $\text{NO}_x$ , constituting an economical cost and affecting the output of the  $\text{NO}_x$  sensors based on  $\text{ZrO}_2$  (see Section 2.3.1.2).

DPF and DOC are standard in EURO V cars, and  $\text{deNO}_x$  systems will also be with the application of EURO VI. Anyway, the trade-off between SCR and LNT is not solved [29], and different factors influence the final decision. Thinking of future regulations, manufacturers struggle between getting lower removal efficiencies (around 85%) by using LNT or allowing the engines to work on much higher efficiencies and then higher  $\text{NO}_x$  production but with the use of a SCR with a removal efficiency nearly 98% [25].

The selection of the layout of the ATs in the engine is another important trade-off and influence the correct operation of the systems. Some possibilities are later shown in Figure 5.1; however this is a trending discussion. It seems logical that the DOC should be located first because it is required for a proper performance of DPF and SCR, but the trade-off between DPF and SCR or LNT present advantages and drawbacks. Indeed, there are some research on moving the DPF upstream of the turbine for preventing turbine and compressor damage (if it is located before the EGR loop) [30] or doing something similar with the DOC [31]. There are also some alternatives that combine some of the systems for reducing the packaging and final weight [32].

With respect to the feedback, the knowledge of the concentrations upstream and downstream of the AT systems allows to diagnose the performance, and is required for direct control of the final emissions. The temperature also influences the regeneration of the DPF and LNT as well as it is critical for the SCR light-off (key reactions are inhibited at low temperatures). The pressure drop in the line could also be used for detecting failures or when the DPF or LNT are full. The decision on the sensor set for a proper control and diagnosing is a challenging problem that requires a specific study out of the scope of this work. The use of sensors and models for estimating exhaust gas concentrations, especially  $\text{NO}_x$  and exhaust oxygen, is discussed in section 2.3.

#### 2.2.4 The control system

As presented in previous subsections, the modern diesel engine has evolved until reaching today an important and increasing complexity, which is reflected with the introduction of different subsystems for improving the performance

and minimising the emissions. The vehicle electronics in the 80s were reduced to the radio plus simple engine controllers. At the same time, these were based on mechanical and hydraulic devices with a simple electronic assistant, and dedicated to specific functions such as the ignition in the case of SI engines. However, the evolution of the engines has been closely related with the rise of complexity of the modern embedded electronics in automotive, as it is well depicted in [33]. Nowadays, safety performance, fuel efficiency and comfort are major demands and these are achieved by installing proper control units. Modern cars have between 20 and 70 control units with a size of  $10^8$  in object instructions, while in the first 90s this was in the order of  $10^6$ , according to Ebert and Jones [33].

From all the control units in the vehicle, the ECU is in charge of managing the engine. Others work with the braking systems, the air conditioning in the cabin or the gear shifting management [1]. In the beginning of the 00s, the ECU EDC16 model by Bosch for diesel engines changed the view of the diesel control by using the final torque demand on the wheels as the feedback for a proper air and fuel path management. This ECU also included a fully flexible control of the common-rail (pilot, main and post-injections), the VGT and the EGR, all in the same processor. The current EDC17 [34] model coordinates a bigger number of data from sensors, actuators and models in order to get a full management of the diesel engine.

From all the requirements and purposes of the ECU, the on-board diagnostics (OBD) standards were applied to engines in order to implement safety routines for detecting sensors or engine systems malfunctions [35]. Currently, about 40% of the automotive electronics are devoted to OBD functions. The gradual introduction of the AT systems in diesel engines, especially with the current trade-off between selecting SCR (more complex due to the urea dosing but effective) or LNT (more simple but less effective), is requiring a renewal of the current diesel air path control [29], and new functions and managing strategies are being developed; see for instance Yang *et al.* [36] who propose an independent control unit for the AT line separated from the main ECU.

#### **2.2.4.1 Summary on the information required for a global diesel management**

The commercial diesel engine management focuses on references measured upstream of the combustion chamber, with no direct feedback from the exhaust line. This makes that emissions are controlled in open loop and possible changes and the ageing are not taken into account. The sensors set used in

the automotive in the beginning of 00s [37] had hardly changed until the end of the last decade. A major factor is the unnecessary of increasing the number and the type of sensors due to the current control logic. Furthermore, other important factor is the low cost of pressure, temperature and rotational motion sensors (bellow 3\$ per unit) against the higher cost (minimum >10-20\$) of more sophisticated sensors, such as NO<sub>x</sub> or in-cylinder pressure ones (references obtained from [37] and conversations with engine manufacturers).

Table 2.2 summarises the usual inputs and sensors that are installed in current diesel engines, while indicating possible alternatives in a short future [38], which are essentially driven by the emission standards (the need of ATs control and diagnosis). It is rather complicated establishing an universal sensor set and some discrepancies can be found, e.g. GM is starting to include  $p_{cyl}$  sensors in engines and some AUDI and BMW vehicles already install NO<sub>x</sub> sensors before and/or after the SCR system.

**Table 2.2.** Summary of the on-board sensors and control inputs used in a commercial LD diesel engine for control and feedback, and split between the main engine subsystems. The inputs column defines the main actuator signals in the diesel engine. The remaining columns show the variables that are measured by sensors, differentiating between standard for EURO V cars and the possible expectations in a short future. Admittedly, some discrepancies can exist between this distinction depending on the considered car and sector (luxury, utility, etc.). The symbol \* represents sensors that can be found in some actual engine models but with less probability. The injection settings are defined by the rail pressure control  $u_{prail}$ , the injection timing command ( $u_{soi}$ ) and the injection duration  $u_{id}$ , and these settings can be specified for pilot injections main injections and possible post injections (variable  $u_{t_{mi}}$  makes reference to the duration of the main injection only). A detailed summary on the symbols and acronyms used in this work appears in the beginning of this document. The interested reader is also referred to Turner [38] for finding basic comprehensive explanations about the principles of measurement of the main automotive sensors.

System	Inputs	Standard sensors	Expected sensors
Injection	$u_{prail}$ $[u_{soi}, u_{id}]_{pilot, main, post}$	$n, p_{rail}$	$p_{cyl}^*$ fuel quality
Air path	$u_{vgt}, u_{hp}, u_{lp}$ $u_{it}, u_{bp}$	$p_{boost}, \dot{m}_a, T_{cool}$ $T_{boost}, p_{amb}, T_{amb}$	$\dot{m}_{egr}, T_{egr}$ $n_{vgt}, \lambda_{int}^*$
AT	$[u_{soi}, u_{id}]_{post}, u_{urea}$		NO <sub>x</sub> <sup>*</sup> , soot, $\lambda_{exh}^*$

Alternatively, model based strategies present different advantages: models can be used for engine calibration saving costs [39]; models can easily be

linked in control and diagnosis algorithms as they present faster responses than sensors; and the fusion of sensors and models permits to estimate variables or update parameters on real time solving the drift and ageing and improving the dynamic responses of the original signals.

This section has emphasised the importance of the information about engine states and how to correctly retrieve them on-board. Next section deals with the topic of exhaust gas concentration estimation in CI engines, stressing the variables  $\text{NO}_x$  and  $\lambda$ .

## 2.3 Dynamic exhaust gas concentration estimation

Previous section has made a review on the different technologies in order to manage the diesel engine and has underlined the need of reliable methods for estimating variables online. Test-bed measurement systems are usual for testing and research but these present in general deficient dynamic responses and overall limited on-board capabilities. However, in the last years, and forced by the stringent emission laws, developments of on-board gas concentration sensors have proliferated and some of them are a reality today, such as  $\text{NO}_x$  or wide-band lambda sensors. Furthermore, COMs and data fusion (DF) techniques that combine different information sources can also be used for real-time (RT) estimations.

Current section gives a review on the different methods for the estimation of the exhaust gas concentration in CI engines, emphasising the on-board methods. Therefore, in addition to being feasible from the point of view of cost and engine implementation, the accuracy and dynamic responses of these estimators should be issued. Hereinafter the following three main possibilities are discussed:

- sensors,
- models and/or virtual sensors, and
- DF techniques for combining different information sources.

### 2.3.1 Sensors

A subsequent division is made between test bed measurements systems and on-board gas concentration sensors, such as exhaust gas oxygen (EGO), universal exhaust gas oxygen (UEGO) sensors, fast soot sensors and finally



innovative solution for monitoring e.g.  $\text{NO}_x$  by using the substrate materials of the AT systems.

### 2.3.1.1 Test bed sensors

Some of the most common test bed sensors are the following: gas analysers [40] for different gas concentrations; opacimeters based on Beer-Lambert law [41] and smoke meters for soot; and gravimetric devices or spectrometers for PM [42], among others. But these are expensive, require from specific installations and their dynamic responses are questionable. Nevertheless, there have been some developments for fast response emissions analysers for  $\text{NO}_x$ ,  $\text{CO}_2$  or particles and portable emission measurement systems (PEMS) [43] are being generalised for on-board inspections but with limited dynamic accuracy [44] and packaging. Finally, standard procedures for measuring emissions in test-cells can show representative errors in the dynamic cycles with respect to the actual transient emissions [42,45] and this should be taken into account when processing the data [46–49] (see Section 2.3.2.3).

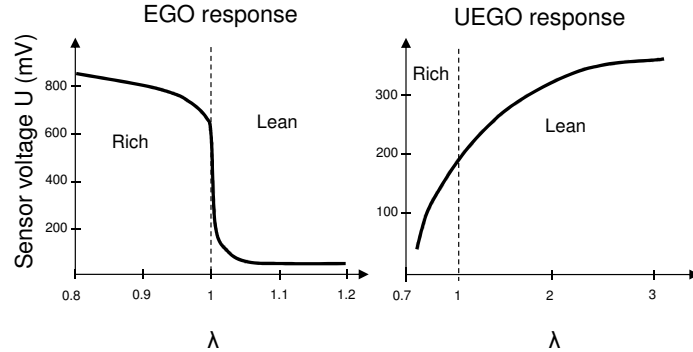
### 2.3.1.2 On-board gas concentration sensors

**Oxygen sensors.** In 1968, the introduction of EGO sensors [50] was the key for the implantation of the primary three way catalyst (TWC) control in SI engines. These sensors offered a binary lean/rich resolution for measuring oxygen concentration and thus the equivalent air-to-fuel ratio or lambda, which represents the excess air factor with respect to the stoichiometric air-to-fuel ratio (approximately 14.5 in diesel engines)

$$\lambda = \frac{\dot{m}_a}{\dot{m}_f} \frac{1}{14.5} \quad (2.4)$$

In the end of the 90s, planar UEGO sensors were implemented for SI and CI engines. It is common to use directly the term lambda probe for referring to these sensors. The term lambda is due to the characteristic shape of the signal output in narrow-band lambda sensors, which is similar to the  $\lambda$  letter from the Greek alphabet (see Figure 2.3). Furthermore, these are also known as switching-type sensors. UEGO sensors present full and linear resolution on oxygen over a wide range, hence the term wide-band. UEGO and EGO response characteristics are shown in Figure 2.3.

$\lambda$  measurement in CI engines is used for correcting injector drift. In the case of low temperature combustion (LTC) processes [51], intake charge

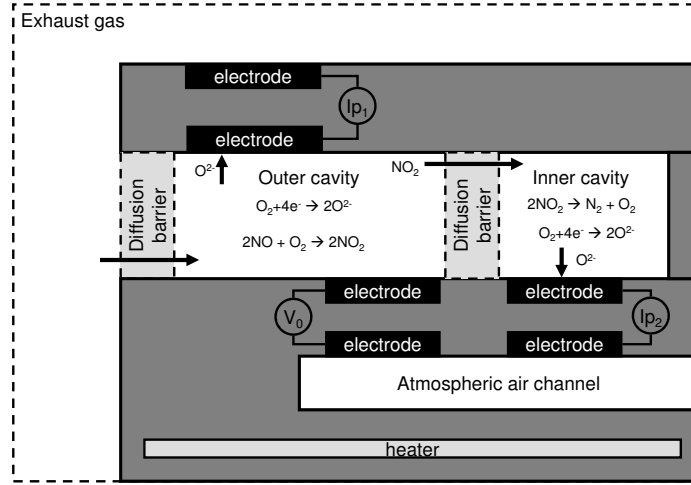


**Figure 2.3.** Characteristic outputs of oxygen lambda sensors. Left: Narrow-band lambda or EGO sensors. Right: Wide-band lambda or UEGO sensors.

composition, and hence  $\lambda$  is important for the combustion stability and control [52, 53]. Some studies about wide-band lambda sensors performance can be found in [54, 55].

**NO<sub>x</sub>-ZrO<sub>2</sub> sensors.** The development of on-board NO<sub>x</sub> sensors has been driven by those which are based on ZrO<sub>2</sub> layers [56, 57] and able to measure gas at wet condition (i.e. without removing exhaust gas water steam content as gas analysers do). These sensors have suffered an evolution over the last 15 years [58] and now are manufactured using the planar ZrO<sub>2</sub> multilayer technology [59], which combines thick film screen printing and ceramic tape casting [60]. Modern versions offer reduced warmup time, smaller size, lower weight and cost-effective production, which encourage their implementation on commercial engines. This kind of sensors simultaneously provides a measurement of the relative air-to-fuel ratio ( $\lambda$ ) and NO<sub>x</sub> concentration and must play a major role in the SCR control and diagnosis [61–64].

The working principle is well explained in Riegel *et al.* [60] and layout is schematised in Figure 2.4. This sensor presents two cavities with membranes for measuring oxygen and NO<sub>x</sub>, respectively. In a first outer cavity, an electrochemical pump adjusts the oxygen concentration from the diffusing gas to a predefined value (pump intensity  $I_{p1}$ ) by reducing O<sub>2</sub> to O<sup>2-</sup> with an electrode (usually Pt), thus providing a linear amperometric measurement of  $\lambda$ . At the same time, a faster binary output is provided, differentiating between rich and lean conditions in the exhaust gases (similar to EGO sensor output). The measurement principle of this first cavity is similar to that of a wide-band UEGO sensor. In addition, NO is oxidised to NO<sub>2</sub>, and then the NO<sub>2</sub> is diffused to a second inner cavity, where the oxygen produced through the



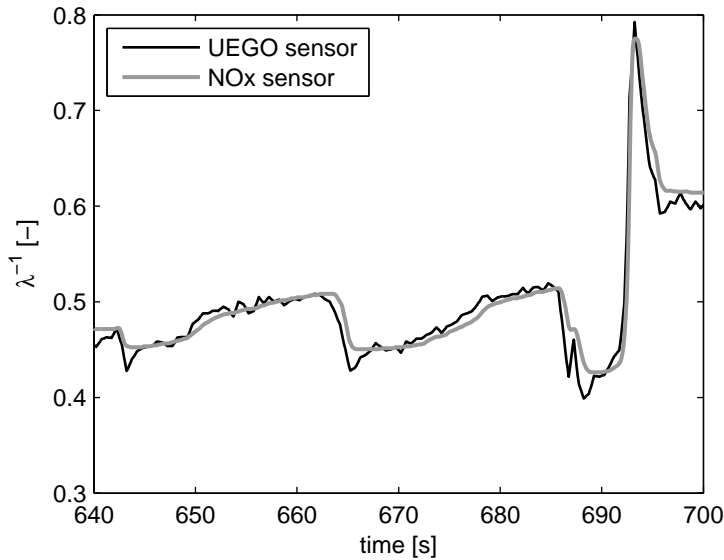
**Figure 2.4.** Schematic representation of a  $\text{NO}_x$ - $\text{ZrO}_2$  sensor, where the main reactions in the chambers and diffusion barriers are shown.

dissociation of the  $\text{NO}_2$  is pumped out in a similar way by means of a second electrochemical pump ( $I_{p2}$ ). The output of this pump is proportional to the  $\text{NO}_x$  concentration in the exhaust gases. A heater is necessary for keeping the sensor at high temperatures, where the device gets good resolution, i.e. the sensor needs some lag for starting to measure when switching on (the same occurs for UEGO sensors). Therefore, these sensors provide three signal outputs: a fast binary  $\lambda$  (around 100 ms), lean or rich, suitable e.g. for TWC diagnosis, a slower (around 500 ms of response time) full resolution  $\lambda$  and  $\text{NO}_x$  (with a response around 750 ms).

**$\text{NO}_x$  and UEGO sensor limitations.** Because of the importance of  $\text{NO}_x$  and UEGO sensors in the present work, response limitations of these sensors are underlined. Several studies have evaluated the accuracy and response times of real time  $\text{NO}_x$  sensor measurements for different applications, such as [65, 66].

Conventional UEGO sensors exhibit fast response times around 70 ms with sufficient accuracy [67]. This can be improved, as proposed in [68], moving the  $\text{O}_2$  sensor upstream of the turbine and using a Kalman filter [69] for taking into account pressure effect on the output signal. If the  $\text{NO}_x$  sensor is placed downstream of the AT systems, its response is affected by a considerable transport delay and filtering. Figure 2.5 compares  $\lambda^{-1}$  signals from a UEGO sensor located upstream of the AT systems, and from a  $\text{NO}_x$  sensor located

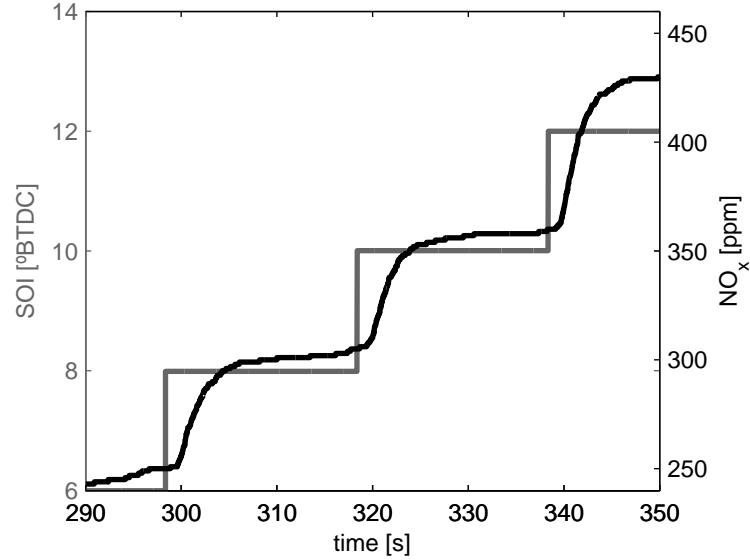
downstream of an AT line including DOC, DPF and SCR in a turbocharged diesel engine. The  $\text{NO}_x$  sensor signal is significantly slower and more filtered than that of a UEGO sensor, although the steady-state accuracy seems to be sufficiently accurate in both devices at a glance. Due to the measurement principles of both UEGO and  $\text{NO}_x$  sensors, the pressure in the runners also compromises the steady-state accuracy of the sensor.



**Figure 2.5.**  $\lambda^{-1}$  measurement from a UEGO sensor located at the turbine outlet and from a  $\text{NO}_x$  sensor downstream the after-treatment systems.

Concerning  $\text{NO}_x$  output from  $\text{NO}_x$  sensors, the delay is variable, depends on operating conditions, and is in the order of 1 s. This delay is caused for both transport and physical motives, which are difficult to model and can vary with the engine operating conditions. This issue makes difficult to use the raw sensor signals for RT critical functions. For illustrating sensor response, see Figure 2.6: the actual  $\text{NO}_x$  is expected to respond instantaneously when performing start of injection (SOI) steps where the delay and response time in the sensor are attributed to the sensor. Several studies have evaluated the accuracy and response times of RT  $\text{NO}_x$  sensor measurements [44, 65, 66].

Finally, since sensors are subjected to unit-to-unit manufacturing discrepancies, and can be affected by significant drift during their lifespan, methods for the online characterisation of sensors would improve the application of these for automotive control purposes. Usual calibration methods consist of



**Figure 2.6.**  $NO_x$  sensor response to SOI steps (black-right axis).  $NO_x$  presents a clear delay and filtering with respect to the SOI signal (gray-left axis). The dynamic response could be fitted with a first order delayed discrete filter. SOI units are crank angle degrees before the top dead centre ( $^{\circ}$ BTDC), while  $NO_x$  is measured in ppm. The delay is in the order of 1s while the response time is around 0.75s.

a specific test rig which generates known composition synthesis gas which is used for the static and dynamic calibration of the sensor, while high speed valves are needed for dynamic calibration [70]; or some solutions based on openly mounted sensors with valves allocated before the sensor, where gas composition is changed within milliseconds [71]. However, these calibration methods are restricted to laboratory use and cannot be performed during the operation phase on the engine. Chapter 3 presents a contribution for the on-line characterisation of  $NO_x$  output by means of SOI steps; furthermore,  $\lambda^{-1}$  output from wide-band lambda sensors may be characterised by means of load steps.

**Fast soot sensors.** According to Kasper [72], particles with lower diameters affect more to alveolar deposition after healthy tests with persons, which underline the risk of underestimating the effect of the residual soot downstream of the DPF. This carries out a sensitive problem for sensors: in addition to robustness, packaging and dynamic limitations, on-board soot sensors must measure with high resolution at low-soot values (this sensitive problem is also

linked with NO<sub>x</sub> sensors). Another problem with soot sensors is the selection of an appropriate metric as different metrics can be used depending on the physical measuring principle: particle number, total mass, sizes or light absorbing properties (opacity). A common solution for fast soot sensors, described in [73], is measuring the electrical conductivity of the carbonaceous fraction of the PM, which acts as a resistance between two electrodes located in the sensor. One of them is feed with a high voltage DC, while the current generated in the other is transformed to voltage, conditioned and then acquired for being converted into a soot estimation. This layout is similar to other fast soot sensors, see e.g. [74, 75]. Even though these sensors have not reached the maturity of UEGO or NO<sub>x</sub> sensors, some companies have already introduced them in the market or are targeted for start of production in 2013.

**Sensors in development, other alternatives.** A relevant problem of fast NO<sub>x</sub>-ZrO<sub>2</sub> sensors is the low sensitivity when the NO<sub>x</sub> concentration is small, i.e. downstream of SCR or LNT systems; and the related cross-sensitivity to ammonia. In that field, it can be underlined the work by Ralf Moos [76–81].

An alternative is using the catalysts and/or substrate materials of the AT devices for determining the NO<sub>x</sub>, soot or ammonia by means of measuring the impedance variance of the material. This allows to profit the proper AT device for measuring, developing an integrated device. These are often called *in situ* monitoring devices. Moos *et al.* [78] show results for *in situ* sensors in order to detect loadings of oxygen in TWC, NO<sub>x</sub> in LNT, NH<sub>3</sub> in SCR-catalyst and soot in DPF, among others. The principal advantage of the approach is the simple and inexpensive setup required as the devices themselves are used as sensors.

Concretely for NO<sub>x</sub> sensing, see [79] and Groß *et al.* [81] who extend the use of integrating-type NO<sub>x</sub> sensors for measuring instantaneous NO<sub>x</sub> concentration, especially in the low ppm range (where NO<sub>x</sub> concentration is low), incorporating temperature based corrections in the measurements. Groß *et al.* also present results with respect to NO/NO<sub>2</sub> differentiation and cross-sensitivities with other gas components. Other alternatives for innovative sensors could be using electromagnetic waves [78] or radio-frequency signals which coincide with the oxidation/reduction rate of a TWC for determining oxygen loading without using lambda sensors [76].

NH<sub>3</sub> on-board sensors [77] are also being developed for closed loop urea dosing control, based on the same principles of the wide-band lambda sensors: the mixed-potential principle between electrodes, separated by a solid electrolyte, where an electric force is created. Delphi has already marketed

a commercial version where the reference electrode is in direct contact with exhaust gas [82]. Schönauer *et al.* [80] show results with mixed potential ammonia sensors and suggest some improvements in the technology.

### 2.3.2 Models and virtual sensors

Modelling or virtual sensing<sup>2</sup> is an important tool in the design, analysis and control of IC engines. It is not in vain that model-based strategies in ECU have each time higher responsibilities, overall in control and OBD functions [83]. Furthermore, models are quite helpful in the analysis and design phases: hardware and software in the loop (HIL and SIL) are usual because of the reduced cost and time-to-implement with respect to a pure experimental approach. However, the accuracy, reliability, computational cost and time effort of models depend highly on the structure, data used for calibration and the dynamic conditions. A classification on models for automotive is made in [84] and these are compared in Table 2.3:

- Mean value models (MV), which average the physical quantities over a time range, neglecting in-cycle variations. They can be further subdivided in data driven models (DDM) and physical mean value engine models (MVEM).
- Emptying-and-filling models (E&FM), based on first-principle equations and with certain capability for estimating in-cycle variations. They are usually crank-angle solved.
- Wave Action Models (WAM), able to quantify variables in 1D space and based on an Euler equation balance. These allow to represent wave effects (useful e.g. for valve lifting tuning).
- Computational fluid dynamics models (CFD), based on a detailed geometrical description of the engine volumes and the application of the Navier-Stokes equations. Due to the 3D nature of this approach, this tool is effective for simulating mixing models, i.e. EGR distribution around the cylinders or injection spray models. However, CFD characteristic times of computation can range from seconds to days, preventing its use for RT purposes.

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<sup>2</sup>The term virtual sensing is often used for RT on-board models for estimating variables that cannot be measured easily or their measurements are not reliable, in contrast with using physical sensors.

**Table 2.3.** Summary of the characteristics of different type of models used in the automotive, reproduced from [84]. MV: mean value, acq.: acquisition, CAS: crank angle solved, CAE: computer aided-engineering.

	DDM	MVEM	E&FM	WAM	CFD
Spatial resolution	0D	0D	0D	1D	3D
Time resolution	MV	MV	CAS	CAS	CAS
Typical acq. rates	0.5-100 Hz	1-200 Hz	>1 MHz	>10 MHz	1-100 MHz
Physical description	no/very low	low	low	medium	complex
Prediction capabilities	no	low	medium	medium	high
Computational cost	very low	low	medium	high	very high
Main application	control	control	CAE	CAE	CAE

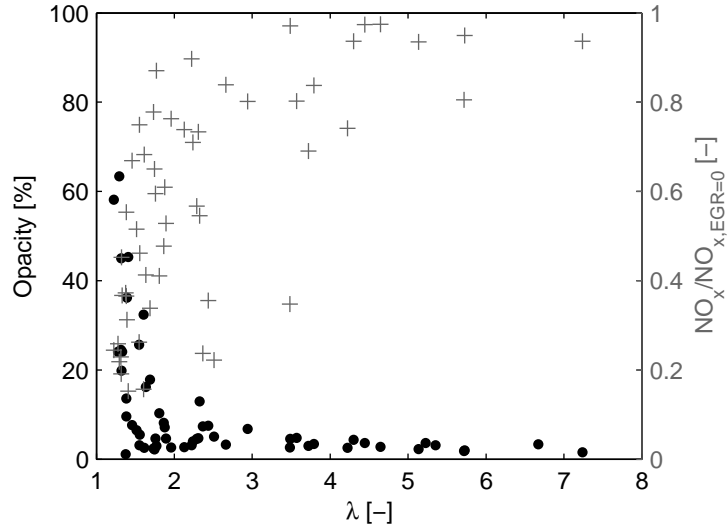
In order to design control oriented models (COMs), MVEM and DDM are the predominant approaches [85], basically for the relatively simple computation and the good accuracy achieved, especially when the training data set is appropriate, getting indeed higher accuracy than more complex models. MVEM and DDM rely on two main hypothesis: variables are averaged over the time, with computation (and acquisition) rates about 50 Hz<sup>3</sup>, and the number of states that these models can manage are reduced due to the limited computational resources available (mainly the ECU limitations). A straightforward classification of COMs is:

- Physical models,
- gray box models, which combine first principle equations with parameters, curves and maps to be fitted with engine data, and
- black box models, which are fitted only by data and with no physical based structure.

These approaches present advantages and drawbacks. In the following, COMs for  $\lambda$  and  $\text{NO}_x$  are discussed, as being core variables of the present work. The interested reader can find a review on proposed models for these variables in the dissertation by Schilling [86] with application to failure detection and isolation (FDI).

<sup>3</sup>For an engine that is spinning at 3000 rpm, that is equivalent to a rate of 1-100 Hz. A typical frequency for getting changes inside the combustion chamber could be in the order of 1° if talking about crank angle, and this makes a required frequency (at 3000 rpm) of  $50 \times 360 = 18$  MHz, quite higher than the typical rates used in MV models.





**Figure 2.7.** Opacity (left axis and black points) and nominal  $NO_x$  (gray +) emissions depending on  $\lambda$  for nominal tests in the TCCI engine used in Chapter 3.  $NO_x$  is normalised with the maximum  $NO_x$  value for each operating point, which corresponds to  $u_{egr} = 100$  (fully closed) for each pair of  $[n, m_f]$ .

### 2.3.2.1 $\lambda$ models

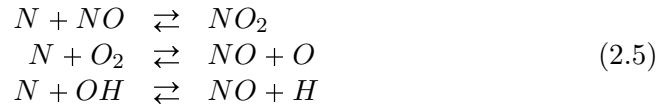
$\lambda$ , although not measuring emissions itself, is a key variable in order to control emissions in CI engines. Figure 2.7 shows the opacity and  $NO_x$  with respect to  $\lambda$  showing the clear opposite trade-off between both variables.  $\lambda$  can be directly estimated in the basis of  $\dot{m}_a$  and  $\dot{m}_f$ , see (2.4). Both signals can be measured from the ECU or estimated by a model. In the first case,  $\dot{m}_a$  is measured by a sensor and  $\dot{m}_f$ , inferred by the  $p_{rail}$  and  $u_{t_{mi}}$  (2.1a). This computation is simple and no specific model is necessary.

Alternatively, MVEMs can be proposed for calculating  $\lambda$ , which will be reduced to estimate the air mass flow at intake as  $m_f$  is often considered as an input for control models. Basic literature for principle equations of CI engines is in [16, 87, 88], while Wahlström [89] develops and tune a MVEM for control of EGR and VGT in diesel engines. In addition, other model types with higher or lower ability to compute on RT can be proposed, Tunestål and Hedrick [90] estimate  $\lambda$  using net heat release data and Arsie *et al.* [91] use neural networks (NN) (with SI application), among other examples [92, 93].

### 2.3.2.2 NO<sub>x</sub> models

The case of emissions, such as NO<sub>x</sub>, is more complex than for λ. NO<sub>x</sub> can be inferred by using predictive models that can be implemented on-engine for RT purposes such as engine or AT control and diagnosis. Furthermore, such models can be useful in an off-line basis for NO<sub>x</sub> simulation and prototype design. The physical based approach presents difficulties to be applied on RT but can be the basis for grey box approaches [94]. First, some words on NO<sub>x</sub> physics.

**NO<sub>x</sub> physical-based models.** The term NO<sub>x</sub> includes all nitrogen oxides but the nitric oxide (NO) is the predominant in diesel engines [87]. NO<sub>x</sub> formation is affected by three different mechanisms: thermal, prompt and from fuel-bound nitrogen [95]. The thermal mechanism is the most important in diesel engines where high temperatures benefit reaction of N and O<sub>2</sub> from air producing NO<sub>x</sub>. NO<sub>x</sub> formation physics in combustion and explosion processes were modelled by Zeldovych [96] in 1946, and formulated for IC engines by Lavoie *et al.* [97] in 1970 with the well-known extended-Zeldovych mechanism



According to Heywood [87], the typical characteristic times of the NO formation in diesel engines combustion is in the order of seconds and thus under the hypothesis of equilibrium of certain species, the  $dNO/dt$  can be fitted with the initial NO formation rate by the Arrhenius equation

$$\frac{dNO}{dt} = \frac{k_1}{T^{0.5}} e^{-k_2/T} [O_2]^{0.5} [N] \quad (2.6)$$

The strong dependency with the temperature  $T$  is clear: when  $T$  increases, NO (and thus NO<sub>x</sub>) increases exponentially. The other mechanisms can be relevant in some specific conditions such as LTC [98,99].

Cylinder conditions such as temperature, pressure and oxygen concentration [100–102] are the most important variables for determining NO<sub>x</sub> concentration.

Because of the cylinder severe conditions, price and signals problems have prevented to use in-cylinder sensors in commercial vehicles. Although in-cylinder pressure cost is one of the main factors burdening its application for automotive engines, the continuous improvements of pressure sensors and its

applications are justifying its nearby implementation [103]. The case of the temperature sensors is more complicated [37], and their use is not foreseen in applications. Therefore, the solution for estimating the in-cylinder conditions on RT is then using virtual sensors.

Physical models for NO<sub>x</sub> estimation are often based in the heat release, using pressure sensor signal for estimating flame temperature ( $T_f$ ) in the cylinder. The problem is not trivial and a multi-zone discretisation is advisable. Furthermore, the residual gases (internal EGR) affect the process and these are not always easy to estimate. An usual solution to the problem is applying mass and energy conservation equations to each zone (walls, injector neighbourhood, etc.) in conjunction with heat transfer equations, among them it stands out the heat transfer to cylinder wall [104], which can be approached by using Woschni equation [105]. Finally, NO<sub>x</sub> emissions for both Diesel and gasoline engines are calculated by using the extended Zeldovych mechanism. Good examples for estimating NO<sub>x</sub> by using heat release are [106–109]. In the case of SI engines, some authors use the ionisation current on the spark for approaching the pressure [110, 111], but this is not available on the diesel engine. Nevertheless, the accuracy of the prediction is not as satisfactory as expected, and the drift cannot be eliminated.

There is an open discussion about using time or crank angle based models for the NO<sub>x</sub> prediction. For the case of pressure based models, crank angle sampling seems more logical as volume can be easily linked with pressure trace. However, these models require heavy calculations and big memory resources. The time required for completing one engine cycle is often bigger than the characteristic time of the engine. In order to overcome this limitation some authors have proposed simplifications. Guardiola *et al.* [108] develop a semi-physical discrete event model based on the heat release calculation but considering only one zone as the main contributor to the NO<sub>x</sub> formation; and the process is supposed adiabatic, approaching  $T_f$  with the adiabatic temperature of the process. This approach requires specific corrections, especially when the combustion temperature is low. Other examples are Westlund and Åmströng [112] who present a fast physical model for NO<sub>x</sub> and soot, or Arsie *et al.* [94] who present a hierarchical model structure for engine control design with different models and layers, ranging from physical based to mean value approaches.

**Control-oriented models for NO<sub>x</sub>.** Literature about COMs for NO<sub>x</sub> is extensive. Schilling [86] gives a short review on different COMs for NO<sub>x</sub> and presents a NO<sub>x</sub> virtual sensor (see also references therein). His work is based

on deriving maps from complex models [67, 113] or engine data [114, 115], building a model that is easy to integrate into the engine ECU, but with limited extrapolation capabilities. In commercial ECUs the prevailing approach is to use look-up tables to model nonlinear and operating point dependent behaviours because of the simple programming in spite of the intensive tuning effort; even though other possibilities can be commented. Winkler *et al.* [116] design a virtual NO<sub>x</sub> sensor for SCR control and diagnosis. Black box models rely on system identification [117] but are often operating point dependent and its adaptation is not an easy task. Hirsch *et al.* [118] present a gray box model for NO<sub>x</sub> and PM. Other non-linear approaches are Takagi-Sugeno fuzzy models [114, 119], Hammerstein-Wiener (HW) [120], or NN [121–124].

### 2.3.2.3 Identification of sensor and physical dynamics

Dynamic models describe the evolution of a state vector  $x$  along the time  $t$  and subjected to a specific set of inputs  $u$

$$\dot{x} = f(x, u, t) \quad x \in \mathbb{R}^{n_x}, u \in \mathbb{R}^{n_u} \quad (2.7a)$$

$$y = g(x, u, t) \quad (2.7b)$$

The identification is the technique of designing and tuning the dynamic models by means of a data set  $\{u, y\}_{t_1}^{t_{end}}$  over a certain time  $t_{end}$ ; see [125, 126] for comprehensive references on system identification. A common procedure for identification is defining different black box linear model structures (AR, ARX, ARMAX, etc.) based on polynomial relationships with a certain order and solving the linear regression problem, usually by means of a least squares (LS) formulation. In addition to the model parameters, the selection of an appropriate order of the system is key for an adequate solution. If the system is non-linear, specific structures should be utilised, e.g. piece-wise affine (PWA), linear parameter varying (LPV) or HW models, among others.

**Sensors identification for on-board strategies.** Most of the dynamic models with physical meaning used in engineering, as for instance sensor models, respond to n-order linear filters with a certain delay  $\tau$  [127], at least in the selected areas of performance. In the case of exhaust gas concentration sensors, 1st order linear models can be sufficient for representing the sensor response in the s-domain [44]

$$G(s) = e^{-s\tau} \frac{k}{1 + Ts} \quad (2.8)$$

where  $T$  is the response time and  $k$  is the gain. Parameters can be fitted by LS. However, actual sensor responses are non-linear and parameters usually vary, especially the delay or dead-time  $\tau$ , due to the hardware itself, the engine operating conditions and ageing. These variations can be modelled by scheduling strategies, where parameters are stored in look-up tables, or functions [128] typically depending on  $n$  and  $m_a$ .

Another possibility is using model structures with physical insight: e.g. Wang [129] identifies the oxygen sensors dynamics by means of a physical based model and Zhuiykov [130] describes and models the electrochemistry of  $ZrO_2$  gas sensors. Adaptive filtering can also be used for taking into account the ageing and dispersion. Furthermore, if the signal processing is made off-line, non-causal deconvolution techniques may also be used, see e.g. [131].

DDMs are often based on static relationships such as look-up tables and algebraic equations. In addition, the dynamics are often modelled by linear filters and the correct identification is critical for the adaptive filtering algorithms presented hereinafter.

### 2.3.3 Adaptive filtering

Sensors present problems when using the raw output signal in RT functions, in particular the delay from engine to sensor and the response time of the sensor. With respect to models, two problems can be underlined when working with COMs. On one hand, the model accuracy is driven by the collection of the appropriate data and calibration of all the parameters. This is a hard and time consuming task. In fact, ECU has a big number of maps and parameters for engine and vehicle management. On the other hand, independently of how well the model has been calibrated there is inevitably a drift between the system and the model as the surrounding conditions changes and the engine ages. DDMs are highly sensitive to the calibration data set and will have problems with the ageing, manufacturing discrepancies, slowly varying parameters and other non-modelled variables. These affects also to dynamic models identified for both sensor and physics filtering.

**Data fusion.** Data fusion (DF) utilises information from different sources for providing a better estimation of a given variable. These information sources can come from sensors, models and/or other DF algorithms. DF has been used for many years in different engineering fields, such as e.g. inertial navigation of satellites and missiles [132] or automotive applications in intelligent transportation systems [133, 134] related with traffic and driver/road assistance,

among others. Different DF methods can be used as El Faouzi *et al.* [133] propose: statistical, ranging from the simplest arithmetic mean to weighted combinations or data mining techniques [131]; probabilistic, such as Bayesian approaches, maximum likelihoods methods and Kalman filter (KF); and artificial intelligence including NN or genetic algorithms. A more exhaustive classification and discussion on challenging problems of multisensor DF is in [135]. From DF methods, those which are applied adaptively for bounding error of parameters, signals and parameters can be included as methods for adaptive filtering, fusion or learning [136].

The present work centres on the use of Kalman filter (KF) [69] based tools in order to observe exhaust gas concentration variables in CI engines. The KF is probably the most extended adaptive filtering algorithm used in engineering [137] and provides a systematic and simple way of manipulating linear dynamic models and engine variables by supposing *a priori* knowledge of the measurement and noise statistics, modelled by Gaussian distributions. This property allows to track the estimation error by means of a covariance matrix  $\mathbf{P}$  that also tracks the ageing of the states, i.e. if the error in a given state is foreseen to be higher or lower depending on when was updated the last time and how much. The KF minimises this expected error by solving an iterative Riccati matrix equation and setting out a Kalman gain ( $K$ ) for correction. The linear structure of the filter makes it appropriate to ECU-oriented approaches, even though simplifications shall be considered when manipulating large state vectors. Furthermore, depending on the available data, different fusion structures can be programmed as Gao and Harris [137] show.

Other filtering alternatives, such as RLS (recursive least squares) or proportional methods can also be applied and in some cases lead to similar solutions. Input observers by means of measurements of the states and outputs can also be addressed for estimating engine variables [138,139]. The interested reader is referred to Simon [140] and Höckerdal [141], the former for finding a complete and broader view on optimal state estimation and KF-based algorithms for engineering applications, and the latter for a key precedent and motivation for the current work.

In the next, and due to the importance of the KF in this work, main equations are recalled although the reader familiarised with control engineering could skip this paragraph.

**Kalman filter.** In the setting the data are assumed to be generated by the following discrete time system

$$x_k = f(x_{k-1}, u_k) + w_k \quad (2.9a)$$

$$y_k = g(x_k, u_k) + v_k \quad (2.9b)$$

where  $x_k \in \mathbb{R}^{n_x}$  represents the state vector,  $u_k \in \mathbb{R}^{n_u}$  the input vector,  $y_k \in \mathbb{R}^{n_y}$  the output vector. If  $f$  and/or  $g$  are non-linear, a previous linearisation step is required for the filter and then, elements  $ij$  of  $\mathbf{F}_k$  and  $\mathbf{H}_k$  are obtained

$$\mathbf{F}_{k,ij} = \left. \frac{\partial \mathbf{f}_i}{\partial \mathbf{x}_j} \right|_{\mathbf{x}=\hat{\mathbf{x}}_k} \quad \mathbf{H}_{k,ij} = \left. \frac{\partial \mathbf{g}_i}{\partial \mathbf{x}_j} \right|_{\mathbf{x}=\hat{\mathbf{x}}_k} \quad (2.10)$$

being  $\mathbf{F}$  the linearised process matrix and  $\mathbf{H}$  the linearised output matrix. From now, the discussion is valid both for linear and non-linear systems and KF will be used referring to Kalman filter based methods, including the non-linear Extended Kalman Filter (EKF) [140] and the standard one.

Noises in (2.9)  $w_k \in \mathbb{R}^{n_x}$  and  $v_k \in \mathbb{R}^{n_y}$  are assumed to be independent and both generated by Gaussian distribution with zero mean and covariance matrices  $\mathbf{Q}_k$  resp.  $\mathbf{R}_k$ , defined by

$$E[w_k w_k^T] = \mathbf{Q}_k \quad (2.11a)$$

$$E[v_k v_k^T] = \mathbf{R}_k \quad (2.11b)$$

In many applications these are often chosen to be constant, i.e.  $\mathbf{Q}$  and  $\mathbf{R}$ , and diagonal.

Then,  $\hat{x}_k \in \mathbb{R}^{n_x}$  is the observation of the state vector  $x_k$

$$\hat{x}_{k|k-1} = f(\hat{x}_{k-1}, u_k) \quad (2.12a)$$

$$e_k = y_k - g(\hat{x}_{k|k-1}, u_k) \quad (2.12b)$$

$$\hat{x}_k = \hat{x}_{k|k-1} + K_k e_k \quad (2.12c)$$

where Kalman gain  $K_k$  is solved by the following iterative equation

$$\mathbf{P}_{k|k-1} = (\mathbf{F}_k \mathbf{P}_{k-1} \mathbf{F}_k^T + \mathbf{Q}) \quad (2.13a)$$

$$K_k = \mathbf{P}_{k|k-1} \mathbf{H}_k^T (\mathbf{H}_k \mathbf{P}_{k|k-1} \mathbf{H}_k^T + \mathbf{R})^{-1} \quad (2.13b)$$

$$\mathbf{P}_k = (\mathbf{I} - K_k \mathbf{H}_k) \mathbf{P}_{k|k-1} \quad (2.13c)$$

and matrix  $\mathbf{P}_k$  is the covariance matrix of the state estimate error [126]

$$\mathbf{P}_k = E[x_k - \hat{x}_k][x_k - \hat{x}_k]^T \quad (2.14)$$

**About the methods in this work.** From all the possible applications of the KF for the gas concentration estimation, the following are exploited in the present work:

- Dynamic estimation of engine variables by means of fusing a fast reference, often a model, and a slow but accurate sensor, for developing robust virtual sensors or bias tracking algorithms; and
- on-line adaptation of models, especially the online updating of look-up tables as basic structures in COMs.

These algorithms, although applied in the present work to  $\lambda$  and  $\text{NO}_x$ , can be used for estimating other relevant engine quantities, such as the intake air mass flow or the volumetric efficiency [142, 143], just considering the appropriate state-space model. In the following, the state of the art related with these applications on automotive is reviewed.

### 2.3.3.1 Adaptive estimators based on the KF

Figure 2.8 shows a model signal for  $\lambda^{-1}$ , based on the injected fuel mass  $\dot{m}_f$  estimated by the electronic control unit (ECU) and the air mass flow  $\dot{m}_a$  determined from a hot wire anemometer. Comparing to the sensor, model is faster and non-delayed but presents a bias with respect to the sensor steady-state value.

Algorithms based on bias tracking take a fast model that keeps high frequency components of the considered variable while a slow but steady-state accurate sensor permits to correct model drift. The vector  $x_m(t) \in \mathbb{R}^{n_x}$  contains model signals and  $z(t) \in \mathbb{R}^{n_z}$  the sensor signals. The drift is modelled with a vector  $\theta(t) \in \mathbb{R}^{n_x}$ , which contains model biases and varies with time often slowly. If  $x_r(t) \in \mathbb{R}^{n_x}$  is the vector containing the actual values of the states, then the true bias  $\theta$  is

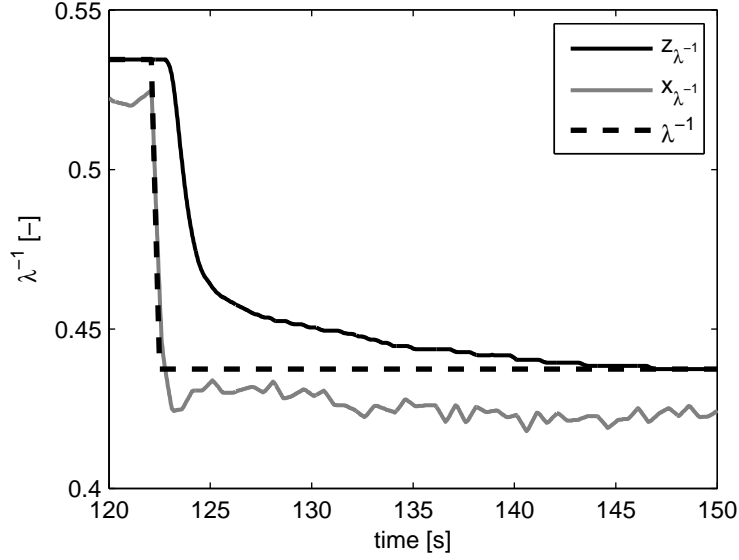
$$\theta(t) = x_r(t) - x_m(t) \quad (2.15)$$

but the problem here is that  $x_r$  is usually measured by a sensor  $z$

$$z(t) = g(x_r(t - \tau)) \quad (2.16)$$

and in general will have a certain delay  $\tau$ . The discussion on identifying the function  $g$  and delay  $\tau$  was made before in Section 2.3.2.





**Figure 2.8.** Different  $\lambda^{-1}$  estimations during a load transient in the tested diesel engine (see Section 3.2):  $z_{\lambda^{-1}}$  is the output of a  $\text{NO}_x$  sensor located at the turbine outlet,  $x_{\lambda^{-1}}$  is the output of the model presented in Section 4.2 and  $\lambda^{-1}$  is the foreseen actual value.

In order to design a drift correction algorithm, an augmented state-space model is written. An extra-state is used for tracking the given error ( $\theta$ ), creating a wide state vector  $x^w$  with drift vectors and the original state vector  $x \in \mathbb{R}^{n_x}$ , which contains the estimations of  $x_m$ ,

$$x^w(t) = \begin{bmatrix} \theta(t) \\ x(t) \end{bmatrix} \quad (2.17)$$

and the state-space model

$$\dot{x}^w = f(x^w, x_m, W, t) \quad (2.18a)$$

$$z = g(x^w, x_m, V, t) \quad (2.18b)$$

is solved by adding noises  $W(t) \in \mathbb{R}^{2n_x}$  and  $V(t) \in \mathbb{R}^{n_z}$  to the model and sensor output respectively.

This formulation, with different variations, is made by different authors in order to observe the bias between two references, usually a model and a sensor. Alternatively, the bias can be applied to the model and/or sensor by shifting

the references when proceeding; for the latter the sensor is not steady-state accurate and a steady-state accurate reference is required [144].

The classical approach for modelling the drift is supposing a slow variation of the state by the dynamic equation

$$\dot{\theta} = 0 \quad (2.19)$$

and supposing that states noise  $W$  is low with respect to output noise  $V$ , for having a smooth correction.

There exist different references with similar models as described above for estimating engine variables. The majority of them uses the EKF due to the non-linear equations involved, some references from 2005 are commented below:

- Hsieh and Wang [62] use an EKF in order to estimate  $\text{NO}_x$  up, in between and downstream of a two-series SCR configuration and consider the ammonia cross sensitivity of a  $\text{NO}_x$  sensor located in between.
- Höckerdal [142] uses an augmented model for observing the sensor bias with an air mass flow sensor application.
- Alberer and del Re [68] use a KF for correcting an oxygen measurement made by a UEGO sensor located upstream of the turbine in a TCCI engine, where pressure effects in the sensor are taken into account.
- Polòni *et al.* [145] compare two different sensors configuration in a TC CI engine for correcting states in a MVEM of the air path, designing a closed loop MVEM. They use an EKF for tracking the bias on the considered states and then simulate the model with the corrected variables. References included in this work are interesting and show other different observer designs.
- Other examples are [17, 128, 146–151] who design observers for different relevant engine quantities, such as intake manifold temperature, soot, individual in-cylinder air to fuel ratio or engine torque among others, with diagnosis or control applications.

**Drift models.** However, there are cases where the considered bias, e.g. model bias  $\theta$  variation does not depend only on time

$$\frac{d\theta}{dt} = \frac{\partial\theta}{\partial t} + \frac{\partial\theta}{\partial n} \frac{dn}{dt} + \frac{\partial\theta}{\partial m_f} \frac{dm_f}{dt} + \dots \quad (2.20)$$

and although the bias variation associated with the system drift ( $\partial\theta/\partial t$ ) is expected to be slow, the actual variation of the bias may be very fast, due to the ability of the engine of performing fast transition between operating conditions (usually defined by  $n$  and  $m_f$ ). This can be solved by using specific models for the bias, where a possible solution may be using look-up tables.

In the following, the use of adaptive virtual sensors is discussed.

### 2.3.3.2 Online adaptation of models

In addition to the bias and ageing, the basic structure of models lead to errors due to uncertainties and non-considered inputs. Therefore, the model error might be tracked and distributed between the model parameters. This could be made on an offline basis if the variables are stored, and then the model signals and parameters might be updated by means of executing an optimisation method. An alternative is using adaptive filtering for directly updating the model parameters online. Schilling [86] makes a selection on parameters from a virtual sensor for  $\text{NO}_x$  and  $\lambda$ ; and designs an adaptive virtual sensor for their online adaptation. Another possibility is introduced by Pòloni *et al.* [145] where a set of individual observers are designed for some states of the model. However, MVEMs often pose a mixed structure with different parameters, curves, tables and dynamic equations. The problem of updating all elements is not straightforward and must be faced carefully. The design of observers for updating parameters and tables can be used for tuning or updating DDMs and MVEMs, although stability and robustness of the solution should be issued. Due to the relevance on the present work, the updating of look-up tables is considered.

Look-up tables<sup>4</sup> allow engineers to easily model systems that present complex expressions by means of mapping outputs with a set of nD heuristic array structures. Look-up tables are used in automotive for different purposes [152]: maps of engine parameters as function of the operating-point conditions, often  $m_f$  (or torque) and  $n$ ; in order to provide set-points for the controllers; or for gain scheduling, among others. Furthermore, the manipulation and interpretation of look-up tables is simple and the usual offline procedures for calibration are based on LS methods. Linear interpolation is usual for computing the outputs. Anyhow, the tables are of course subjected to drift and

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<sup>4</sup>Usual dimensions of look-up tables are 2D, while 1D tables are often named as curves, but in fact they can be considered as 1D look-up tables. Using higher dimensional grids is not usual due to the involved computational burden, excepting specific cases such as for representing different combustion modes.

ageing when running online. Adaptive filtering can be used for correcting the drift and/or tuning parameters in an offline basis.

**EKF for updating look-up tables.** Höckerdal *et al.* [143] propose an EKF for updating table elements where they are treated as states to be observed and updated by measuring the considered output. In this approach,  $\mathbf{P}$  represents the ageing of the states, defining a proper  $K$  for updating in such sense that locally inactive elements see how their related variance grows monotonically and active elements see how this is modified depending on the weighted distance to the considered element. The interpolation principle weights the correction and  $\mathbf{P}$  updating: more excited elements will tend to have lower variance than the less excited ones. A bounded limit on  $\mathbf{P}$  should be applied for engineering applications in order to avoid robustness problems.

The first limitation of this algorithm is that although inactive elements do not really affect the updating in a given iteration, EKF manipulates all elements of  $\mathbf{P}$ . The global stability of the method relies on the observability (linked with activeness) of the states and although during each iteration only 4 elements (in a 2D example) are locally observable and the remaining are unobservable, the calculation involves all grid areas. A second limitation is that system matrices vary with time making impossible to derive a steady-state Kalman filter [153]. This forces to solve the Riccati equation at every instant for inferring  $K$ , involving a huge memory and computational resources. By utilising a numerical example, an EKF for updating a look-up table with 20-by-20 elements gives rise to a covariance matrix of 400-by-400 elements and thereby also a significant computational burden when solving the Riccati equations, which grows rapidly with the grid dimensions. Therefore an important discussion of the work is how the KF can be used or modified to get an efficient updating procedure for look-up tables without an important loss of properties.

**RLS and other approaches.** An alternative solution for estimation is the use of LS techniques, see e.g. Peyton and Muske [154] who use a RLS with a forgetting factor for updating look-up tables. The programming of a RLS is fairly simple and requires that table grid and data is well distributed for avoiding robustness problems (as the authors literally claim). Vogt *et al.* [152] present the normalised least mean square (NMLS) method optimised for online-adaptation of tables and propose an interesting simplification on this method by separating non-active and active elements of the considered table, aspect that is also exploited in this work but for EKF-based methods.

The results are good with much lower computation involved but with a slower convergence with respect to the standard RLS.

Another interesting contribution due to the simplicity of the formulation is made by Wu [155], who treats the problem of table updating as a reverse interpolation problem. A proportional weighting is then used for updating all the elements that were involved in the previous interpolation calculation. The author cites it literally as *multiple nodes proportional distribution*. This weighting is calculated only from the inputs and previous outputs. In that way, the method does not take into account any noise in the measurement, and can be considered as sub-optimal if comparing with the KF, but with a much lower computation and memory weight because only the table values related with observable elements are updated.

This work presents two approaches (SKF and SSKF) with some characteristics of the commented above in order to update look-up tables efficiently.

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# Chapter 3

## System setup and sensors characterisation

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### 3.1 Introduction

This chapter is divided into two parts. The first part, which comprises Sections 3.2 and 3.3, is devoted to the description of the experimental configuration and tests used for this work. Section 3.2 describes the experimental set-up used in the present work, including the relevant characteristics of the

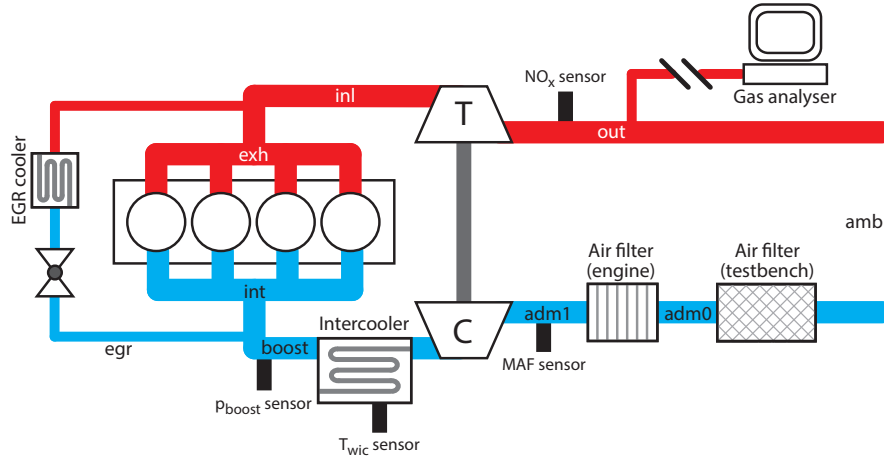
engine, sensors and test cell equipment, while the Section 3.3 presents the steady-state and dynamic tests performed to tune and validate the methods.

In the second part of the chapter, the use of online methods for characterising  $\text{NO}_x$  and  $\lambda^{-1}$  output from exhaust gas concentration sensors is emphasised.  $\text{NO}_x$  output is characterised by a novel method based on SOI steps, while  $\lambda^{-1}$  output is characterised by performing injection steps.

## 3.2 Experimental set-up

Test engine is an automotive 2.2-litre 4-cylinder common rail DICI engine with a sequential parallel turbocharger [1], certified according to EURO IV standard (model DW12B from PSA). Engine specifications are shown in Table 3.1. A bypass-allowed ECU permits varying injection parameters, such as  $p_{rail}$ , SOI and main injection duration ( $t_{mi}$ ), and also boost and EGR control set-points. The engine is installed on an engine test bench and coupled to a variable frequency eddy current dynamometer that allows carrying out dynamical tests. Standard air intercooler is substituted by a water intercooler for a proper heat extraction on the intake air mass flow. The usual experimental set-up used for generating the data in this work is shown in Figure 3.1, where the AT devices are removed and the main sensors are sketched. A back-pressure valve is installed downstream of the turbine outlet and commanded for replicating the pressure drop in the exhaust line of series engines. The nomenclature for the engine positions is clarified in the Figure 3.1 and in the list of variables at the beginning of the document as well. In some specific tests, the sensor layout is modified and AT devices are included, however they are indicated where proceed. The main actuators on the engine are described in the Table 3.2.

**The rapid prototyping system.** In order to avoid the shifting and causality problems with the signals, the system layout shown in the Figure 3.2 is used for acquiring from and commanding signals to the engine. A real time (RT) PXI system (from National Instruments) is used for the acquisition of data from series and experimental sensors. The PXI is connected via CAN with a ES910 hardware from ETAS for acquisition and bypass of the ECU signals, allowing to design control routines and skipping ECU logic. In addition to series sensors (already connected with the ECU), other sensors are installed on engine and linked with the PXI for acquisition and providing feedback for control strategies. Labview software is used for managing PXI, while



**Figure 3.1.** Engine layout with the main sensors used in this work (pressures, temperatures in different positions are also measured but are not shown for clarity).

**Table 3.1.** Main characteristics of the DW12B engine from PSA.

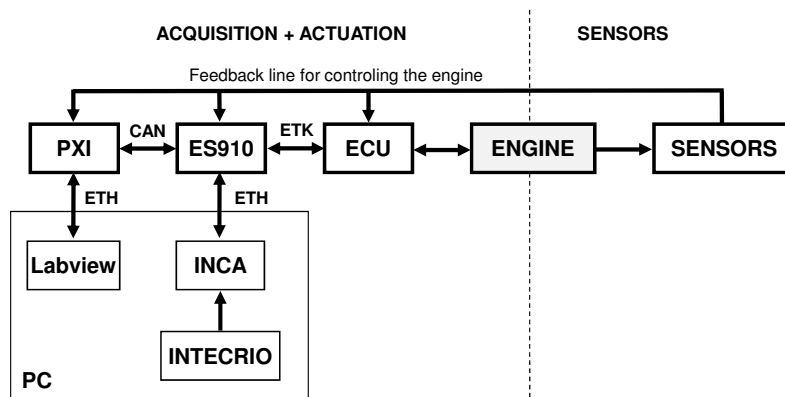
Stroke ( $S$ )	96 mm
Bore ( $D$ )	85 mm
$S/D$	1.129
Number of cylinders	4
Displacement	2179 cm <sup>3</sup>
Turbocharging system	Sequential parallel
EGR	HP
Injectors	Solenoid (common-rail)
Valves by cylinder	4
Maximum power	125 kW@4000 rpm
Compression ratio	17:1

INCA software is utilised for interfacing with the ES910. INTECRIO software permits to program the engine variables for the bypass.

When bypassing, the time needed by the PXI software for writing into the ECU channel is 20 ms, and the same for reading in the other direction. Then, the maximum expected error between commanding and reading a signal in the PXI is 40 ms for the round path. Lags are around 20 ms when storing data by INCA software (used for sensors characterisation and connected with the ES910). Acquisition frequencies are 100 Hz for integrated sensors and between

**Table 3.2.** Main actuators used in the experimental setup.

EGR command	$u_{egr}$	%
WG command	$u_{wg}$	%
Rail pressure command	$u_{prail}$	bar
Main Injection duration (MI)	$u_{t_{mi}}$	$\mu s$
Start of injection of the MI	$u_{soi}$	$^{\circ}BTDC$

**Figure 3.2.** System layout for rapid prototyping (fast acquisition, bypass and control) and integration of new sensors for testing the engine.

10-100 Hz for ECU signals (fixed by manufacturer). Commanding frequencies depend on the test and variable.

### 3.2.1 Engine sensors

Besides the commercial sensors, a full set of sensors is installed on-engine in order to measure full operating conditions (pressures, temperatures, mass flows, engine and turbo speeds, torque and fuel consumption), atmospheric conditions and emissions. Some of them are measured redundantly for calibration. All sensors are tested and calibrated with standard procedures, both in steady and transient-state. Table 3.3 shows the main series and test cell sensors harnessed for measuring engine conditions. Due to the relevance in this work, gas concentration sensors are described separately.

**Table 3.3.** *Sensors used in the experimental setup (gas concentration sensors are discussed below).*

Series sensors	Symbol	Location
Air mass flow (MAF) [kg/h]	$\dot{m}_a$	<i>adm1</i>
Boost pressure [bar]	$p_{boost}$	<i>boost</i>
Engine speed [rpm]	$n$	<i>crank</i>
EGR valve position [%]	$x_{egr}$	<i>EGR valve</i>
WG position [%]	$x_{wg}$	<i>WG</i>
Rail pressure [bar]	$p_{rail}$	<i>CR</i>
Test cell sensors	Symbol	Location
Torque [Nm]	$M_e$	<i>crank</i>
Air mass flow (sensyflow) [kg/h]	$\dot{m}_a^{cell}$	<i>adm0</i>
Fuel (gravimetric) [kg/h]	$\dot{m}_f^{cell}$	<i>external</i>
Pressures [bar]	$p$	<i>runners</i>
Temperatures [°C]	$T$	<i>runners</i>
In-cylinder pressure [bar]	$p_{cyl}$	<i>cyl</i>

### 3.2.1.1 Exhaust gas concentrations sensors

Different gas concentration sensors are installed for designing and validating the algorithms. These can be summed up in research grade and on-board (experimental and commercial) sensors. From the first type, a gas analyser is used for measuring gas concentrations and an AVL 439 opacimeter for measuring the opacity at the exhaust (not relevant in this work). From the second type,  $\lambda$  sensors and  $\text{NO}_x$  sensors are highlighted. When necessary, redundancy is applied to sensors for comparing signals in different locations and indeed different units, e.g. the measurement of  $\lambda^{-1}$  before and after the turbine. Appendix 3.A summarises the technical data of these sensors according to the suppliers.

**Gas analyser.** A Horiba MEXA 7100DEGR gas analyser [2] is used as calibration standard and among others, the following relevant concentrations are measured:  $\text{CO}_2$  at the intake and exhaust,  $\text{NO}_x$  at the turbine outlet and  $\text{O}_2$  at the intake manifold and turbine outlet. The engine-out  $\lambda^{-1}$  and  $\text{NO}_x$  are inferred by locating the gas probe downstream of the turbine.  $\lambda^{-1}$  is calculated from the  $\text{O}_2$  measurement, estimated by means of a magneto-pneumatic detection (MPD), while  $\text{NO}_x$  is measured on a dry basis, by means

of a heated chemiluminescence detector (HCLD) with a NO<sub>2</sub>/NO converter. Calibration gases are used for the calibration of the system (zero and span).

Three key factors do not allow to use this system as a dynamic calibration method for on-board sensors:

1. Measurement system response time, which includes the signal rising time, is slower than the tested on-board sensors.
2. A long line is used for providing exhaust gas to the gas analyser (12m). This long line causes not only a transport delay, but also can cause diffusive effects.
3. Soot-free gas is needed in order to prevent sensor damage; therefore a filter is used in the line [3]. In transient engine operation, filter contributes to substantial emission signal distortion.

It is worth noting that some studies combine experimental data and phenomenological modelling approach to synchronise [4,5] or to restore [6–8] the signals from slow response analysers during transient engine operation. This is done by analytical means, using simple gas diffusion models and perfect gas mixing models, or through a control theory approach to the problem. Section 3.4 shows static and dynamic results by using the gas analyser for characterising the NO<sub>x</sub> sensor.

**UEGO sensor for measuring  $\lambda^{-1}$ .** This on-board sensor may be located in different positions in order to evaluate the possibilities of using measured  $\lambda^{-1}$  in models and adaptive algorithms. The proposal in this work for estimating engine-out  $\lambda^{-1}$  is utilising a NO<sub>x</sub> sensor, capable of measuring also  $\lambda^{-1}$ , while avoiding the use of a UEGO sensor (used for some production engines). Anyway, the sensor is used for helping some explanations and as a reference for comparison with the NO<sub>x</sub> sensor. Section 3.2.1.2 presents different alternatives for measuring intake  $\lambda^{-1}$  from UEGO sensors and the gas analyser.

**ZrO<sub>2</sub>-based sensor for measuring  $\lambda^{-1}$  and NO<sub>x</sub>.** The availability and the trade-off for locating NO<sub>x</sub> sensors in production diesel engine cars varies depending on the final application. In HD engines, the NO<sub>x</sub> sensor is becoming usual, and the normal solution is locating one NO<sub>x</sub> sensor downstream of the SCR for diagnosing and managing the system, while exhaust NO<sub>x</sub> is observed (saving the need of using more than one sensor unit). In LD, the sensor will

become necessary for EURO VI diesel engines with a similar trade-off to that of HD (although LNT might be used instead of SCR), while some luxury cars already mount two  $\text{NO}_x$  sensors for providing a valid reference before and after the AT devices.

In this work, the AT devices are removed in the standard configuration and the sensor is located at the turbine outlet. The capability of measuring  $\lambda^{-1}$  permits to work without UEGO sensors. It should be clarified that even though  $\lambda^{-1}$  is not really necessary for the control of diesel engines, it is advised for emissions control (soot and  $\text{NO}_x$ ) and may be used for correcting injector drift or bounding injection (on the basis of the smoke limiter actuation). Anyway, the sensor response must be assessed and this is performed in Section 3.4.

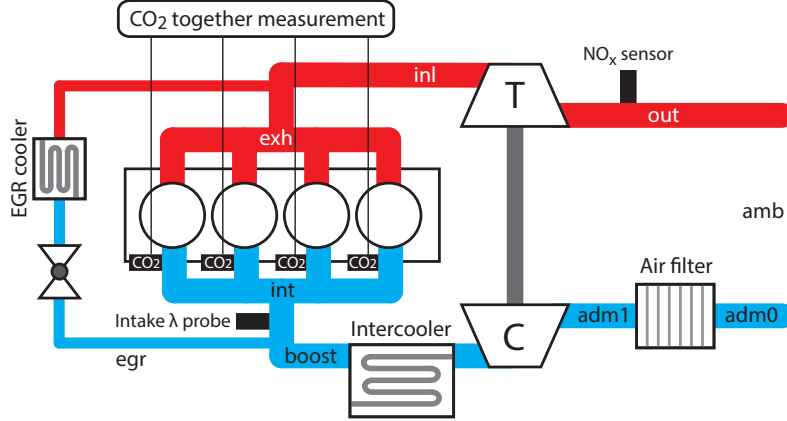
### 3.2.1.2 Intake gas concentration sensors

Due to the strong dependency of  $\text{NO}_x$  emissions on  $[\text{O}_2]_{int}$ , the measurement of  $[\text{O}_2]_{int}$  is of major importance. However, the lack of homogeneity of the intake mix stands a major problem in some engines. In this work, the methodology presented in [9] has been used for analysing the homogeneity of the mix. The method is based on feeding the gas analyser from different probes, as sketched in Figure 3.3:

- Measurement of the  $[\text{CO}_2]_{int}$  just before the separation of the manifolds at the intake by the gas analyser,
- extractions in the four individual cylinder intake runners and measurement by the gas analyser (cylinder 1 to 4  $[\text{CO}_2]_{int}$  independent measurements),
- measurement of a sample result of the mix of the 4 extractions and made by the gas analyser (1 measurement, from a rail where gas coming from all the probes is mixed).

When using the gas analyser, the  $[\text{CO}_2]_{int}$  [%] is measured with a non-dispersive infrared method (NDIR). As the intake gas is a mix of ambient air and the results of the combustion of this air with fuel, the  $[\text{O}_2]_{int}$  is linearly correlated with the  $[\text{CO}_2]_{int}$  (the higher the  $[\text{CO}_2]_{int}$ , the lower the  $[\text{O}_2]_{int}$ ).

Additionally, and as a real time alternative, the measurement of the intake  $\lambda$  by means of the UEGO sensor located in the intake manifold has been implemented. The sensor has been located with a certain distance upstream



**Figure 3.3.** Engine layout showing the main sensors used for measuring  $[O_2]_{int}$  or equivalently  $[CO_2]_{int}$ .

the separation between individual cylinder runners for avoiding separation effects (1 measurement).

UEGO sensor provides the ion pump current  $I_p$  [mA]. The  $I_p$  signal is previously corrected with the effect of  $p_{boost}$  which affects to the oxygen partial pressure measurement. Furthermore,  $I_p$  is related with  $\lambda^{-1}$  (and thus  $[O_2]_{int}$  and  $[CO_2]_{int}$ ) by a curve given by the manufacturer.

A set of measurements is made in order to evaluate the different alternatives and to determine the best method to estimate the intake gas concentration. In total 4 operating points defined by the pairs  $[n \ m_f]$  are studied:  $[2500,35]$ ,  $[2500,7]$ ,  $[1500,20]$  and  $[1500,7]$  (units are rpm and mg/str respectively) with the rest of inputs fixed by the standard engine calibrations.

The results of the study are summarised in Table 3.4 for the 4 operating points. Even though there exists an important dispersion among the cylinder samples (due to not utilizing a mixer that guarantees the homogeneity of the fluid at intake), the comparison of the mean value for the 4 cylinders with the measurement from the rail (Cyl 1-4 probe) is satisfactory. Nevertheless, the  $[CO_2]_{int}$  measured just before the cylinder manifolds is highly affected by the lack of homogeneity in the mix and the signal presents a significant error when comparing with the rest of alternatives. The results obtained by the UEGO sensor are acceptable, despite of a slight deviation lower than that of the gas analyser intake extraction. However, the Bosch sensor presents a saturation when  $I_p > 2.3$  mA (low  $[CO_2]_{int}$  and high  $[O_2]_{int}$ ) as shown



**Table 3.4.** Comparison of different sensor outputs for determining the effective intake  $[CO_2]_{int}$  (and thus  $[O_2]_{int}$  and  $\lambda^{-1}$ ) in the cylinder. std: standard deviation.

Operating point $[n, m_f]$	[2500,35]	[2500,7]	[1500,20]	[1500,7]
$[CO_2]_{int}$ [%]	1.37	1.07	1.27	2.30
$[CO_2]_{int}$ 1 [%]	1.14	0.98	2.38	2.62
$[CO_2]_{int}$ 2 [%]	1.65	1.09	2.46	2.92
$[CO_2]_{int}$ 3 [%]	2.07	1.11	1.89	2.64
$[CO_2]_{int}$ 4 [%]	1.89	1.10	1.79	2.55
$[CO_2]_{int}$ 1-4 [%]	1.62	1.07	2.27	2.73
$[CO_2]_{int}$ mean 1-4 [%]	1.69	1.07	2.13	2.68
std 1-4 [%]	0.40	0.06	0.34	0.16
$I_p$ [mA]	1.93	2.12	1.91	1.78
$[CO_2]_{int}$ [%] from $I_p$	1.99	0.95	2.10	2.80

in Figure 3.4. Chapter 4 completes the discussion comparing the different possibilities (sensors and models) for estimating the  $[O_2]_{int}$ .

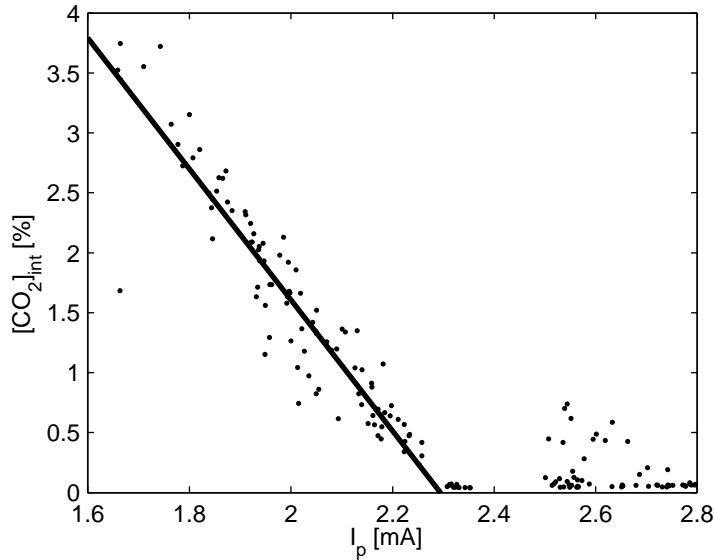
### 3.3 Engine tests

Steady-state and dynamic tests are performed for obtaining the results presented in this work. A summary on these is shown in this section.

#### 3.3.1 Steady-state tests

A full set of steady-state tests are performed in order to characterise the engine and fit the tables and parameters of the models described in Chapter 4. This includes different pairs of  $[n, m_f]$  values covering:

- nominal tests, defined by the manufacturer ECU settings, with warm conditions and variables stabilised,
- variations of the EGR rate (EGR),
- variations of the engine coolant temperature ( $T_{cool}$ ),
- variations of the water temperature of the intercooler ( $T_{wic}$ ) to influence the intake temperature ( $T_{int}$ ), and
- variations of  $p_{boost}$  to influence on  $\dot{m}_{int}$ .



**Figure 3.4.** Equivalency between  $[CO_2]_{int}$  and the pressure corrected signal  $I_p$  from the UEGO sensor. The saturation of the intake probe is evident for values of  $I_p$  around 2.3, i.e. equivalent to a low  $[CO_2]_{int}$  concentration.

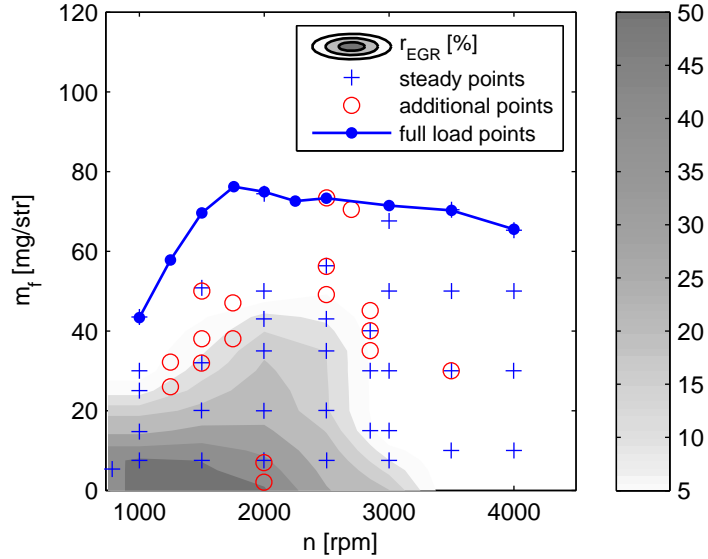
Ranges are selected in order to produce appreciable  $NO_x$  variations but taking into account the limits for engine operation. 363 steady measurements are made in total. Covered engine areas are shown in Figure 3.5.

### 3.3.2 Transient tests

#### 3.3.2.1 Steps

Step-like tests are used for identifying sensor and system dynamics (flow, temperatures and pressures). Table 3.5 shows specific tests performed for different engine operating points.

Furthermore and for demonstrating the effect of sharp variations in SOI on the  $NO_x$  emissions, 19 specific tests consisting in several consecutive steps were performed at different engine speeds (from 1000 to 3000 rpm) with different initial  $NO_x$  concentrations (i.e. different initial engine operating conditions) and step sizes (ranging from  $1^\circ$  to  $4^\circ$  in SOI). Regarding the fuel-to-air ratio, specific load transients are performed by keeping the EGR valve fully closed in



**Figure 3.5.**  $[n, m_f]$  set of values run for testing the engine in steady-state. In each set, different variations of  $u_{egr}$ ,  $T_{cool}$ ,  $p_{boost}$  and  $T_{int}$  are made for characterizing the engine. Some additional points are measured for completing the engine characterization. Nominal EGR map  $EGR_0$  (white area corresponds to EGR valve closed,  $EGR = 0$ ) and full load line are also plotted.

**Table 3.5.** Steps performed in the engine for system and sensors characterisation. BTDC: before top dead centre.

$n$ [rpm]	$m_f$ [mg/str]	$\Delta u_{soi}$ [ $^\circ$ BTDC]	$\Delta u_{egr}$ [%]	$\Delta u_{wg}$ [%]
1000	7.5	$\pm 2$	$\pm 10$	-30
1000	43	$\pm 2$		-3
2500	35	$\pm 2$	$\pm 5$	$\pm 3$
4000	10	$\pm 2$		$\pm 5$
4000	50	$\pm 2$		$\pm 3$

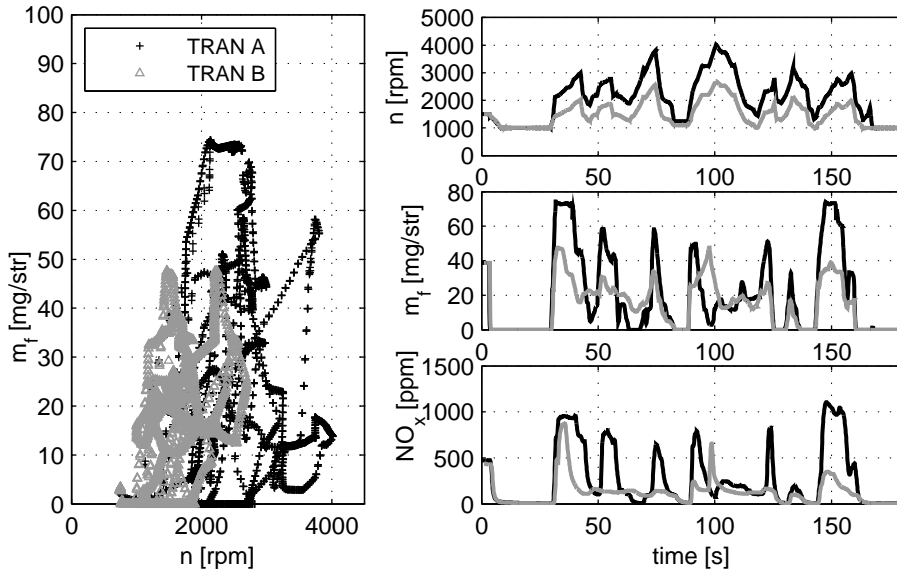
order to avoid possible interactions. Data is later used for the sensor dynamics identification in Section 3.4.

### 3.3.2.2 Cycles

Different engine cycles are designed on the basis of the homologation cycle and *ad-hoc* modifications for proving and testing the models and methods

developed in this work. In the following, the cycles are described and the test campaign is summed up in Table 3.7.

**Adapted transient cycles (TRAN).** A set of dynamic tests are designed on the basis of two different modifications of the FTP cycle for HD engines (named here as TRAN A and TRAN B), and are used for training and evaluating a  $\text{NO}_x$  model in transient operation. These consist of a re-escalation of the central part of the standard FTP and are compared in Figure 3.6. Test proposition is a full factorial design by performing modifications on set-points for  $u_{egr}$  (by modification on the  $\dot{m}_a$  maps),  $p_{boost}$  (by fixing  $\dot{m}_a$  set-point and modifying  $p_{boost}$  map) and  $T_{cool}$  (modifying the engine coolant reference temperature at the test cell controller). The sequence is presented in Table 3.6, totalling 36 different tests (27 TRAN A and 9 TRAN B, both in the test campaign A). For each test, three repetitions of the sequence shown in Figure 3.6 are made.



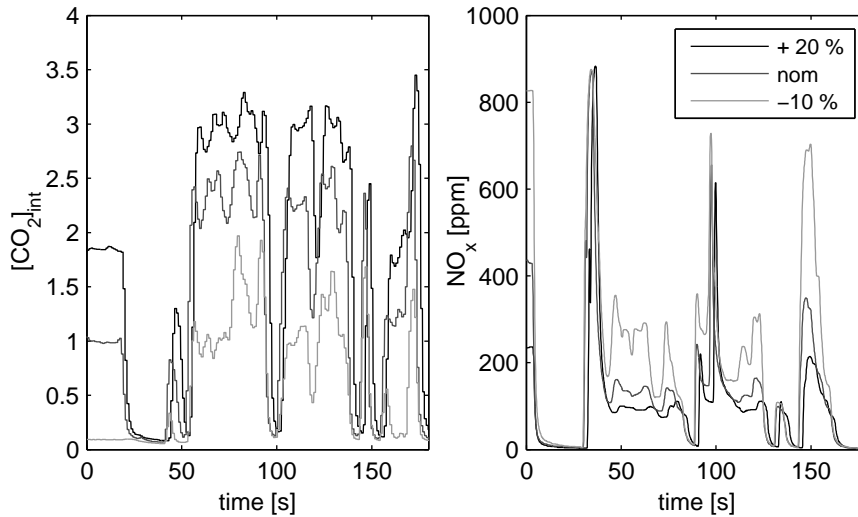
**Figure 3.6.** TRAN nominal cycles designed for the model validation. TRAN A is the black line, while TRAN B is the grey line.

The  $\text{NO}_x$  sensitivity to variations on inputs and parameters is considerably lower for transient than static conditions. This fact is due to the systems and sensors dynamics which filter the effects. For instance, an increasing of 10%

**Table 3.6.** Experimental plan of the modified TRAN cycles.

cycle	$T_{cool}$ [°C]	$\Delta u_{egr}$ [%]	$\Delta p_{boost}$ [mbar]
TRAN A	[75,85,95]	[-5,0,+5]	[-100,0,+100]
TRAN B	nom	[-10,0,+20]	[-100,0,+100]

in  $p_{boost}$  (100 mbar) in the TRAN A cycle leads to an increase  $NO_x$  around 10%, while in steady conditions produces an increase in the order of 40% for the same conditions (while fixing  $\dot{m}_a$  set-point in the EGR area of the engine). In all cases, the higher  $NO_x$  sensitivity is found for EGR variations, especially for TRAN B cycles, which present sharper EGR transients. Figure 3.7 shows the effect of modifying EGR rate (by plotting  $[CO_2]_{int}$ ) over the final  $NO_x$  emissions measured by the sensor.

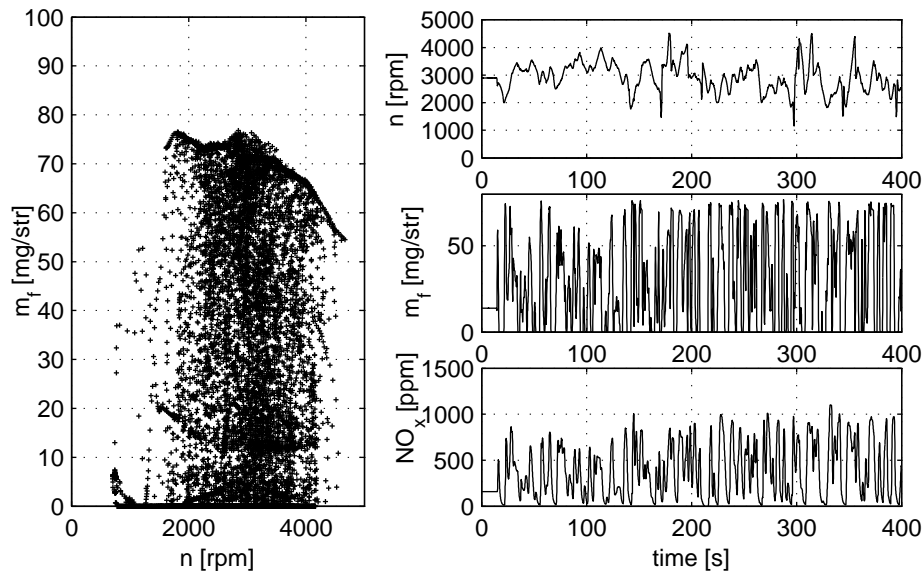


**Figure 3.7.**  $[CO_2]_{int}$  at intake and  $NO_x$  emissions for TRAN B after varying the EGR set-point. The legend in the right plot shows the values of  $u_{egr}$ . Note that the higher  $u_{egr}$  the lower  $\dot{m}_{egr}$ .

**Sportive driving mountain profile (SDMP).** The SDMP is designed for covering a wide operating range of the engine by sharp variations on  $n$  and  $m_f$ . A zoom on 400 seconds (total is 1200) of the SDMP is shown in Figure 3.8 (left plot indicates the whole cycle points). Since the SDMP presents sharp load transients, the EGR valve is mostly closed during the cycle, making that

$EGR = 0$  during the major part in spite of the engine might be run in the EGR area. This circumstance distorts the closed loop nominal engine operation.

**New European driving cycle (NEDC).** The NEDC is tested in cold and warm-start conditions (at the beginning of the test). The cycle consists of a urban part with accelerations/decelerations and idling, and an extra-urban part that lasts 400 s. The urban part has 4 equal repetitions of a sequence that lasts 200 s; Figure 3.9 shows last repetition of the urban part and the extra-urban section. Cold-start cycle presents lower  $NO_x$ , indeed when  $m_f$  is slightly higher because of cold starting strategies. Anyway, differences are not considerable and can also be associated with cycle differences in operating points ( $n$  and  $m_f$  have dispersion, see left plot of Figure 3.9).



*Figure 3.8. SDMP cycle performed on the engine.*

**Common Artemis driving cycles (CADC).** CADC cycles are designed upon the basis of an statistical analysis of European real world driving patterns, developed by the European Artemis<sup>1</sup> project. The cycle includes three

<sup>1</sup>ARTEMIS: Assessment and Reliability of Transport Emission Models and Inventory Systems.

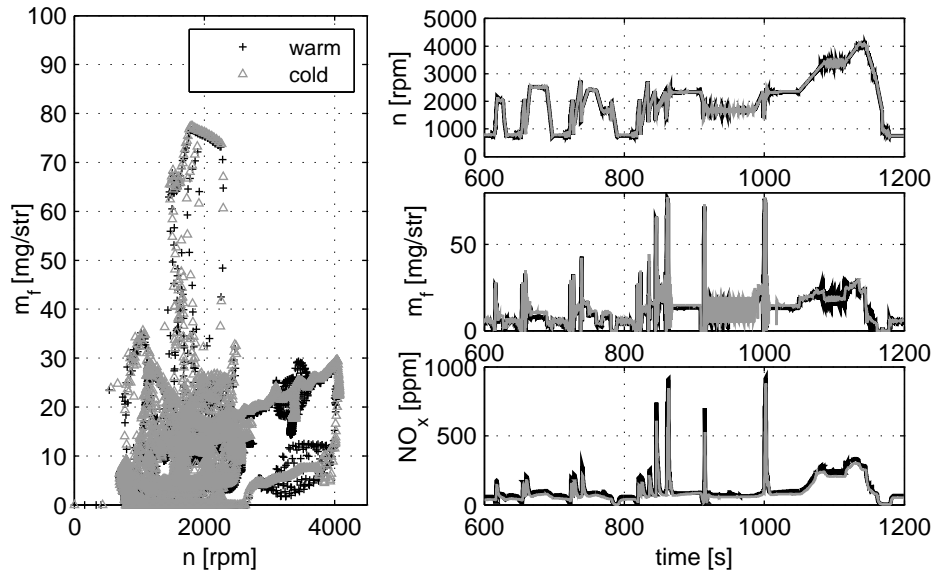


Figure 3.9. NEDC cycle performed on the engine.

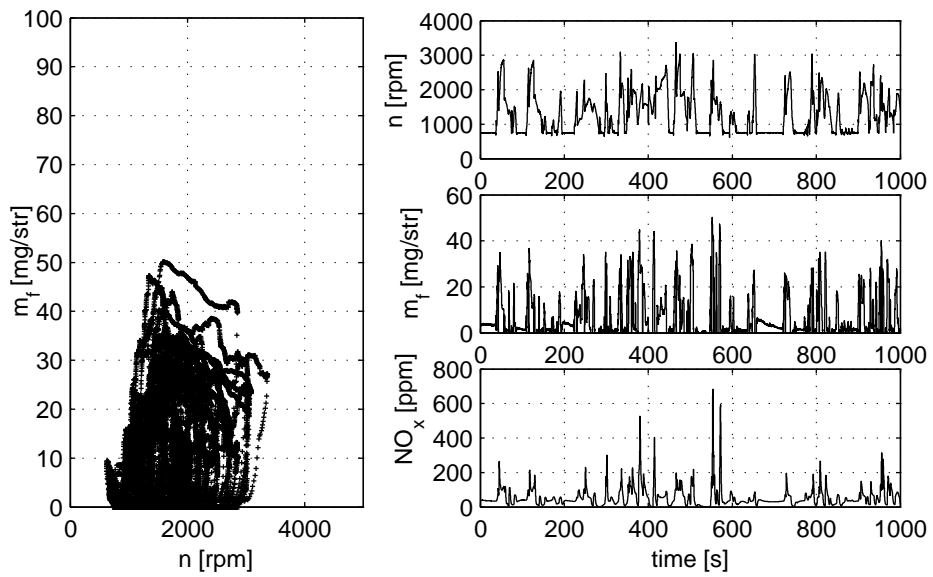


Figure 3.10. Urban CADC cycle performed on the engine.

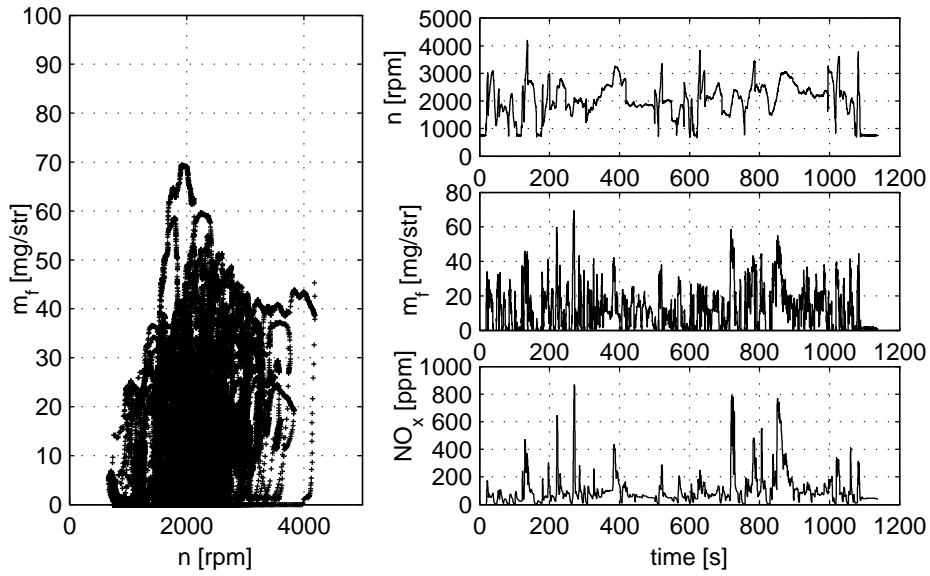


Figure 3.11. Rural CADC cycle performed on the engine.

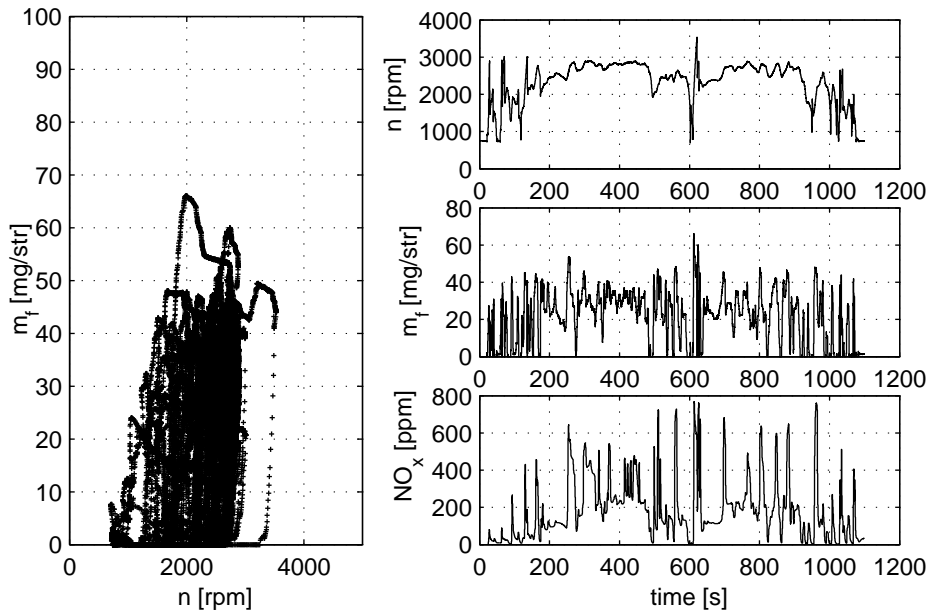


Figure 3.12. Highway CADC cycle performed on the engine.



different variants: urban, rural road and highway. These profiles are more realistic than the NEDC as they cover different real situations and then constitute a good source for experimental testing. Three versions of the CADC are shown in Figures 3.10 to 3.12.

**Summary on the transient cycles.** In total, two different test campaigns for testing dynamic cycles have been performed for the present work; these are separated in test campaign A and B just for remarking that they have been performed in different moments. Table 3.7 sums up the dynamic tests made, where an identifier (ID) names the cycles. A3 to A29, and A30 to A38 stand for the test of the full factorial design summarised in Table 3.6.

*Table 3.7. Dynamic test campaigns.*

ID	Cycle	Calibration
Test campaign A		
A1	NEDC warm	nom
A2	NEDC cold	nom
A3-A29	TRAN A	DoE
A30-A38	TRAN B	DoE
Test campaign B		
B1	NEDC warm	nom
B2	CADC Urban	nom
B3	CADC Rural	nom
B4	CADC Highway	nom
B5	SDMP 1	nom
B6	SDMP 2	nom

### 3.4 Gas concentration sensors characterisation

In the next, the static and dynamic calibration of  $\lambda$  and  $\text{NO}_x$  output from the gas concentration sensors is discussed. The static calibration is based on the linear fitting of sensor outputs by using the steady-state data-set collected from the engine, while the dynamic characterisation is based on generating exhaust gases with different concentrations by means of the step-like variations on the engine injection settings.

With respect to the  $\text{NO}_x$  output, this section presents an easy on-engine experiment for  $\text{NO}_x$  sensors by abrupt changes on the start of injection (SOI).

The results shown in this section corresponds to a  $\text{NO}_x$  sensor and a gas probe located downstream of the turbine, removing the AT line.

Regarding the  $\lambda^{-1}$  output characterisation, variations on the injected fuel mass rate  $\dot{m}_f$  are used for provoking step-like transitions on the  $\lambda$  output and the methodology for identifying the sensor is similar to that of the  $\text{NO}_x$  output. Figure 2.5 already compared  $\lambda^{-1}$  output from the  $\text{NO}_x$  and a UEGO sensor in a configuration different from the standard of this work, where the former is located downstream of the ATs while the latter at the turbine outlet. The  $\text{NO}_x$  sensor provides a slower and more filtered response than the other due to the position and the sensor response time. However, if using a  $\text{NO}_x$  sensor, the UEGO sensor can be removed instead of using two different sensor units. In the following, the results for the  $\text{NO}_x$  sensor output located at the turbine outlet are presented<sup>2</sup>. The conclusions and results of the characterisation are required for the implementation of models and observers in subsequent chapters.

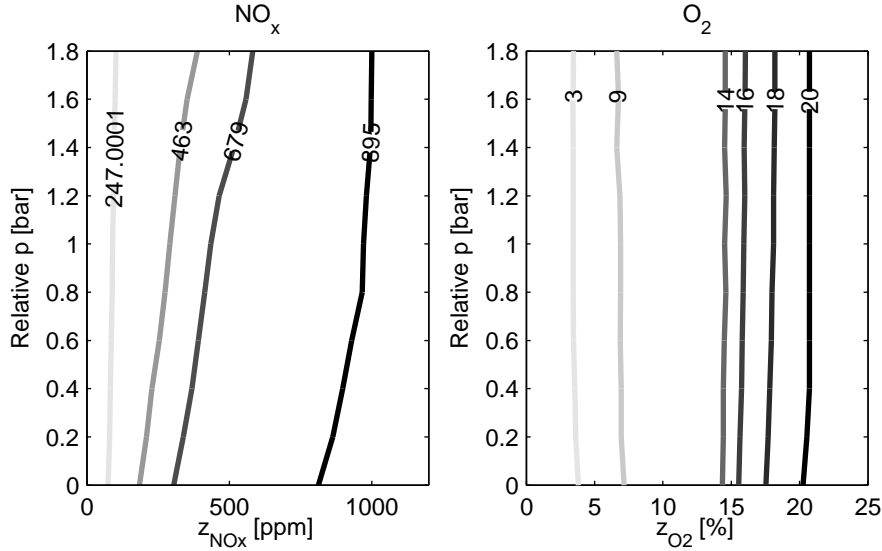
### 3.4.1 Static calibration

Previous to the  $\text{NO}_x$  sensor static calibration, some issues must be considered. On one hand, the measurement procedures of the sensors affect the output itself, especially for  $\text{NO}_x$  where a number of different species (NO and  $\text{NO}_2$  among others) can be distinguished. This discussion is moved to the Appendix 3.B.

On the other hand, the  $\text{NO}_x$  sensor principle is based on the measurement of the partial pressure of oxygen in order to determine both  $\lambda^{-1}$  ( $z_{\lambda^{-1}}$ ) in a first chamber (outer cavity) and  $\text{NO}_x$  ( $z_{\text{NO}_x}$ ) in a second one (inner cavity), as shown in Figure 2.4. The total pressure of the manifolds affects the output, as occurs also on UEGO and MAF sensors, and a correction (by software) is required. A simple analysis by disposing different gas concentrations and pressures (in a static way) in a sealed chamber allows checking the pressure effect over the  $\text{NO}_x$  sensor outputs. Two gas bottles containing pure  $\text{CO}_2$  and a calibrated mix of  $\text{NO}_x$  and  $\text{N}_2$  are used for designing the mixes when added to atmospheric air. A  $\text{ZrO}_2$  sensor is installed and measurements are made in steady-state.

Figure 3.13 shows iso- $\text{NO}_x$  curves (left plot) and iso- $\text{O}_2$  (right plot) for different relative pressures in the chamber and for different  $\text{NO}_x$  levels (ppm) and oxygen levels (%), tested in the sealed chamber with the actual references known beforehand. The curves are obtained discharging the vessel, which

<sup>2</sup>The UEGO sensor is used for initial phases of this work but then removed.



**Figure 3.13.** Effect of the pressure over the NO<sub>x</sub> sensor for the NO<sub>x</sub> output  $z_{NO_x}$  (left) and oxygen estimation  $z_{O_2}$  from  $\lambda^{-1}$  output (right). It seems that the manufacturer has compensated the oxygen output but not the NO<sub>x</sub> output.

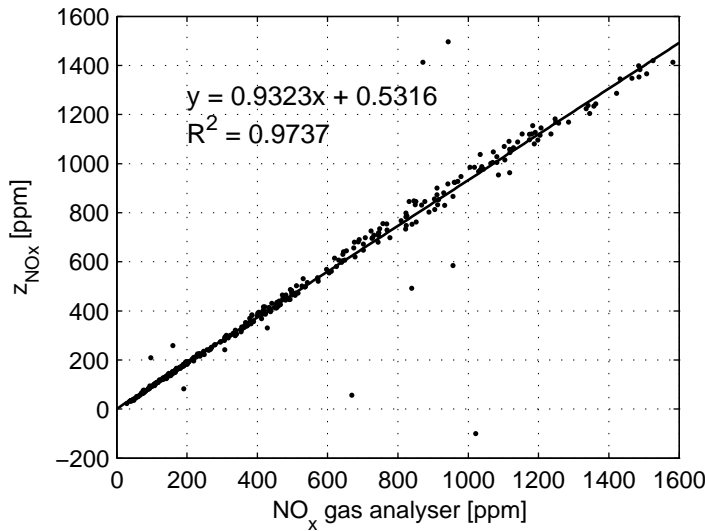
ensures that the concentration is constant along each one of the tests. As it can be noticed in the plot, the oxygen output ( $z_{O_2}$ ) is already pressure compensated by the manufacturer, while the NO<sub>x</sub> output is affected and a specific correction seems necessary if the sensor is to be used with varying gas pressures.

The NO<sub>x</sub> sensor is installed at the turbine outlet. The turbine outlet pressure ( $p_{out}$ ) is in all cases between 1 and 1.6 bar (in the NEDC the maximum  $p_{out}$  is about 1.3 bar). Derived from the limited range in  $p_{out}$ , the possible error due to the pressure effect is lower than the expected error according to the specifications provided by the manufacturer (shown in Appendix 3.A); hence no specific correction has been used in the present work.

### 3.4.1.1 Static calibration of the NO<sub>x</sub> output

With respect to the NO<sub>x</sub> sensor accuracy, the ZrO<sub>2</sub> sensor is calibrated by comparing the signal with the NO<sub>x</sub> output of the gas analyser, used as static standard. The NO<sub>x</sub> output in the gas analyser is measured by a chemiluminescence system (CLD) in a dry basis and corrected by humidity and other

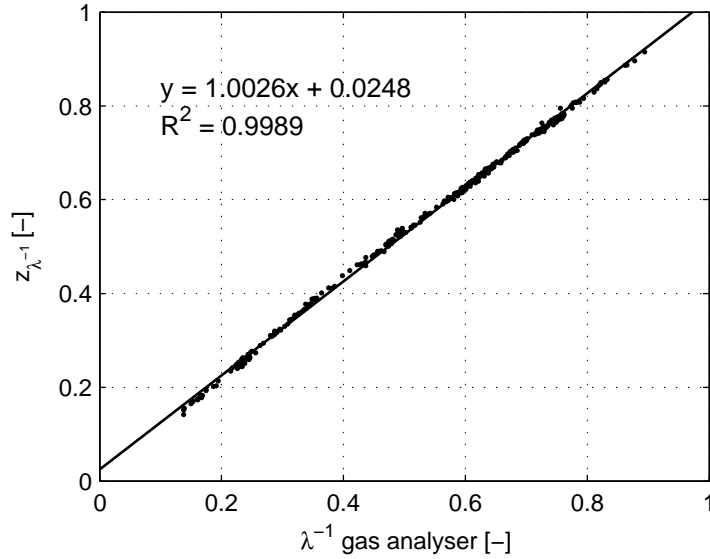
factors. The important dependability of the exhaust  $\text{NO}_x$  concentration when varying the engine operating conditions is profited for the steady calibration. The steady-state tests are used for the calibration, causing  $\text{NO}_x$  concentration variations to cover the sensor measuring range. Figure 3.14 illustrates the results. After removing the outliers (usually caused by human errors in the measurement procedure of the gas analyser) and using the box-and-whiskers method, a linear fit is performed. Mean absolute error after calibrating  $\text{NO}_x$  from  $\text{ZrO}_2$  output results of 10.7 ppm, and the residuum standard deviation of 14.77 ppm.



**Figure 3.14.**  $\text{NO}_x$  from  $\text{ZrO}_2$  sensor ( $z_{\text{NO}_x}$ ) and the gas analyser under different steady operating conditions; the linear fit and correlation coefficient are represented.

### 3.4.1.2 Static calibration of the $\lambda^{-1}$ output

$\text{NO}_x$  sensor provides a very accurate information of  $\lambda^{-1}$  in steady-state operation. Static measurements of the  $\text{NO}_x$  sensor are compared to those of the gas analyser in Figure 3.15 by using the steady-state data-set. Mean absolute error after calibrating  $\lambda^{-1}$  from  $\text{ZrO}_2$  output results of 0.0048, and the residuum standard deviation of 0.004.



**Figure 3.15.**  $\lambda^{-1}$  from  $ZrO_2$  sensor ( $z_{\lambda^{-1}}$ ) and the gas analyser under different steady operating conditions; the linear fit and correlation coefficient are represented.

### 3.4.2 Dynamic calibration

The second step in the calibration procedure consists in the identification of the sensor dynamics. A well-known method consists of using a step transition in the measured quantity [10]. However, usual means for providing step-like transition in gas concentration are complex set-ups with fast valves and synthetic gas mixes [11], which can hardly be implemented on-board.

For solving this issue, on-engine alternatives are next presented for the calibration of the  $NO_x$  and  $\lambda^{-1}$  outputs of the exhaust concentration sensor.

#### 3.4.2.1 Dynamic calibration of the $NO_x$ output

This subsection presents a novel method based on the variation of the SOI, which can be easily done in current engines during their normal operation. In order to decouple as much as possible the effect of SOI variations on the air path of the engine, EGR rate is set to zero (EGR valve fully closed) during the tests. Although other variables could be used for affecting the  $NO_x$  concentration (e.g. EGR valve control), SOI is selected because:

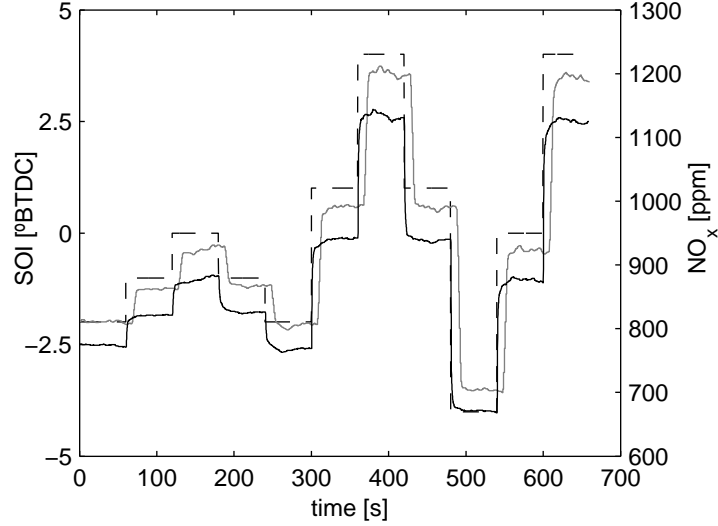
- SOI varies the exhaust gas composition in an important way, producing appreciable  $\text{NO}_x$  concentration variations.
- SOI variations produce low variations for turbine intake temperature and pressure. Hence conditions at the turbine inlet are not importantly varied, and the turbocharged speed and boost pressure are not significantly affected; since the EGR valve is closed, the effect of the slight variation in the exhaust pressure may be neglected. This is considered as a key factor since pressure and temperature variations are associated with slow transients (because of mass and heat accumulation) which would affect the  $\text{NO}_x$  concentration.
- Other control variables correspond to slow acting valves, which will distort the step-like profile in the gas composition. However, the system response to SOI variation can be considered instantaneous: SOI is electronically applied, thus no actuation delay is expected beyond the cycle-to-cycle response time<sup>3</sup>. Additionally, since no EGR is performed,  $\text{NO}_x$  concentration at the cylinder exhaust port is only dependent on the trapped air mass quantity and temperature, and injection settings during the previous cycle ( $\text{NO}_x$  contained in the residual gas fraction is re-burned during the combustion [12]). Finally, the gas transport delay from the cylinder to the  $\text{ZrO}_2$  sensor (located at least 1 m from the exhaust port) is of a few tens of ms, as it will be discussed in Section 3.4.2.3.

Figure 3.16 shows the evolution of  $\text{NO}_x$  measurements during a cycle with SOI steps, while Figure 3.17 illustrates that other significant variables (as intake and exhaust pressure, gas temperatures, coolant temperature and air mass flow) do not vary in an important way during the tests. The maximum variation in these properties is 2.8% for exhaust pressure (considering a moving average filtered value), while the rest of them varies less than 0.5%.

**Sensor identification.** For illustrating the applicability of using data derived from the SOI steps, a dynamic identification is done assuming a linear first order model behaviour (other possible model structures with physical insight could be found in [13]) with the transfer equation in the s-domain

$$G(s) = \frac{e^{-s\tau_{\text{NO}_x}}}{1 + T_{\text{NO}_x}s} \quad (3.1)$$

<sup>3</sup>This delay is in the order of 20 ms maximum for an engine spinning at 3000 rpm, which can be neglected when comparing with typical delays of the sensor, which are above 500 ms. On the other hand, maximum software delay is below 40 ms, which is not additive to the previous one.



**Figure 3.16.**  $NO_x$  concentration during a SOI step-like test at 2000 rpm and 50 mg/stroke. Legend: --- is SOI, — is the  $NO_x$  sensor ( $z_{NO_x}$ ) before the steady-state calibration and — is the gas analyser. The difference in terms of delay, dynamics and steady-state values between both sensors is clearly noticed in the Figure.

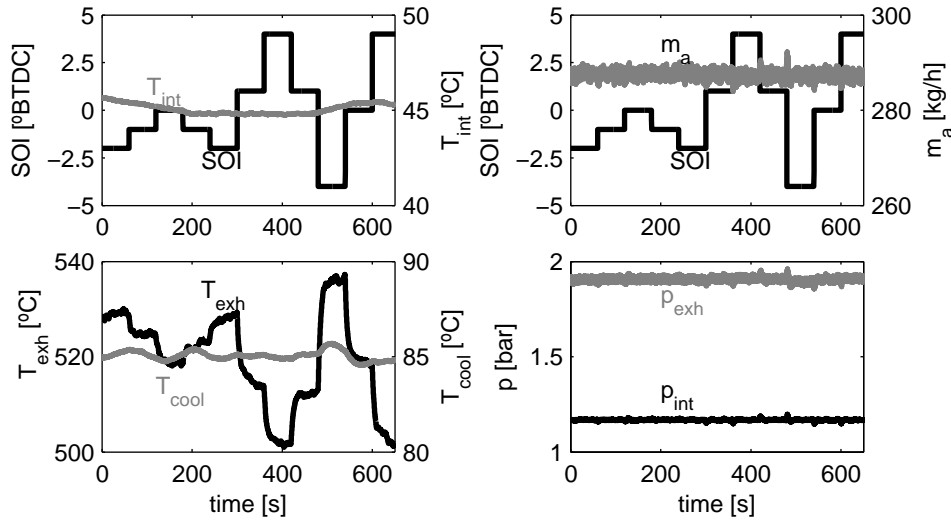
The sensor response to the step is then

$$z_{NO_x}(t) = z_{NO_x}(t_0) + \Delta NO_x \left( 1 - e^{-\frac{t - \tau_{NO_x}}{T_{NO_x}}} \right) \quad t \geq \tau_{NO_x} \quad (3.2a)$$

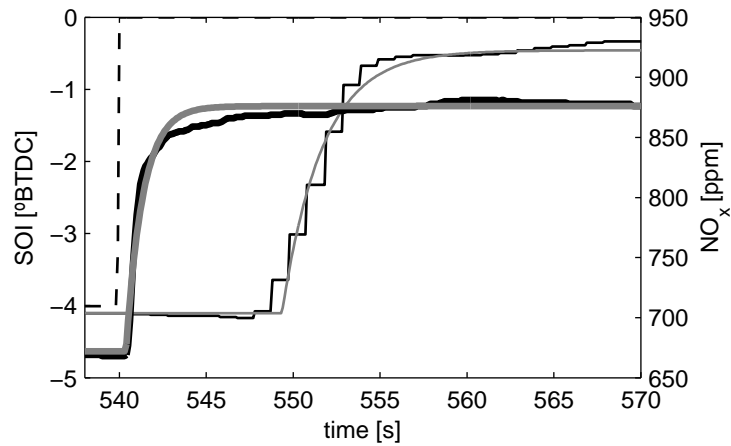
$$z_{NO_x}(t) = z_{NO_x}(t_0) \quad t < \tau_{NO_x} \quad (3.2b)$$

where  $z_{NO_x}(t_0)$  is the  $z_{NO_x}$  initial value before the step,  $\Delta NO_x$  is the step-amplitude and  $t$  the time. The sensor delay ( $\tau_{NO_x}$ ) represents the delay from the application of SOI (judged to be instantaneous) to the start of the sensor output variation, and  $T_{NO_x}$  stands for the sensor time constant. For each individual step and sensor,  $T_{NO_x}$  and  $\tau_{NO_x}$  are obtained. For the coefficient identification, a least-square fit of (3.2) is performed (although more robust recursive identification techniques could be used for the online case [14]). Even though not shown in this work, other different models have been proved, as for instance second order filters and AR or ARMAX models.

Figure 3.18 illustrates an example of the fit obtained with the first order response model. Averaged values of the identified time constant  $T_{NO_x}$  and their deviations are shown in Table 3.8. On the other hand, the standard



**Figure 3.17.** Evolution of different engine variables during a SOI step-like test at 2000 rpm and 50 mg/str.



**Figure 3.18.** Identified models (solid gray) and measured evolution (solid black) for both sensors: thin lines correspond to the gas analyser while thick ones to the  $\text{NO}_x$  sensor. SOI step is reproduced with a black dashed line. The delay of the gas analyser is clear with respect to  $\text{NO}_x$  sensor.

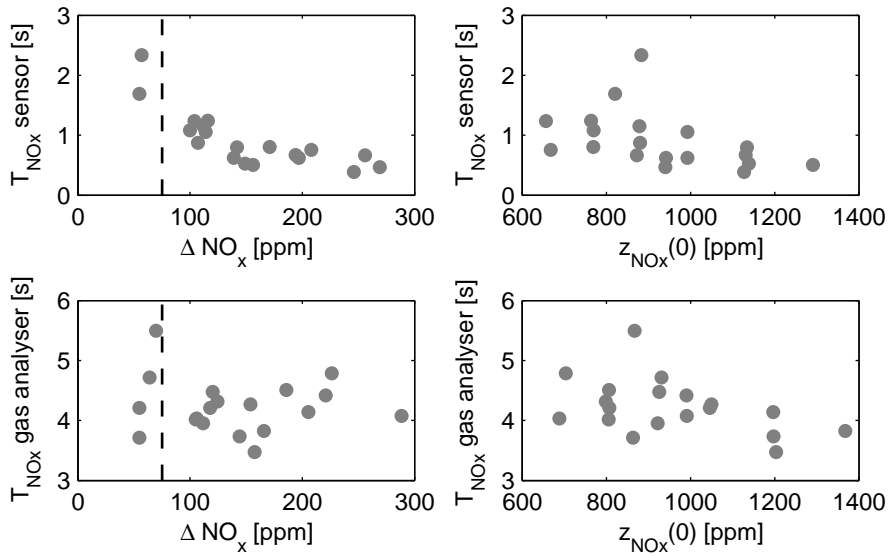


deviations obtained for  $T_{\text{NO}_x}$  estimates merit for additional analysis, as they suggest that the sensor behaviour could be affected by the gas concentration conditions during the test.

**Table 3.8.** Mean and standard deviations for the identified time constant  $T_{\text{NO}_x}$ .

	Gas analyser	NO <sub>x</sub> sensor
$T_{\text{NO}_x}$ [s]	4.25	0.91
$\sigma_k$ [s]	0.46	0.47

No direct influence is found on the sensors response from the gas speed, temperature or pressure, since the identified parameters do not vary significantly with these conditions. In order to analyse the effect of the initial  $z_{\text{NO}_x}$  concentration ( $z_{\text{NO}_x}(0)$ ) and of the step size ( $\Delta\text{NO}_x$ ), scatter plots of  $T_{\text{NO}_x}$  are presented in Figure 3.19.



**Figure 3.19.** Response time  $T_{\text{NO}_x}$  for the NO<sub>x</sub> sensor as function of the concentration step  $\Delta\text{NO}_x$  and the initial concentration  $z_{\text{NO}_x}(0)$ .

In the case of the ZrO<sub>2</sub> sensor, there exists an important variability of  $T_{\text{NO}_x}$ , especially at low  $\Delta\text{NO}_x$ , with an exponential-like shape. One reason is that, in the case of small steps, the concentration profile is not step-like, but affected

by other slower variations (temperature, pressure, etc.) that here are not negligible; another possible reason could be related to the diffusion speed of the gas inside the sensor chambers, which depends on the concentration itself and its evolution along time. Note also that the sensor presents a low sensitivity to low  $\text{NO}_x$  values [15]. This explains the larger deviations obtained in  $T_{\text{NO}_x}$  as presented in Table 3.8. However, the response time of the gas analyser does not present these problems and the trend is more constant (and response quite slower) for all values (see bottom plots in Figure 3.19). For larger step sizes, the response time of the  $\text{NO}_x$  sensor is more stable, as shown in Table 3.9 where tests with  $\Delta\text{NO}_x < 75$  ppm have been removed. Physical modelling approaches can be useful for identifying these behaviours by including dynamic models for diffusion chambers [16].

**Table 3.9.** Identified dynamics for  $\text{NO}_x$  sensors considering tests with  $\Delta\text{NO}_x > 75$  ppm.

	Gas analyser	$\text{NO}_x$ sensor
$T_{\text{NO}_x}$ [s]	4.14	0.79
$\sigma_k$ [s]	0.33	0.27

Furthermore, only low  $\Delta\text{NO}_x$  could be applied in tests with low  $z_{\text{NO}_x}(0)$ , because SOI actuation does not permit significant variations in the gas concentration. Hence  $z_{\text{NO}_x}(0)$  and  $\Delta\text{NO}_x$  are correlated. This partly explains the effect of high deviation presented for low  $z_{\text{NO}_x}(0)$  values shown in right-hand plots of Figure 3.19. In addition, the low sensitivity of  $\text{ZrO}_2$  sensors to low  $\text{NO}_x$  values can be addressed [17].

The discussion about the sensor delay is included in a different subsection. Finally, one of the main problems of automotive sensors is their degradation during their lifespan. For ensuring the correct performance of the control structure, the system has to be capable of checking the sensor periodically. This methodology could be useful for these purposes.

**Correspondence with the discrete filter.** The results above have been presented in the s-domain and the continuous time domain, due to the easy interpretation and for supporting the discussion. However, electronic systems work on the discrete domain. An alternative for converting the obtained results to the z-domain is using the first order discrete filter to model the sensor response

$$G(z) = \frac{1 - a_{\text{NO}_x}}{1 - a_{\text{NO}_x}z^{-1}}z^{-\tau_{\text{NO}_x}/T_s} \quad (3.3)$$

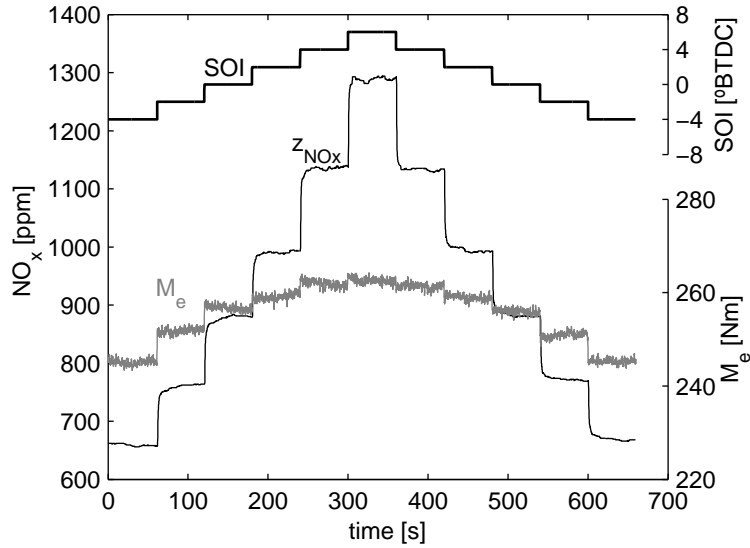
instead of the continuous filter where  $z$  is the  $z$ -transform variable,  $a_{\text{NO}_x}$  represents the response time and  $\tau_{\text{NO}_x}$  is the delay. Depending on the sample time ( $T_s$ ),  $a_{\text{NO}_x}$  can be deduced [10].

**Prospective for the on-board application.** Presented method based on shifting SOI could be easily applied during operation phase of automotive diesel engines. That means that the ECU could use this test for identifying the sensor behaviour, and diagnosing the sensor and system performance. This could be periodically done during specific conditions of the engine run. The knowledge of the dynamic sensor characteristics is needed for data fusion and signal reconstruction techniques [18, 19]. For this, the dispersion of the sensor model might be corrected online using the combination of the mentioned techniques and the sensor measurements by using the proposed method. In this way, the robustness of the control law can be checked regarding this dispersion.

Note that the presented results correspond to a  $\text{NO}_x$  sensor located upstream of the AT devices. Hence the presented methodology will be of direct application for sensors in this position, which can be used for feed-forward control of the SCR system [20]. For the case of the sensor located downstream of a SCR system, the following two aspects must be considered:

- The method can be applied but the model should take into account the effect of the SCR. On one hand, the important removal of the raw  $\text{NO}_x$  concentration (about 80%) compromises the  $\text{NO}_x$  output due to the insufficient accuracy of the sensor for low  $\text{NO}_x$  values. On the other hand, the cross sensitivity to ammonia affect the sensor measurement. Because of these, the dynamic characteristics of the signal can be importantly distorted (also filtering the signal and thus losing the step-like behaviour).
- Even in the case the SCR is not operating, the possibility of exciting the  $\text{NO}_x$  sensor needs to be checked: depending on the signal dynamics and the volume of AT devices, the gas concentration can sufficiently keep its step-like profile. This issue must be verified beforehand by installing two sensors, one up and one downstream of the AT line. Anyway, once established the feasibility of the method, one of them may be removed. Finally, the  $\text{NH}_3$  storage mechanism could also contribute to unexpected variations in the  $\text{NO}_x$  concentration at the SCR outlet, and hence a sufficient delay is needed between the SCR cut-off and the dynamic characterisation checking.

However, the method could still be used for the SCR system diagnosing, or a proper dynamic model as e.g. the one presented in [21] might be used for covering the storage dynamics and  $\text{NH}_3$  effects.

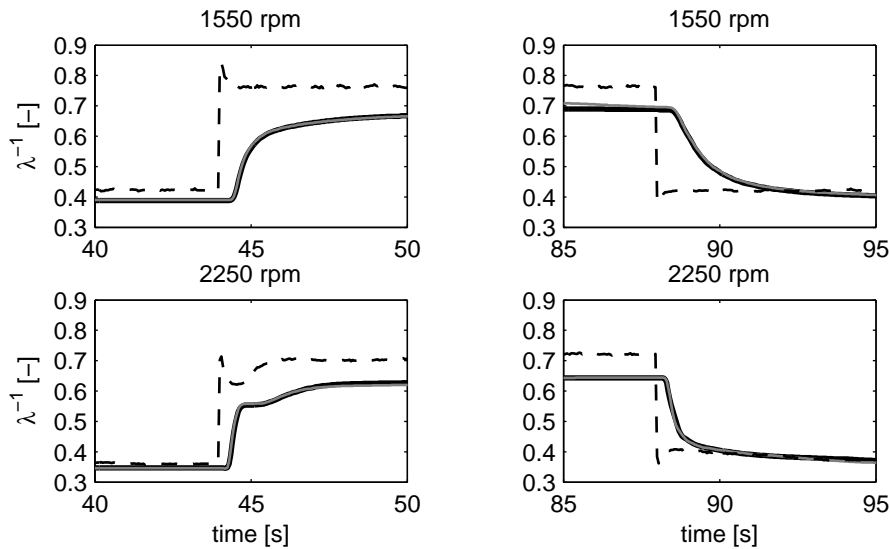


**Figure 3.20.**  $\text{NO}_x$  concentration measured by the  $\text{NO}_x$  sensor ( $z_{\text{NO}_x}$ ) and torque evolution ( $M_e$ ) at 2000 rpm and 50 mg/stroke when varying SOI.

Excellent operating condition for performing the method is highway operation (where torque and engine speed are reasonably constant during long periods and SOI variations can produce appreciate variations). Figure 3.20 shows the variation in  $\text{NO}_x$  concentration and torque when SOI is varied in a 2-degree step profile while keeping constant  $m_f$  and  $n$ . As shown in the central part of the plot, the raw  $\text{NO}_x$  production can clearly be increased with a low impact on torque; that means that the driveability can be ensured while the test is performed (corrections on  $m_f$  can be performed for maintaining the engine torque to its set-point ensuring the driver to be blind to the test). The increase in  $\text{NO}_x$  emission during the sensor testing should not importantly impact the overall emissions, since a few seconds are enough for the test (and also a descendent step could be used). On the other hand, in engines with SCR technology and a  $\text{NO}_x$  sensor located upstream of the SCR, the SCR could cope with the excess of  $\text{NO}_x$ .

### 3.4.2.2 Dynamic calibration of the $\lambda^{-1}$ output

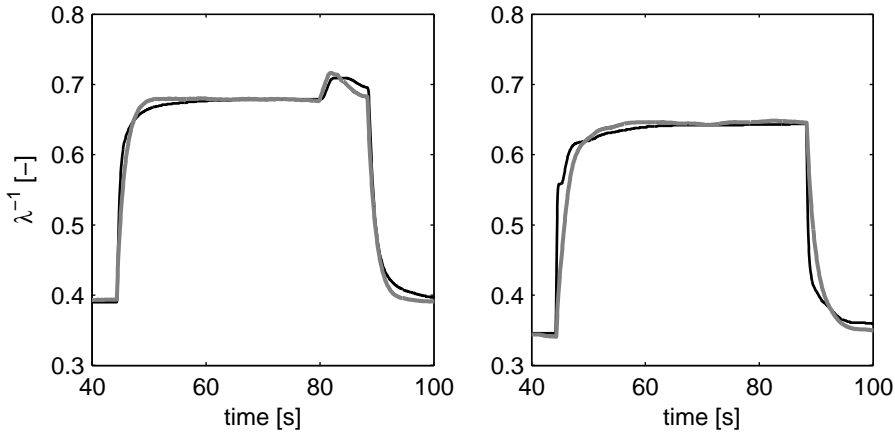
The slow dynamic response of  $\lambda^{-1}$  from the  $\text{NO}_x$  sensor is explained by several reasons: the distance and the system volume between engine exhaust and sensor location affect transport delay and distort the signal; the sensor measurement principle and sensor hardware itself; and the acquisition chain. Although slow, the sensor is quite precise, since several repetitions of the same test provide similar results. This is illustrated in Figure 3.21, where 2 repetitions of the same fuel step are depicted for 2 different engine speeds. Top plots show a step up (left) and step down (right) in injection for 1550 rpm and bottom plots the same for 2250 rpm. Despite the sensor precision, it must be considered that the dynamic response is significantly affected by the operative conditions (e.g. the sensor response in bottom plots in Figure 3.21 is always faster than in top plots). A model based on the fuel-to-air ratio calculation ( $x_{\lambda^{-1}}$ ) is also shown in the plots (detailed explanation is found in the Chapter 4).



**Figure 3.21.** Two sensor responses ( $z_{\lambda^{-1}}$  in solid lines) for injection steps from 15 mg/str to 30 mg/str (left plots) and the opposite (right plots) at 1550 rpm (top plots) and 2250 rpm (bottom plots). The model  $x_{\lambda^{-1}}$  is provided for comparison (dashed black).

A first order filter similar to that of (3.3) is used and parameters identified on the basis of the data obtained by performing steps on  $m_f$ , which provoke step-like transitions on  $\lambda^{-1}$ . Figure 3.22 shows the evolution of the measured

$z_{\lambda^{-1}}$  and the identified sensor model  $y_{\lambda^{-1}}$  for two different operating conditions. Left plot is used for identification while validation is provided by the right plot. The fitting is good (even in the right plot) although a slight error can be appreciated, which suggests that the dependence of the sensor dynamics with the engine operating point may be neglected in a first approach, i.e. the response time may be taken as constant. A similar study to that of the  $\text{NO}_x$  output may be performed in order to obtain characteristic values of the response time. Anyway, the repetitiveness of the parameters is higher than for the  $\text{NO}_x$  case, and in this work the  $\text{NO}_x$  sensor response time for  $\lambda^{-1}$  output is considered constant.



**Figure 3.22.** Example of identification of the sensor behaviour for two engine speeds, where  $z_{\lambda^{-1}}$  and  $y_{\lambda^{-1}}$  are depicted. Top plot: training data at 1550 rpm where parameters  $a_{\lambda^{-1}}$  and  $\tau_{\lambda^{-1}}$  are identified. Bottom plot: validation data at 2250 rpm, where parameters  $a_{\lambda^{-1}}$  and  $\tau_{\lambda^{-1}}$  obtained for the top plot situation are used now for obtaining  $y_{\lambda^{-1}}$ .

This method could be applied when driver performs a sharp tip-on or can profit DPF regeneration modes, even though the effect of post-injections should be studied more deeply.

### 3.4.2.3 Delay in the $\text{NO}_x$ sensor

As expected, the  $\text{NO}_x$  sensor delay is much lower than chemiluminescence system delay. This occurs because the gas analyser delay is mainly caused by the treatment operations and by the transport in the long gas line feeding the measurement system (12 m), which adds a non-negligible delay due to

the low flow allowed. The outputs have been plotted in Figure 3.16, and the significant lower delay for the ZrO<sub>2</sub> sensors with respect to the gas analyser can be clearly noticed.

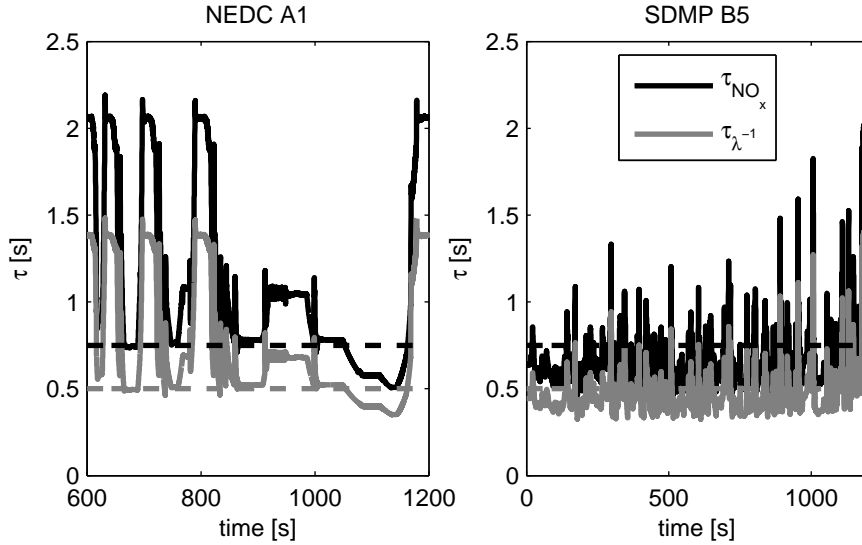
In principle, the ZrO<sub>2</sub> sensor delay (for both  $\lambda^{-1}$  and NO<sub>x</sub> outputs) varies with the sensor actuation possibly by gas concentrations [16] and in a minor way by transport lag. The effect of the gas concentrations and mass flows themselves over the final delay is appreciable and should be analysed but it escapes from the scope of this work. With respect to the transport delay, it could be calculated taken into account the gas flow rate and the exhaust pipe cross section. The transport delay can take a wide range of values depending on the operating conditions. Simulations with a wave action model [22] and for the considered diesel engine after varying injected fuel mass and engine speed show velocities of the exhaust flow between 10 m/s and 140 m/s. The transport delay could be then calculated as the time that the gas takes to reach the sensor from the combustion chamber, and goes with the inverse of the gas velocity. From the simulations, transport delays rarely present values above 100 ms, and according to simulations by wave action models, are generally lower than 25 ms. The influence on the total delay is supposed to be minimum with respect to the hardware delay.

Some authors have proposed functions for modelling the delay. For instance, Trimboli *et al.* [23] design the following function

$$\tau(n, m_a) = \alpha_0 + \frac{\alpha_1}{n} + \frac{\alpha_2}{m_a} \quad (3.4)$$

Figure 3.23 shows the delay modelled by (3.4) in the cycles A1 and B5 for both NO<sub>x</sub> and  $\lambda^{-1}$  outputs. The results show that the delay is quite constant during part and full load operation, while it is slightly bigger when approaching to idling or low loads. The variation can be explained not really by the (minimum) transport lag but mainly by the sensor diffusion processes between the sensor chambers, which are affected by the total oxygen and NO<sub>x</sub> concentrations. In the case of NO<sub>x</sub>, the delay is slightly higher than for  $\lambda^{-1}$  due to the sensing element location is in the inner cavity. It should also be noted that the delay tends to be overestimated during idling (about 2 s in NO<sub>x</sub> against 1.4 in the reality) due to the simple model structure (3.4). In this work, a constant delay is used for the implementation of models and fusion strategies, and values are shown in dashed lines in Figure 3.23 (about 0.75 s in NO<sub>x</sub> and 0.5 s in  $\lambda^{-1}$ , in line with the manufacturer technical data). The results obtained in subsequent chapters justify this decision, in addition to the simple programming of a constant delay against a variable one. Furthermore, the robustness, dispersion and ageing of a delay model should be assessed more in

detail, where scheduling strategies in line with the thesis algorithms might be explored.



**Figure 3.23.** Variable sensor delay modelled for  $z_{NO_x}$  and  $z_{\lambda^{-1}}$  outputs from the  $NO_x$  sensor during the second half of the NEDC A01 (left) and SDMP B5 (right). Higher peaks correspond to the idling operating conditions, which are overestimated. In general, the bigger  $m_f$  and  $n$ , the lower the delay.

### 3.5 Conclusions

This Chapter has described the experimental setup and the test campaign performed in order to design and validate the models and algorithms utilised in the current dissertation. Furthermore and considering the gas concentration sensors used for this work, the  $NO_x$  and  $\lambda^{-1}$  outputs from a  $ZrO_2$  sensor are characterised by means of steady-state tests for static calibration and transient steps for dynamic characterisation. More precisely:

- A simple on-engine experiment for characterising  $NO_x$  output is presented. The method uses the engine as a gas generator and the procedure consists of applying steps in SOI producing sharp variations in  $NO_x$  concentration. With a simple variation of  $2-3^\circ$  in SOI,  $NO_x$  concentration is varied whilst torque values and fluid conditions remain approximately constants.



- A similar procedure based on performing injection fuel rate steps in order to produce sharp variations on  $\lambda^{-1}$  is used for the dynamic characterisation of this signal.

These methods allow the determination of the dynamic response of the sensor (for both  $\text{NO}_x$  and  $\lambda^{-1}$ ) over its entire useful life and the detection of faults in the sensor. A simple mathematical modelling, taking as a reference a first order model response, is proposed, although more complex models might be utilised. According to this, the delay and the response time is modelled for both  $\text{NO}_x$  and  $\lambda^{-1}$ . In both cases, and even though there exists variability in the parameters, the delay and the response time are considered constants for the applications made in subsequent chapters. The good results of the models and fusion strategies for the proposed applications give validity to this decision, provided that a deeper analysis is advised for future robust implementations. Due to its nature, the methods could be implemented on-board for the online dynamic calibration and development of adaptive strategies.

### 3.A Gas concentration sensor specifications

Table 3.10 presents the relevant characteristics of the gas concentration sensors certified by the OEMs.

**Table 3.10.** *Relevant characteristics of the gas concentration sensors used for this dissertation. \*Data not specified by the OEM. FS is Flux Storage.*

	Horiba MEXA 7100DEGR	Bosch LSU 4.9	Uni- $\text{NO}_x$ sensor
Output signals	CO,CO <sub>2</sub> ,NO <sub>x</sub> ,THC,O <sub>2</sub>	$\lambda$	NO <sub>x</sub> , linear $\lambda$ , binary $\lambda$
Measuring range	NO <sub>x</sub> : 0-5000 ppm O <sub>2</sub> : 0-25 % vol. CO <sub>2</sub> : 0-12 % vol.	( $I_p$ ) 0.65 to air	NO <sub>x</sub> : 0-1500 ppm lin $\lambda$ : 0.75 to air bin $\lambda$ : >0.75 V ( $\lambda=0.9$ ) bin $\lambda$ : <0.2 V ( $\lambda=1.1$ )
Accuracy	NO <sub>x</sub> : $\pm 1.5\%$ FS O <sub>2</sub> : $\pm 1.5\%$ FS CO <sub>2</sub> : $\pm 1\%$ FS	$\lambda = 1.016 \pm 0.007$ $\lambda = 0.8 \pm 0.01$ $\lambda = 1.7 \pm 0.05$	NO <sub>x</sub> : 100 and 500 ppm $\pm 10\%$ 0 ppm $\pm 10$ ppm lin $\lambda$ : $\lambda = 1 \pm 6 (1000/\lambda)$ fresh bin $\lambda$ : $1.002 \pm 0.008$
Response times	NO <sub>x</sub> : < 900ms O <sub>2</sub> : < 750 ms CO <sub>2</sub> : < 800 ms	< 100 ms*	NO <sub>x</sub> : 750 ms $\lambda$ : 550 ms

### 3.B Principles of the $\text{NO}_x$ measurement

The term  $\text{NO}_x$  emissions includes the di-nitrogen oxide  $\text{N}_2\text{O}$ , nitrous oxide NO and the nitrogen oxide  $\text{NO}_2$  [24]; the most habitual at the engine exhaust

in diesel engines is NO, with percentages about 70-90% [25, 26], while the remainder is mainly NO<sub>2</sub>.

The gas analyser is capable of measuring NO<sub>x</sub> and separating between NO and NO<sub>2</sub> if necessary. For the NO<sub>x</sub> case, the total composition is first oxidised to NO<sub>2</sub> and then converted by a catalytic reaction to NO, that afterwards reacts with ozone (O<sub>3</sub>) for producing NO<sub>2</sub> and emitting photons. The produced light is detected with the CLD and is proportional to the NO concentration. For detecting NO (before oxidising), the thermal catalytic phase is avoided.

The measurement provided by the gas analyser for NO<sub>x</sub> is made in a dry basis and for converting to a standard wet basis, the correction  $K_w$  is calculated as follows

$$K_w = (1 - F_{fh} \frac{\dot{m}_f}{\dot{m}_a} (1 + 0.001H_a)) - K_{w2} \quad (3.5a)$$

$$K_{w2} = \frac{1.608H_a}{1000 + 1.608H_a} \quad (3.5b)$$

$$F_{fh} = \frac{1.969}{1 + \frac{\dot{m}_f}{\dot{m}_a}} \quad (3.5c)$$

where  $\dot{m}_a$  and  $\dot{m}_f$  are the air mass flow and the injected fuel mass flow [kg/h] respectively, while  $H_a$  is the absolute humidity.

Furthermore, the legislation fixes the factor  $K_{hd}$  for accounting with the ambient conditions

$$K_{hd} = \frac{1}{1 + A(H_a - 10.71) + B(T_{atm} - 298)} \quad (3.6a)$$

$$A = 0.309 \frac{\dot{m}_f}{\dot{m}_a} (1 + 0.001H_a) - 0.0266 \quad (3.6b)$$

$$B = -0.209 \frac{\dot{m}_f}{\dot{m}_a} (1 + 0.001H_a) + 0.00954 \quad (3.6c)$$

where  $T_{atm}$  [K] is the ambient temperature. Therefore, the NO<sub>x</sub> output after corrections is computed

$$\text{NO}_x = K_w K_{hd} \text{NO}_x^{gas \text{ analyser}} \quad (3.7)$$

A deeper explanation on the gas analyser procedures for measuring NO<sub>x</sub> and separating components is in [25].

In the case of the NO<sub>x</sub> sensor, the exhaust gas is diffused first to a first chamber, where the oxygen is measured, and then diffused again to the second

chamber. In the first chamber, NO species are oxidised to NO<sub>2</sub>, while in the second, total NO<sub>x</sub> is reduced to oxygen and nitrogen ions. These oxygen ions are pumped out and are proportional to the total NO<sub>x</sub> concentration. Note that the first oxidation step is required in order to get a correct measurement of total NO<sub>x</sub>, i.e. NO<sub>2</sub> contains two oxygen atoms while NO contains only one and then the oxygen ions resulted from the dissociation could not be directly related with total NO<sub>x</sub>. The static calibration was shown in Figure 3.14, where the fitting between ZrO<sub>2</sub> sensor and gas analyser is quite acceptable.

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# Chapter 4

## Control models for engine-out $\text{NO}_x$ and $\lambda^{-1}$

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### 4.1 Introduction

Control oriented modelling is aimed to predict engine variable with a low computational effort and is the basis of model based strategies, such as diagnosis or MPC [1]. Models usually have the possibility of predicting with acceptable accuracy the transient performance of the engine. Once that the Chapter 2 have already reviewed models for  $\text{NO}_x$  and  $\lambda^{-1}$ , this Chapter is devoted to the design of an ECU signals based virtual sensor for the estimation of engine-out  $\lambda^{-1}$  and  $\text{NO}_x$ . The reported results correspond to the DW12B DICI engine presented in the Chapter 3 with all AT devices removed.

## 4.2 $\lambda^{-1}$ model

Instead to refer to  $\lambda$ , results and discussion are centred in its inverse  $\lambda^{-1}$ , namely relative fuel-to-air ratio or simply richness, as it presents a bounded value (ranging from 0 to 1 in diesel engines, although slightly higher values might be reached during engine load transients).

Depending on the engine configuration, an air mass flow sensor ( $\dot{m}_a$ ) may be present (or mass flow can be estimated through the boost pressure), and the injected fuel mass flow ( $\dot{m}_f$ ) is usually estimated by the ECU. A simple model may be then proposed

$$x_\lambda^{-1} = 14.5 \frac{\dot{m}_f}{\dot{m}_a} \quad (4.1)$$

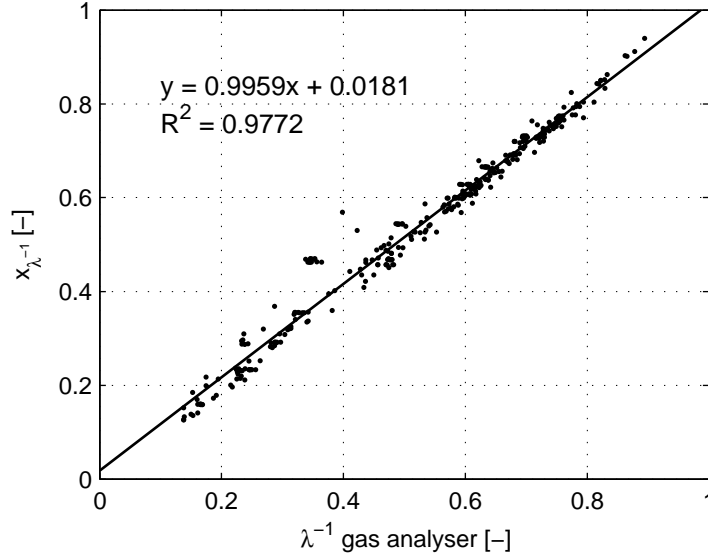
where  $\dot{m}_f$  and  $\dot{m}_a$  quantities are expressed in the same units, i.e. mg/str or kg/s, and 14.5 stands for the stoichiometric air-to-fuel ratio. Note that the model neglects the mass accumulation in the intake and exhaust manifolds (the reader is referred to the Chapter 2 for finding more elaborated models). Despite such effects should be considered for correcting  $\dot{m}_a$  during the engine transients (for example, when boost pressure is increased or when shifting between boosting modes [2]), they will be neglected hereinafter for simplicity.

Figure 4.1 compares model steady-state results ( $x_\lambda^{-1}$ ) with gas analyser measurements. As it can be easily noticed, the model provides a lower accuracy than that of the sensor (shown in Figure 3.15). Regarding the transient behaviour, the model exhibits an almost instantaneous response; the injected fuel mass estimate, although biased is fast; and the air mass flow measurement characteristic time is in the order of milliseconds, while the exhaust gas sensor response time can be of several hundreds and even seconds (depending on the considered set-up).

**$\lambda^{-1}$  model drift.**  $x_\lambda^{-1}$  output presents a significant bias that strongly depends on the operating conditions due to several facts:

- $m_f$  is obtained from internal calculations of the ECU, and response is fast and non-delayed, but is based on tabulated values which rely on rail pressure measurement and injection duration. The error of these tables can be significant when short injections or split injection strategies are applied. On the other hand, injector manufacturing discrepancies and ageing can create a significant unit-to-unit (and cylinder-to-cylinder [3–5]) dispersion in the actual injected quantity.





**Figure 4.1.**  $x_{\lambda^{-1}}$  and  $\lambda^{-1}$  from the gas analyser for different steady operating conditions; the linear fit and correlation coefficient are represented.

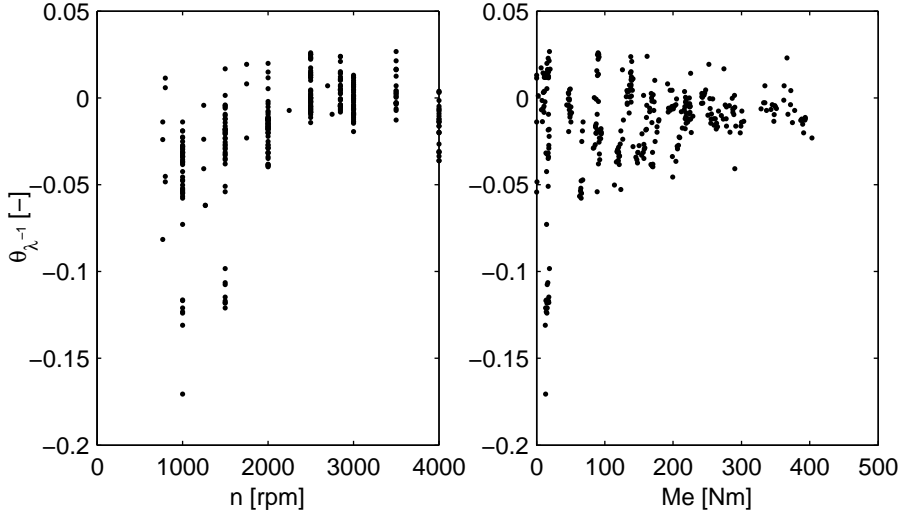
- While  $\dot{m}_a$  sensor response is judged to be fast and non-delayed enough for being directly used as a model input, the precision of the sensor is not very high (for example in [6] the non-systematic error of the air mass flow sensor is quantified with a standard deviation of 3.23%, in part due to flow pulsations in the intake line associated with the engine speed). On the other hand, during transient processes in which a significant variation of the air mass accumulated in the manifolds exists, the air mass flow measured in the intake line is different to that entering the engine [2]. This can be corrected accounting pressure variations in the  $\dot{m}_a$  estimation, although this is not considered in the current work.

With this scenery,  $x_{\lambda^{-1}}$  can be related to the actual value of  $\lambda^{-1}$  considering a bias  $\theta_{\lambda^{-1}}$

$$x_{\lambda^{-1}} = \lambda^{-1} - \theta_{\lambda^{-1}} \quad (4.2)$$

The model bias is not constant and varies with the operating condition, since  $\dot{m}_a$  error significantly depends on the engine speed, and  $\dot{m}_f$  error is affected by the injection profile, which is usually scheduled as a function of  $n$

and  $M_e$ . Figure 4.2 shows bias estimation for the measured steady points and its variation with engine speed and torque.



*Figure 4.2. Model bias for the steady-state tests.*

Additionally, the model bias varies with time due to the system drift, and a slow variation could be expected. At the same time, the model drift may be associated with injection soiling, air mass flow sensor drift, or the variations of leakage in the intake manifold and blow-by. In order to model that, both effects, the dependency with the operative conditions, and the drift, must be considered

$$\frac{d\theta_{\lambda^{-1}}}{dt} = \frac{\partial\theta_{\lambda^{-1}}}{\partial t} + \frac{\partial\theta_{\lambda^{-1}}}{\partial n} \frac{dn}{dt} + \frac{\partial\theta_{\lambda^{-1}}}{\partial m_f} \frac{dm_f}{dt} + \dots \quad (4.3)$$

Although the bias variation associated with the system drift ( $\partial\theta_{\lambda^{-1}}/\partial t$ ) is expected to be slow, the actual variation of the bias may be very fast, due to the ability of the engine of performing fast transition between operating conditions (defined by  $n$  and  $m_f$ ).

### 4.3 $\text{NO}_x$ model

A RT gray box  $\text{NO}_x$  model combining first principle equations with the use of operating-point dependent look-up tables is developed based on simple

correlations of engine available signals. This model describes a quasi-static representation of the engine plus some filters for describing actuators and sensor dynamics, where the scheduling points are  $n$  and  $m_f$ .

### 4.3.1 Previous considerations

Firstly, previous considerations related with the validity of a quasi-static representation and definition of engine operating points, the  $[O_2]_{int}$  estimation and the selection of the appropriate units for modelling  $NO_x$  are discussed.

**Comments on the engine dynamics and definition of the engine operating point.** The engine works in closed loop by mainly following set-point references for air mass flow ( $\dot{m}_a^*$ ) and boost pressure ( $p_{boost}^*$ ), which are controlled by modifying the actuators for EGR ( $u_{egr}$ ) and waste-gate ( $u_{wg}$ ). Other variables such as the coolant temperature ( $T_{cool}$ ) are measured and taken into account in the ECU controller. Speed ( $n$ ) and injection ( $m_f$ ) are used as scheduling variables in order to build the references for the controllers by means of look-up tables. Since the fuel path representative times are in the order of  $\mu s$ , this can be considered as instantaneous if compared with the air path dynamics. However, the major influence of the fuel path over the air path is with the smoke limiter, which limits the injected fuel rate until  $\dot{m}_a$  or  $p_{boost}$  (depending on the manufacturer) reach a certain limit in order to bound  $\lambda^{-1}$ . Therefore, the engine states ( $X$ ) could be modelled with the dynamic equation

$$\dot{X} = f(X, n, m_f, u_{egr}, u_{wg}, T_{cool}) \quad (4.4)$$

The set-point references for air and boost pressure ( $\dot{m}_a^*$  and  $p_{boost}^*$ ) are modelled by look-up tables scheduled by  $n$  and  $m_f$

$$\dot{m}_a^* = f(n, m_f) \quad (4.5a)$$

$$p_{boost}^* = f(n, m_f) \quad (4.5b)$$

The variables  $\dot{m}_a$  and  $p_{boost}$  are measured by sensors in diesel production engines, and if considering them as fast signals and neglecting mass storage effects (MAF sensor for measuring  $\dot{m}_a$  is usually installed upstream of the low pressure compressor), the dependency of the command signals  $u_{egr}$  and  $u_{wg}$  might be skipped as follows

$$\dot{X} = f(X, n, m_f, \dot{m}_a, p_{boost}, T_{cool}) \quad (4.6)$$

If the engine air path states can be modelled by quasi-static representations (dynamics do not have influence on the output) or by adding filters and

delays function only of the scheduling inputs  $n$  and  $m_f$ , then the dependency with  $\dot{m}_a$  and  $p_{boost}$  might also be modelled as function of  $n$  and  $m_f$ . The case of temperatures, associated with slow transients is more difficult and the measured engine variable  $T_{cool}$  is not able to exactly represent for instance in-cylinder conditions. The inner wall cylinder temperature influence deserves a specific consideration, and its influence is included in the model with a dynamic factor (see discussion bellow in Section 4.3). Anyway,  $T_{cool}$  is used as input for switching warm-cold engine strategies. Based upon this reasoning,  $X$  could be represented as

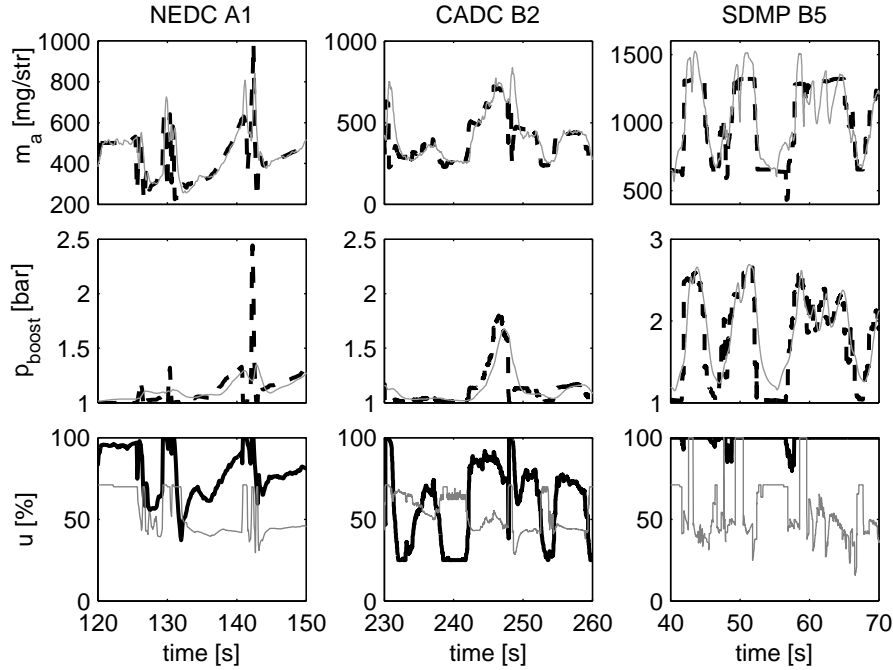
$$\dot{X} = f(X, n, m_f, T_{cool}) \quad (4.7)$$

Putting aside the thermal dynamic effects, further simplifications could be made if the engine performance is slow enough that  $\dot{m}_a$  and  $p_{boost}$  are able to track  $\dot{m}_a^*$  and  $p_{boost}^*$  fast and then, engine performance is simplified to

$$X = f(n, m_f) \quad (4.8)$$

Figure 4.3 shows examples of the set-point references and responses for the air and fuel path and for three different cycles (NEDC A1, CADC B2 and SDMP B5), including also in the bottom plots the variables  $u_{egr}$  and  $u_{wg}$  corresponding to the actuation signals for EGR and WG respectively. The cycles correspond in that order from the slowest to the fastest transitions. In the A1 test, the air path controllers track with a minimum error both  $\dot{m}_a^*$  and  $p_{boost}^*$ , especially for the case of  $\dot{m}_a$  which is logical due to the slower dynamics associated to the turbo inertia. The response for the B2 cycle is also good but transients are a little bit more aggressive than for A1, and the errors are slightly higher. For these two cycles, the quasi-static representation defined by (4.8) gives good results if the engine is already warmed up ( $T_{cool}$  is stable). However, for the case of B5 cycle, the transients are much more aggressive and errors on the  $\dot{m}_a$  and  $p_{boost}$  tracking are appreciable as well as the fact that the EGR valve is mostly closed during the cycle and the second TC is indeed working at some points (see Figure 4.3 when  $u_{wg}$  reaches 100%). For this cycle, the engine is running out of the nominal ( $\dot{m}_a^*$  and  $p_{boost}^*$  set-points differ from the other cycles) and the effect of the thermal loading due to the fast transients is critical.

Thus, the selection of the model structure for a given precision would depend not only on the engine characteristics but on the use cases. In all these cases and examples, it should be considered not only the thermal transient effects but the sensor dynamics. Therefore, (4.8) should consider the sensor model if the output is compared with a sensor output. The model structure



**Figure 4.3.** Air and fuel path response to variations in the driver's command ( $u_\alpha$ ) for different cycles. First and second row plots: --- is set-point  $\dot{m}_a^*$  and  $p_{boost}^*$ , and — is measured value of  $\dot{m}_a$  and  $p_{boost}$ . Last row plots: —  $u_{wg}$  [mg/str], —  $u_{egr}$  [mg/str] is the smoke limit for injection. Note that 100% in  $u_{egr}$  corresponds to the EGR valve fully closed and around 25% the EGR mass flow is maximum (the relation between  $u_{egr}$  command and the effective area is not linear). For the case of  $u_{wg}$ , 100% corresponds to WG fully closed and the two turbo shafts working.

proposed in this Chapter, aimed to be representative of the engine performance under different conditions, is built on the basis of (4.6) as a quasi-static model representation and an extra-state for representing sensor dynamics (now included in the equation):

$$X = f(n, m_f, \dot{m}_a, p_{boost}, T_{cool}) \quad (4.9a)$$

$$\dot{X} = f(X, t - \tau) \quad (4.9b)$$

where  $t$  is the time and  $\tau$  the sensor delay. This model structure, considering an state for including the thermal transient effect, is used after along the present Chapter for developing a  $NO_x$  model. Alternatively, in Section 6.3.1 a similar model to (4.8) will be considered for modelling  $NO_x$  in engine cycles.

**Intake oxygen estimation.** In the interest of developing a RT model for  $\text{NO}_x$ , the Arrhenius equation (2.6) can be arranged with two operating point dependent factors  $C_1$  and  $C_2$

$$\frac{d\text{NO}_x}{dt} = C_1 e^{-C_2/T} \quad (4.10)$$

where  $C_1$  includes gas intake concentrations (nitrogen and oxygen) while  $C_2$  is the exponential factor. According to [7], oxygen rate in the cylinder and operating conditions ( $n$  and  $m_f$  as discussed above) seem to be suitable variables for predicting  $\text{NO}_x$  emissions. Hence it is possible to find a new generic expression whose structure is equivalent to 4.10:

$$x_{\text{NO}_x} = f([\text{O}_2]_{\text{int}}, x_1, \dots, x_n) \quad (4.11)$$

where  $x_{\text{NO}_x}$  is the model output. This expression relates  $\text{NO}_x$  with  $[\text{O}_2]_{\text{int}}$  and different parameters ( $x_1, \dots, x_n$ ) including temperature and time reactions effects. The model, which is based on the thermal  $\text{NO}_x$  formation mechanism, might be calibrated for accounting for the  $\text{NO}_x$  formation by other mechanism.

Figure 4.4 shows  $\text{NO}_x$  versus  $[\text{CO}_2]_{\text{int}}$  in the intake manifold when varying EGR valve position for different speed  $n$  and injected fuel rate  $m_f$ .  $\text{NO}_x$  is normalised by  $\text{NO}_{x, \text{EGR}=0}$ , which corresponds to the maximum  $\text{NO}_x$  value at every operating point (equivalent to the EGR valve fully closed). The exponential fit in Figure 4.4 is clear, in spite of operating point dependency of the exponential factor is not modelled yet.

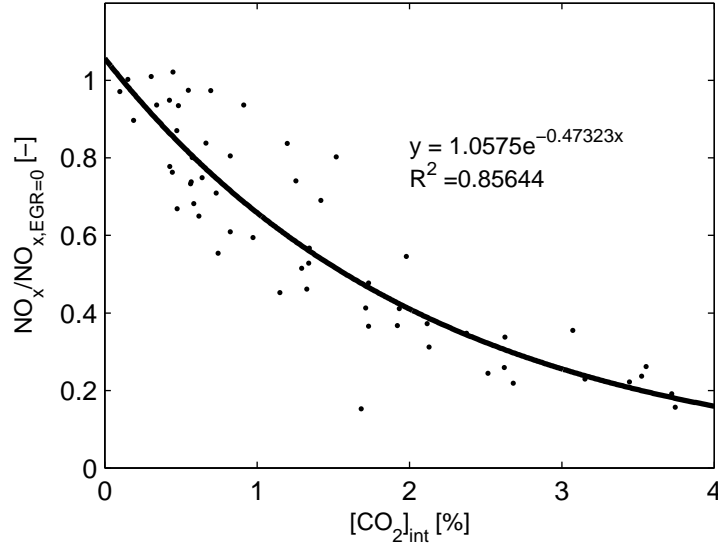
$[\text{O}_2]_{\text{int}}$  can be measured, but intake lambda probes are not always available in diesel engine production cars, and measurements errors related with the gas non-homogeneity may appear [8,9]. Considering the test engine, Section 3.2.1.2 already compared the intake oxygen estimation by a UEGO sensor with different samples in different intake positions measured by a gas analyser. Although the steady-state accuracy is acceptable, the UEGO sensor presents important problems linked with the pressure effects over the output signal, which drive to saturation and the need of compensation.

Alternatively,  $[\text{O}_2]_{\text{int}}$  can be modelled by relating  $[\text{O}_2]_{\text{int}}$ ,  $EGR$  and  $\lambda^{-1}$

$$[\text{O}_2]_{\text{air}}(1 - \lambda^{-1}) = [\text{O}_2]_{\text{exh}} \quad (4.12)$$

By the oxygen balance at the intake junction

$$[\text{O}_2]_{\text{int}} = [\text{O}_2]_{\text{air}}(1 - EGR) + [\text{O}_2]_{\text{exh}}EGR \quad (4.13)$$



**Figure 4.4.** Normalised  $NO_x$  emissions as function of  $[CO_2]_{int}$  for different nominal operating points in the diesel engine explained in Chapter 3 and coming from the steady-state campaign. Engine calibration is the standard given by the manufacturer.

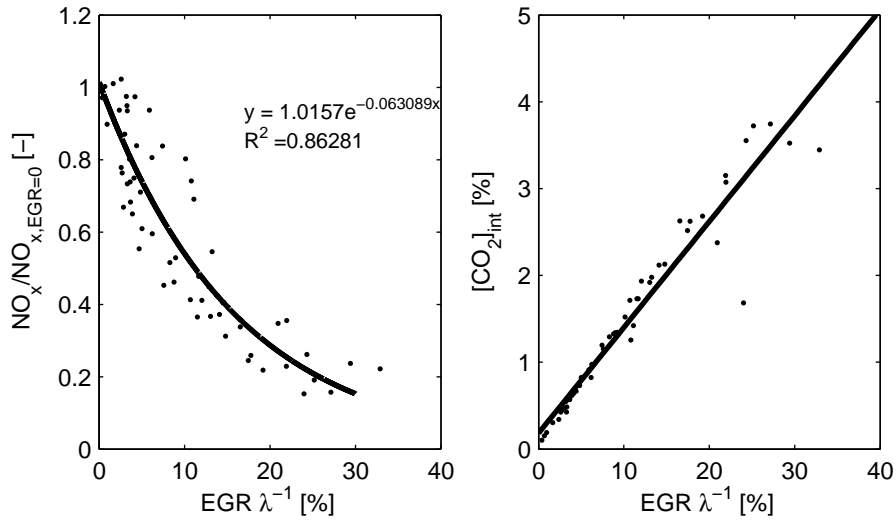
and substituting the former in the latter

$$[O_2]_{int} = [O_2]_{air}(1 - EGR\lambda^{-1}) \quad (4.14)$$

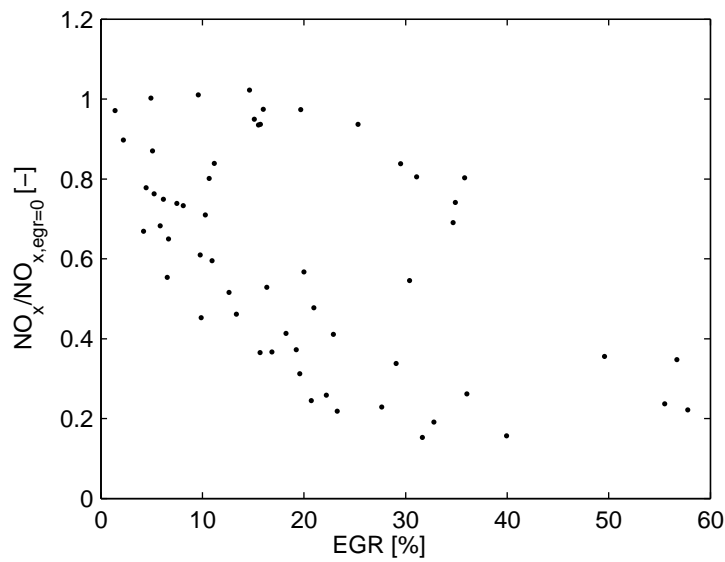
that shows a direct relationship between  $[O_2]_{int}$  and  $EGR\lambda^{-1}$ . The product  $EGR\lambda^{-1}$  is often called the inert gas rate and represents the actual portion of exhaust gas that has reacted with the injected fuel and then contains no available oxygen; e.g. Andersson *et al.* [10] use  $EGR\lambda^{-1}$  for designing a fast  $NO_x$  model. Figure 4.5 shows the  $NO_x$  fitting with  $EGR\lambda^{-1}$  and the fitting between  $EGR\lambda^{-1}$  and  $[CO_2]_{int}$ , where the expression (4.14) is proved.

On the other hand, Figure 4.6 shows the fitting of  $NO_x$  by using  $EGR$ . Results with  $EGR$  are much worse than with  $EGR\lambda^{-1}$  and it is difficult finding a tendency, which proves that  $EGR\lambda^{-1}$  seems a good alternative for estimating  $[O_2]_{int}$  (or  $[CO_2]_{int}$ ) when an intake lambda probe is not installed, although the uncertainties propagated from the different involved measurements are to be checked [9].

**About the  $NO_x$  model output units.** Sensor raw signal is measured in ppm, which indicates the relative molar concentration of  $NO_x$ , but the final



**Figure 4.5.** Left plot shows normalised  $\text{NO}_x$  emissions as function of  $\text{EGR}\lambda^{-1}$  for different operating points in the diesel engine explained in Chapter 3. Right plot shows the fitting of  $[\text{CO}_2]_{\text{int}}$  and  $\text{EGR}\lambda^{-1}$ .



**Figure 4.6.** Normalised  $\text{NO}_x$  emissions as function of EGR for different nominal operating points of the DW12B engine.



NO<sub>x</sub> mass depends on the exhaust mass flow and time. The conversion from ppm to g/h used in this work is

$$\dot{x}_{\text{NO}_x}^{g/h} = 0.001587 x_{\text{NO}_x}^{\text{ppm}} \cdot (\dot{m}_a^{kg/h} + \dot{m}_f^{kg/h}) \quad (4.15)$$

considering that molecular mass of NO<sub>x</sub> is 46.1 g/mol and exhaust mass flow density is 1.293 kg/m<sup>3</sup>, measured at 273 K and 101.3 KPa. The conversion of mg/str to kg/h is defined as

$$\dot{m}^{kg/h} = 1.2 \cdot 10^{-4} n \cdot m^{mg/str} \quad (4.16)$$

in a 4 cylinder basis. The use of mass flow variables is appropriate for taking into account total emissions and calculating the real impact on the environment. In addition, AT devices such as the LNT require a count of the accumulated NO<sub>x</sub> mass for regenerating the trap. Then, a mass flow seems also suitable for OBD purposes.

For computing total NO<sub>x</sub> emissions,  $\dot{x}_{\text{NO}_x}^{g/h}$  can be integrated in the considered cycle

$$M_{\text{NO}_x}^g = \int_{t_1}^{t_2} \dot{x}_{\text{NO}_x}^{g/h} dt \quad (4.17)$$

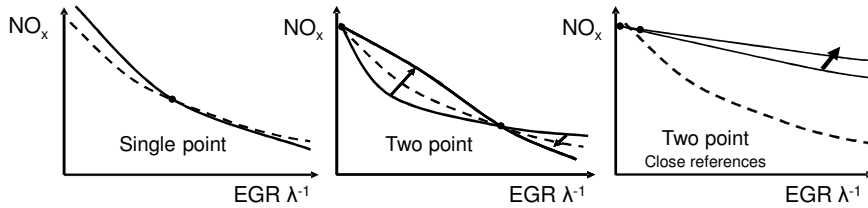
where  $M$  stands for total mass.

Commercial NO<sub>x</sub> sensors measure the relative volume flow of NO<sub>x</sub> in parts per million (ppm) and the conversion to g/h can induce some error, especially if  $\dot{m}_a$  presents some drift or filtering effect. Anyway, the use of g/h for online NO<sub>x</sub> tracking is justified for two reasons: (1) dilution effects are avoided as ppm is a relative measurement and (2) the use of g/h is coincident with the actual emissions. Dilution effects can distort the ppm output and can lead to sign shifts in the gradient when comparing ppm output and g/h at high load conditions (no EGR), i.e. an increase of g/h could result in a decrease of ppm if  $\dot{m}_a$  increases. On the other hand, minimum variations of g/h can lead to huge variations in ppm even though the relative gradient could be high in both two, especially at low load conditions (related with low  $\dot{m}_a$ ), leading to overestimate NO<sub>x</sub> emissions when tracking ppm. In this work, the NO<sub>x</sub> model output is calibrated for minimising the error on g/h, even though ppm output is also considered.

**Selection of the model structure.** In order to find a suitable model equation for a RT NO<sub>x</sub> model, different structures may be used [11,12]. Figure 4.7 shows a visual example of point dependent fittings: single (left plot) and two-point functions (centre and right plot). On one hand, if using a single point

function, model variations are centred around one reference. The point should be chosen as representative of the engine operation:  $\text{NO}_x$  emissions under nominal conditions (defined with the standard calibration of the OEM) or maximum emissions ( $EGR = 0$ ), which is used in previous figures. On the other hand, the two point function interpolates in a two reference basis, which in diesel engines could be around nominal EGR actuation and closed EGR (at least for the EGR area of the engine).

Both correlations have advantages and drawbacks: single point fitting is more flexible than two point for mathematical adjustment but with a smaller accuracy in points far from the nominal, however the single point one is easier to calibrate as only one map or reference must be tuned. Admittedly, extrapolation capability for the two-point function output when input values are out of the two-point area is compromised. See for instance right plot of Figure 4.7, the two references are quite close to each other and extrapolation is not fair. The same occurs in diesel engines in areas where nominal  $EGR\lambda^{-1}$  is low. Here, the single reference point model is used due to simplicity and robustness.



**Figure 4.7.** Single point reference model (left plot) VS two-point reference models (center and right plots).

### 4.3.2 A real Time $\text{NO}_x$ model

A RT  $\text{NO}_x$  model is proposed using  $EGR\lambda^{-1}$  for representing the intake oxygen jointly with maps and corrections fitting the operating point dependency [11]. The model scheme is outlined in Figure 4.8. In the top right and for clarifying the reader, engine positions subscripts are reminded: *boost* is downstream the intercooler, *int* the intake manifold, *egr* the EGR manifold at the intake junction and *exh* the exhaust. The main RT  $\text{NO}_x$  model requires two additional blocks: a mean value engine calculation of the EGR flow for obtaining  $EGR$  and the calculation of the fuel-to-air ratio  $\lambda^{-1}$ . The model

design is flexible and other structures for the EGR flow and  $\lambda^{-1}$  model or other measurements might be utilised just replacing the corresponding blocks.

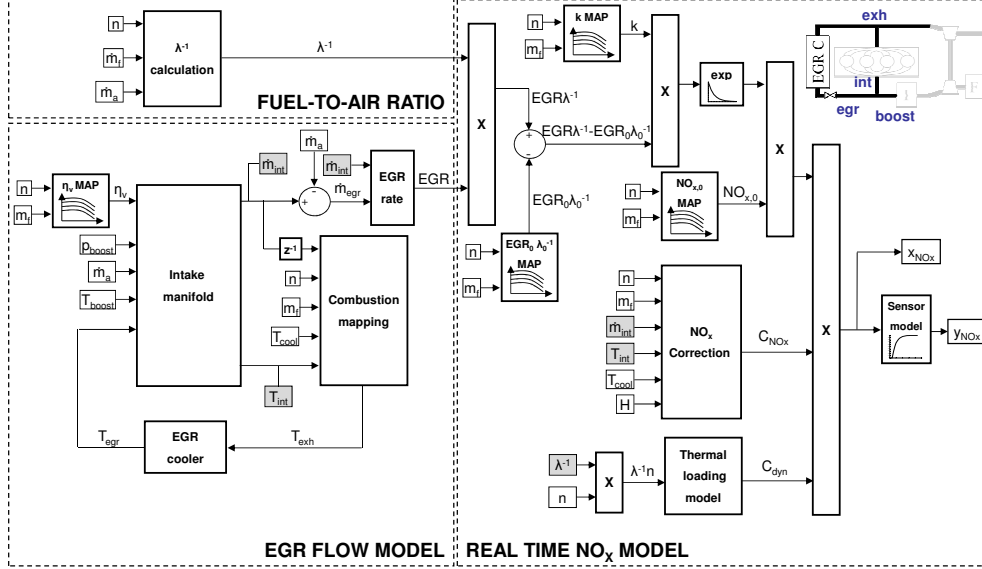


Figure 4.8. Structure of the RT  $NO_x$  model.

Following the discussion presented in the previous section about representation of engine states, an engine-out  $NO_x$  model is designed considering the state vector  $X$

$$X = [x_{NO_x} \quad y_{NO_x} \quad C_{dyn}]^T \quad (4.18)$$

where  $x_{NO_x}$  is the  $NO_x$  model output (actual  $NO_x$ ),  $y_{NO_x}$  is the filtered and delayed  $NO_x$  output (for comparing with  $NO_x$  sensor) and  $C_{dyn}$  is a dynamic factor for coping with in-cylinder temperature and explained after. The input signals contained in the vector  $U$  are available in the ECU and are shown in squares in the Figure 4.8

$$U = [n \quad m_f \quad \dot{m}_a \quad p_{boost} \quad T_{cool} \quad H]^T \quad (4.19)$$

where  $H$  is the humidity and the rest of variables have already been defined. Model output vector  $Y$  is the modelled sensor response

$$Y = y_{NO_x} \quad (4.20)$$

in order to have a comparable signal with  $NO_x$  sensor output during dynamic tests.  $x_{NO_x}$  indicates the raw or actual  $NO_x$ , which must be filtered and

delayed. The delayed first order discrete model identified in Chapter 3 is used

$$y_{\text{NO}_x} = z^{-\tau_{\text{NO}_x}/T_s} \frac{1 - a_{\text{NO}_x}}{1 - a_{\text{NO}_x} z^{-1}} x_{\text{NO}_x} \quad (4.21)$$

where  $a_{\text{NO}_x}$  is the sensor response time and  $\tau_{\text{NO}_x}$  is the total sensor delay (see Chapter 3 for more information).

Finally, constants are also marked with squares and gray squared variables are used as shortcut for avoiding crossed lines in the Figure 4.8.

**Model hypotheses.** The following hypotheses are assumed:

- The model is programmed and calibrated in discrete form and frequency used for simulation is 50 Hz.
- The model is based on a set of static look-up tables, curves and parameters with the addition of discrete filters and lag blocks in order to consider system dynamics.
- Intake conditions ( $\dot{m}_{int}$  and  $T_{int}$ ) are calculated on the basis of the volumetric efficiency and an EGR flow model by assuming constant pressures upstream and downstream of the intake junction and an adiabatic mixing model for the temperature.
- The engine combustion is tabulated by using look-up tables related to nominal conditions and different variations of the intake conditions.
- The EGR path is modelled with a cooler model and a first order filter for accounting with the manifold dynamics.
- Turbocharger effects are bypassed by directly using sensor signals  $\dot{m}_a$  and  $p_{boost}$ , the former located upstream of the compressor and the latter downstream of the intercooler.

In the next, the model and sub-model equations are described.

**RT  $\text{NO}_x$  model equation.** The  $\text{NO}_x$  model is presented in (4.22) and consists in a single function referenced to conditions with nominal EGR for calculating  $x_{\text{NO}_x}$ , which is affected by an exponential  $EGR\lambda^{-1}$  correction and additional factors  $C_{\text{NO}_x}$  and  $C_{dyn}$ . This structure guarantees that nominal tests corrections are eliminated minimising the nominal error and subscript 0

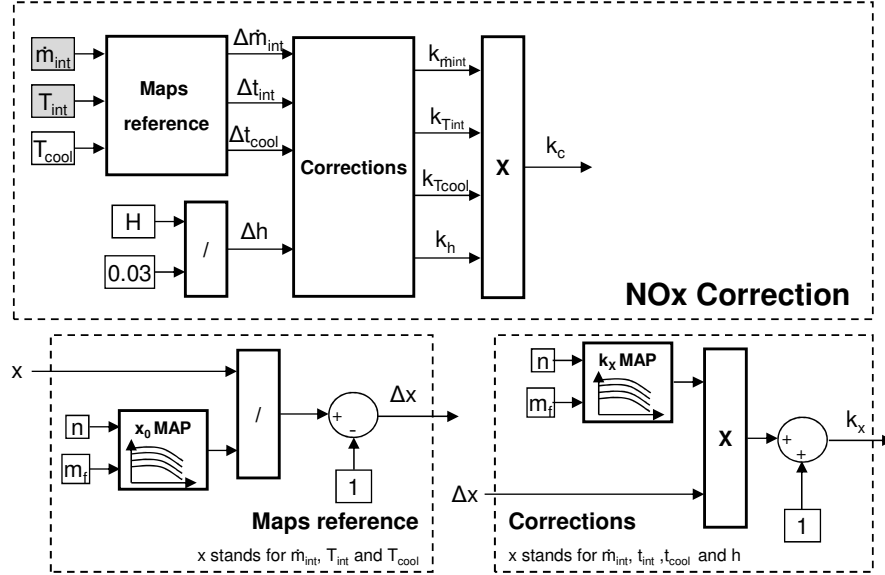


Figure 4.9. Structure of the factor  $C_{NO_x}$  calculation for correcting  $NO_x$  output.

goes with nominal conditions. Then, the model  $x_{NO_x}$  output is calculated as follows

$$x_{NO_x} = NO_{x,0} \cdot e^{-k_{NO_x} \cdot (EGR\lambda^{-1} - EGR_0\lambda_0^{-1})} \cdot C_{NO_x} \cdot C_{dyn} \quad (4.22)$$

where  $C_{NO_x}$  and  $C_{dyn}$  are used as correction factors to the model structure, and are explained in subsequent paragraphs.

**Correction factors.** The factor  $C_{NO_x}$ , defined with the product

$$C_{NO_x} = C_{\dot{m}_{int}} \cdot C_{T_{int}} \cdot C_{T_{cool}} \cdot C_h \quad (4.23)$$

is calculated on the basis of previous experience and tests made in the engine (see [13] for a sensitivity study of  $NO_x$  to relevant engine inputs). For that, 4 important effects are included in  $C_{NO_x}$ : (4.24a)  $C_{\dot{m}_{int}}$  indicates variations of the intake mass flow  $\dot{m}_{int}$  [8]; (4.24b)  $C_{T_{int}}$  copes with intake temperature  $T_{int}$ ; (4.24c)  $C_{T_{cool}}$  relates engine temperature (especially for distinguishing cold and warm conditions [14]); (4.24d)  $C_h$  considers Humidity  $H$  (if a humidity sensor is available). The expected  $NO_x$  sensitivity for the different parameters is depicted at the right side of (4.24). Figure 4.9 shows the blocks diagram of the  $C_{NO_x}$  model.  $k_{\dot{m}_{int}}$ ,  $k_{T_{int}}$ ,  $k_{T_{cool}}$ ,  $k_h$  are calibrated by  $[n, m_f]$  dependent

maps. If engine is operating at nominal conditions then  $C_{\dot{m}_{int}}, C_{T_{int}}, C_{T_{cool}} = 1$ . Equations are shown in the following.

$$C_{\dot{m}_{int}} = 1 + k_{\dot{m}_{int}} (\dot{m}_{int}/\dot{m}_{int,0} - 1) \quad \uparrow \dot{m}_{int} \downarrow \text{NO}_x^1 \quad (4.24a)$$

$$C_{T_{int}} = 1 + k_{T_{int}} (T_{int}/T_{int,0} - 1) \quad \uparrow T_{int} \uparrow \text{NO}_x \quad (4.24b)$$

$$C_{T_{cool}} = 1 + k_{T_{cool}} (T_{cool}/T_{cool,0} - 1) \quad \uparrow T_{cool} \uparrow \text{NO}_x \quad (4.24c)$$

$$C_h = 1 + k_h H/0.03 \quad \uparrow H \downarrow \text{NO}_x^2 \quad (4.24d)$$

**Thermal loading model.**  $T_{cool}$  represents the engine coolant temperature and is measured on-board.  $T_{cool}$  is used on-board for defining warm and cold strategies. However, neither the engine block temperature nor the in-cylinder temperature ( $T_{cyl}$ ) are measured on-board (and rarely in test benches) and are factors of great importance for  $\text{NO}_x$  formation in diesel engines. If the engine is working on steady-state operation and the variables are already stabilised,  $T_{cyl}$  might be approximated as a function only of the operating point conditions and  $T_{cool}$  as follows

$$T_{cyl} = f(T_{cool}, n, m_f) \quad (4.25)$$

But if the engine is in transient operation, e.g. a tip-in, it is necessary to consider an extra degree-of-freedom since  $T_{cyl}$  will be related with the historic of the engine load. Note that  $T_{cool}$  itself is not capable of representing actual  $T_{cyl}$  and in addition, it is controlled in closed loop for tracking a reference making that

$$T_{cyl} \neq f(T_{cool}, n, m_f) \quad (4.26)$$

An option is installing a temperature sensor at the exhaust but even though this option should be interesting for improving the model accuracy, these sensors are not usually available on-board.

An alternative is considering the heat conductive problem between the cylinder walls and the coolant. The temperature wall affects the in-cylinder temperature and should be considered as a state of the model. If analysing

<sup>1</sup>If  $p_{boost}^*$  set-point is increased for a constant air mass flow ( $\dot{m}_a$ ) set-point, then  $u_{egr}$  will be opened and more  $\dot{m}_{egr}$  will enter the cylinder (consequently  $\dot{m}_{int}$  will be higher), producing that  $\text{NO}_x$  should be lower. This effect might be shifted if  $u_{egr}$  is already opened, especially at low load and regime areas, or engine is running out of the EGR area.

<sup>2</sup>Moisture increases the air specific heat  $c_p$ , diminishing in-cylinder temperature  $T_{cyl}$ .

the problem, as the gas temperature variation ( $\Delta T_{cyl}$ ) state defines the heat power released ( $\dot{Q}$ )

$$\dot{Q} \propto \Delta T_{cyl} \quad (4.27)$$

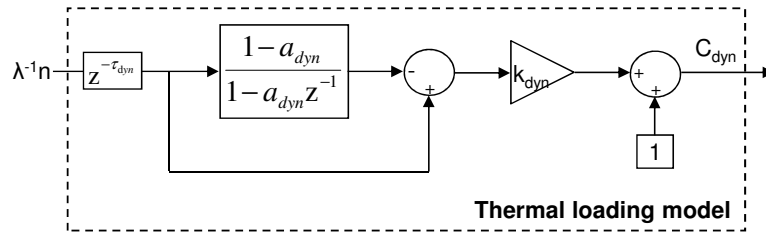
while the total heat dissipation between the cylinder and the coolant depends on  $\Delta T_{cyl}$  and the considered time, which is inversely proportional to  $n$ , then

$$Q = k \Delta T_{cyl} \frac{1}{n} \quad (4.28a)$$

$$\Delta T_{cyl} \propto \lambda^{-1} n \quad (4.28b)$$

which shows that  $\lambda^{-1} n$  is related with the thermal load in the cylinder and could be used as an estimator for the problem.

Therefore, a multiplicative factor  $C_{dyn}$  to the NO<sub>x</sub> model output is proposed. This factor augments the state vector  $X$  and copes with the engine loading and indirectly  $T_{cyl}$  effects. The correction factor is designed for correcting the NO<sub>x</sub> output only during transient, since steady state effects due to changes in  $T_{cool}$  are yet corrected through  $C_{T_{cool}}$ . The main equation is presented in (4.29) and block diagram shown in Figure 4.10.  $C_{dyn}$  is built with the filtered difference of  $\lambda^{-1} n$ , which represents engine thermal load variation (other options have also been checked, but this has given the best results for the model, which is in line with the discussion above), and a calibrated gain  $k_{dyn}$ . When the engine is steady, so  $\lambda^{-1} n$  is,  $C_{dyn} = 0$ . This factor is set to one for steady-state simulations.

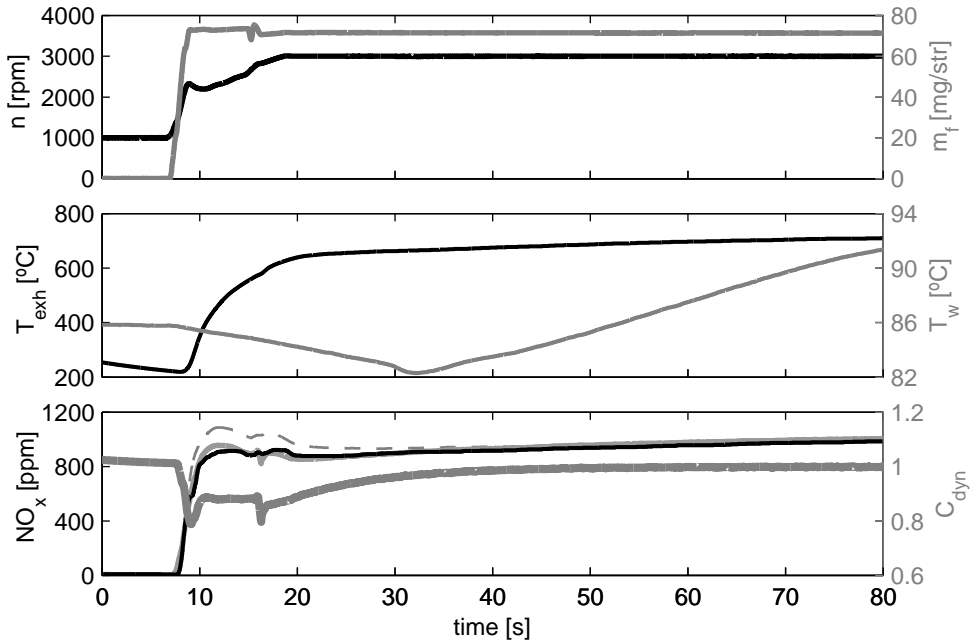


**Figure 4.10.** Block diagram of the thermal loading factor  $C_{dyn}$ .

$$C_{dyn} = 1 + k_{dyn} (\lambda^{-1} n) \left( 1 - z^{-\tau_{dyn}} \frac{1 - a_{dyn}}{1 - a_{dyn} z^{-1}} \right) \quad (4.29)$$

Figure 4.11 shows relevant signals during a sharp engine transient in  $n$  and  $m_f$ .  $T_{exh}$  is measured with a sensor and compared with  $T_{cool}$  showing two effects: First,  $T_{cool}$  is much more filtered and delayed than  $T_{exh}$ , and second,

the sensitivity of  $T_{cool}$  is not sufficient for representing the actual thermal transient occurring on-engine, i.e.  $T_{cool}$  hardly varies from 82 to 92 °C while  $T_{exh}$  goes from 200 to 700 °C. Bottom plot shows the value  $C_{dyn}$  tuned for the  $\text{NO}_x$  model, being lower than 1 during the transient but finally converging to 1. This is for representing that engine comes from a colder state. In that plot,  $\text{NO}_x$  sensor ( $z_{\text{NO}_x}$ ) is compared with  $y_{\text{NO}_x}$  model output (see (4.21)) without  $C_{dyn}$  (it overestimates  $\text{NO}_x$ ) and with  $C_{dyn}$ .



**Figure 4.11.**  $\text{NO}_x$  and thermal response during a sharp transient. Top plot:  $n$  and  $m_f$  evolution. Middle plot: Exhaust temperature ( $T_{exh}$ ) and coolant temperature ( $T_{cool}$ ) during the transient. Dynamics of both signals, especially for  $T_{cool}$ , are slow, which affects final  $\text{NO}_x$ . Bottom plot and left axis: —  $z_{\text{NO}_x}$  by sensor, - - -  $y_{\text{NO}_x}$  with  $C_{dyn} = 1$ , —  $y_{\text{NO}_x}$  using  $C_{dyn}$ . Bottom plot and right axis: —  $C_{dyn}$  factor calculated by the thermal model.

**The effect of the injection parameters.** The injection duration ( $t_{id}$ ), the start of injection (SOI) and the common rail pressure ( $p_{rail}$ ) are critical for  $\text{NO}_x$  emissions [15] and so is the engine speed ( $n$ ) that is coupled with the others, which here is a model input. For the current structure, the engine



commercial calibration for the injection settings is used being the injected fuel mass ( $m_f$ ) the representative variable. For that reason, all nominal conditions (marked with subscript 0) are related to the set  $[n m_f]$ , i.e every pair  $[n m_f]$  defines nominal  $u_{soi,0}$ ,  $u_{prail,0}$  and  $u_{id,0}$ . The engine torque ( $M_e$ ) could be used instead of  $m_f$ , but even though both  $M_e$  and  $m_f$  are estimated by the ECU,  $m_f$  is preferred as used in the ECU of the considered engine. SOI or  $prail$  could be considered as exponential corrections if different from the nominal, and such effects should be taken into account in the combustion model (this could be useful for controlling the fuel path).

#### 4.3.2.1 The EGR flow model

The EGR rate

$$EGR = \frac{\dot{m}_{egr}}{\dot{m}_{int}} \quad (4.30)$$

is calculated on the basis of a mean value model of the volumetric efficiency ( $\eta_v$ ) using boost pressure ( $p_{boost}$ ) and air mass flow ( $\dot{m}_a$ ) inputs coming from commercial sensors. The model block diagram is shown in Figure 4.12.

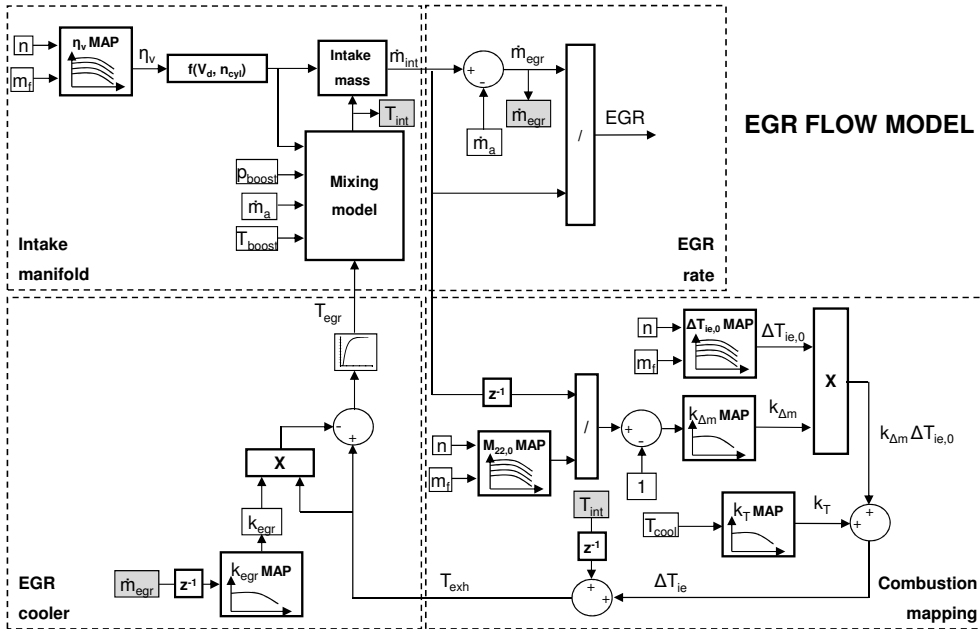


Figure 4.12. Structure of the EGR flow model.

The intake manifold mass flow ( $\dot{m}_{int}$ ) is obtained from  $\eta_v$  (mapped with the steady-state tests).  $p_{boost}$  is assumed constant in the manifolds that confluence at the intake ( $egr, boost, int$ ) while  $T_{boost}$  is measured and  $T_{exh}$  calculated by means of a combustion model based on maps and corrections over  $T_{cool}$  and  $\dot{m}_{int}$ .

$$\dot{m}_{int} = \frac{\eta_v n_{cyl} V_d n p_{boost}}{2 R T_{int}} \quad (4.31)$$

$n$ ,  $m_f$ ,  $T_{boost}$  and  $\dot{m}_a$  are measured by sensors, the unit displaced volume ( $V_d$ ) and the number of cylinders ( $n_{cyl}$ ) are known beforehand and a model is proposed for  $T_{int}$ . For instance, Wahlström *et al.* [16] propose an isothermal model ( $T_{int} = T_{boost}$ ); for the current work a mixing model based on the energy conservation at the intake junction is utilised, i.e. intake volume is small and mixing is so fast that no heat is transferred to the walls.

$$\dot{m}_{int} T_{int} = \dot{m}_a T_{boost} + \dot{m}_{egr} T_{egr} \quad (4.32)$$

and from (4.31), the product  $\dot{m}_{int} T_{int}$ , defining the intake enthalpy flow, can be inferred as

$$\dot{m}_{int} T_{int} = \frac{\eta_v n_{cyl} V_d n p_{boost}}{2 R} \quad (4.33)$$

and rewriting the mass equilibrium at the intake junction

$$\dot{m}_{egr} = \frac{\eta_v n_{cyl} V_d n p_{boost}}{2 R T_{int}} - \dot{m}_a \quad (4.34)$$

and finally substituting in (4.32),  $T_{int}$  can be solved

$$T_{int} = \frac{\eta_v n_{cyl} V_d n p_{boost} T_{egr}}{\eta_v n_{cyl} V_d n p_{boost} + 2 R \dot{m}_{boost} (T_{egr} - T_{boost})} \quad (4.35)$$

Two additional models (see Figure 4.12) are needed in order to calculate  $T_{egr}$  (not usually measured on commercial engines): a  $T_{exh}$  model and an EGR cooler efficiency model.

**$T_{exh}$  model.** Temperature increase between the intake and the exhaust  $\Delta T_{ie}$  is calculated from a steady map and corrections

$$T_{exh} = T_{int} z^{-1} + \Delta T_{ie} \quad (4.36)$$

adding the discrete delay  $z^{-1}$  for coping with causality.

$\Delta T_{ie}$  can be calculated on the basis of heat release functions (e.g. Wiebe for Watson functions) and solving the principle equations of the cylinder, or fitted to experimental values. Since the results of such physical approach may be calculated beforehand, a map based model coherent with the full RT  $NO_x$  model is proposed with two corrections based on  $\dot{m}_{int}$  and  $T_{cool}$

$$\Delta T_{ie} = \Delta T_{ie,0} \cdot k_{\Delta m} + k_T \quad (4.37)$$

Maps depending on  $[n, m_f]$  are used for generating references for  $\Delta T_{ie,0}$  and  $\dot{m}_{int,0}$ . Then  $\Delta m$  is calculated

$$\Delta m = \frac{\dot{m}_{int} z^{-1}}{\dot{m}_{int,0}} - 1 \quad (4.38)$$

and defining the calibrated map output

$$k_{\Delta m} = f(\Delta m) \quad (4.39)$$

Finally, term  $k_T$  represents the steady-state effect of the wall cylinder temperature effect and is calculated from the curve

$$k_T = f(T_{cool})$$

noting that  $T_{cool}$  adds a degree of freedom to the problem as there exists a closed loop control that manages the coolant mass flow and the set-point for  $T_{cool}$ . The difference with respect to the thermal factor  $C_{dyn}$  is that the latter copes with transients, while  $k_T$  with the static effect.

If SOI and  $p_{rail}$  are different from the nominals used for generating the nominal maps, then the additional factors  $k_{SOI}$  and  $k_{p_{rail}}$  should be calibrated in the combustion model.

**EGR cooler model.**  $T_{egr}$  is calculated by a temperature drop and a first order discrete filter with response time  $a_{egr}$  and a transport delay  $\tau_{egr}$

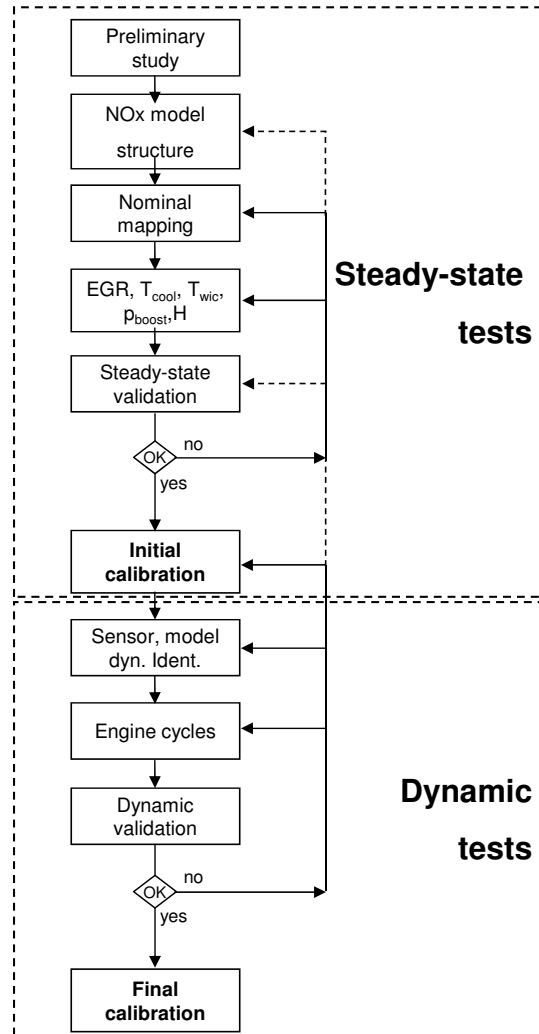
$$T_{egr} = z^{-\tau_{egr}} \frac{1 - a_{egr}}{1 - a_{egr} z^{-1}} (T_{exh} - k_{egr} (T_{exh} - T_{cool})) \quad (4.40)$$

where  $k_{egr}$  is calculated by the curve

$$k_{egr} = f(\dot{m}_{egr}) \quad (4.41)$$

and  $\tau_{egr}$  is function of the engine characteristic time

$$\tau_{egr} = \frac{k_{\tau_{egr}}}{n} \quad (4.42)$$



**Figure 4.13.** Scheme of the calibration and validation procedure for the  $\text{NO}_x$  model.

### 4.3.3 Tuning methodology

The tuning and validation methodology is shown in Figure 4.13, splitting the steady-state and dynamic testing. Dashed lines show critical paths, e.g. once that  $\text{NO}_x$  structure has been fixed after preliminary studies that include  $[\text{CO}_2]_{\text{int}}$  and sensitivity analysis, changes in the  $\text{NO}_x$  structure are expected to be minimum.

**Table 4.1.** Mean errors when fitting the model with different input variables.  $[\text{CO}_2]_{\text{int}}$  from the gas analyser is the mean of the 4 extractions of the intake manifolds as discussed in Section 3.2.1.2.  $EGR\lambda^{-1}$  is calculated on the basis of the EGR mean value flow model and the  $\lambda^{-1}$  model. Signal from the intake lambda probe ( $I_p$ ) is corrected by pressure and some data points with faulty measurements have been removed because they present problems with the sensor saturation.  $\text{NO}_x$  output is adjusted in g/h.

	$EGR\lambda^{-1}$	Corrected Ip	$[\text{CO}_2]_{\text{int}}$ from gas analyser
g/h	9.31	9.78	8.54
%	8.09	8.96	7.58

**Static tuning.** The steady tests are considered for having a representative sample of the operative engine range and calibrating nominal maps, main parameters and curves by using LS, with the following sequence:

- From all tests,  $\eta_v$ ,  $\dot{m}_{egr}$  and temperatures differences are calculated and used for fitting the maps as function of the engine speed and fuel mass.
- From nominal tests, maps are built for  $\text{NO}_{x,0}$ ,  $EGR_0\lambda_0^{-1}$  and  $T_{cool,0}$  as function of the engine speed and fuel mass.
- From all tests, the RT  $\text{NO}_x$  model is simulated in order to estimate the results for  $\dot{m}_{int}$  and  $T_{int}$ . The results are used for identifying  $\dot{m}_{int,0}$  and  $T_{int,0}$ .
- The ratios  $\dot{m}_{int}/\dot{m}_{int,0}$ ,  $T_{int}/T_{int,0}$  and  $T_{cool}/T_{cool,0}$  are calculated from all tests.
- Curves  $k_{\text{NO}_x}$ ,  $k_m$ ,  $k_{T_{int}}$ ,  $k_{T_{cool}}$  and  $k_h$  are fitted by LS of the main model equation (4.22). This adjustment is made independently by using the steady-state data set. In addition to  $EGR\lambda^{-1}$  and in order to compare results with other possibilities for estimating  $[\text{O}_2]_{\text{int}}$ ,  $I_p$  from an intake UEGO sensor and  $[\text{CO}_2]_{\text{int}}$  from the gas analyser are evaluated. The results are presented in Table 4.1. The best fit is for  $[\text{CO}_2]_{\text{int}}$  but this is not available on-board, while it can be used for tuning and validation. Therefore, the minimum error is achieved with the variable  $EGR\lambda^{-1}$ , calculated in (4.14).

**Dynamic tuning.** Dynamic tests are used for adjusting the  $C_{dyn}$  model, the path dynamics (e.g. EGR path) and for the fine tuning of the maps and curves fitted with the steady-state tests. Furthermore, the step tests are used for calibrating the sensors; and the cycles for validating the model in dynamic

state. If the validation is not satisfactory, the dynamic cycles can be repeated and identifications checked. If the error persists, it could be advisable to come back to the static calibration and re-tune the unreliable parameters (see Figure 4.13).

Even though not included in the scheme, tests from different campaigns could lead to drift because of the ageing or other effects that are not well modelled. Anyway, the repetition of a specific DOE will not avoid the model drift, and for sure, it is not a cost-effective solution. For that, adaptive strategies are useful as discussed in next chapters.

#### 4.3.4 $\text{NO}_x$ model results

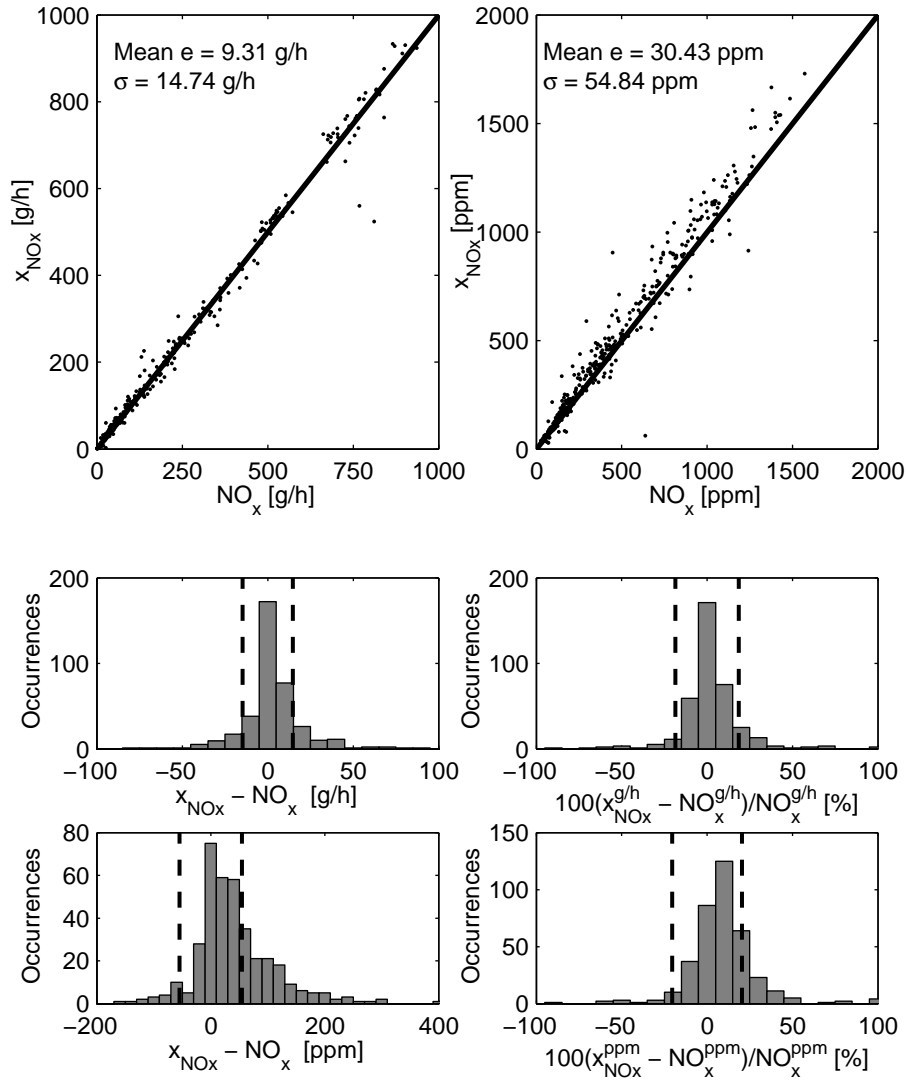
In the following,  $\text{NO}_x$  model results with static and dynamic tests are presented. Furthermore, error metrics are built in order to determine the model accuracy.

##### 4.3.4.1 Steady-state results

Top plots of Figure 4.14 show  $x_{\text{NO}_x}$  and  $z_{\text{NO}_x}$  (from the  $\text{NO}_x$ -ZrO<sub>2</sub> sensor) for both ppm and g/h units as well as histogram plots of relative and absolute errors. The mean error is 9.31 g/h, with a minimum deviation of  $\sigma=14.74$  g/h proving the validity of the calibration. The results by evaluating the ppm output present a mean error of 30.43 ppm and  $\sigma = 54.84$  ppm. It must be noted that errors of 1% in input variables can produce final  $\text{NO}_x$  errors of 33%, measured in a predictive and physics-based  $\text{NO}_x$  model (see the  $\text{NO}_x$  sensitivity study published in [13]). The maximum error in a DDM, such as the one presented here, is bounded by the tables grid and only the ageing and dispersion errors can affect, problems that exist also in more complex models. This fact and the results presented in this subsection justify the validity of a data-driven  $\text{NO}_x$  model, as well as the efficient computation and simple programming.

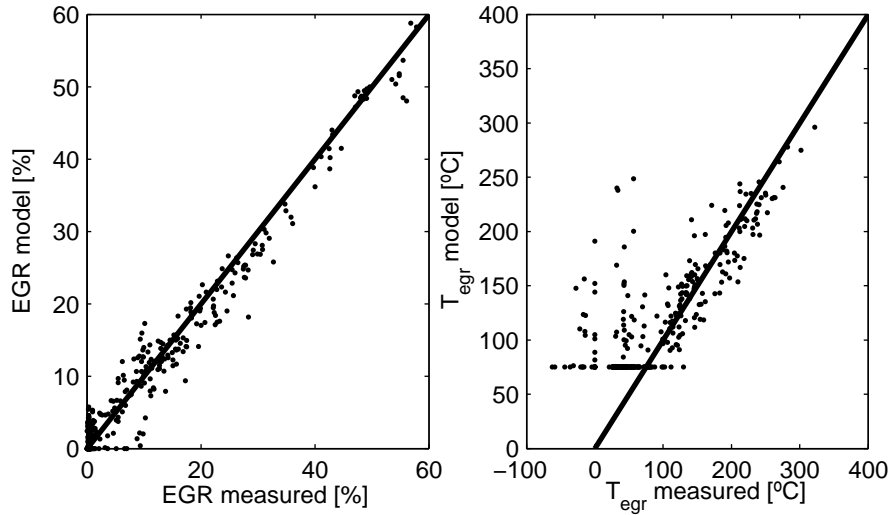
Medium and bottom plots of Figure 4.14 show absolute and relative error metrics for both outputs in g/h and ppm. Again, only a minimum percentage of points (total are 363) are out of the  $\pm\sigma$  boundary, also plotted in those plots. If a higher accuracy on ppm output is required, tuning objective can be easily shifted to ppm.

Additional information from the model can be extracted to compare intermediate variables (calculated by the model) and the corresponding measurements, as for instance for the mean value EGR flow model. A multi-objective



**Figure 4.14.** RT  $\text{NO}_x$  model results with steady tests. Top plots: Model fitting with respect to the  $\text{NO}_x$  sensor output. Medium plots: Absolute and relative error in percentage with respect to the  $\text{NO}_x$  sensor output and for g/h output ( $x_{\text{NO}_x}^{\text{g/h}}$ ) of the model. Bottom plots: Absolute and relative error in percentage with respect to the  $\text{NO}_x$  sensor output and for ppm output ( $x_{\text{NO}_x}^{\text{ppm}}$ ) of the model.

optimisation could be programmed if other model states should be important for control. However, here  $x_{\text{NO}_x}$  output error has been minimised. The left plot of Figure 4.15 compares  $EGR$  calculated by the model and by using  $[\text{CO}_2]_{int}$  averaged from the 4 extractions of the in-cylinder manifolds. The fitting is acceptable in the whole range. Results for  $T_{egr}$  are shown in the right plot of Figure 4.15, and the fitting in general is also acceptable, but there exists a high dispersion around the centre line. This is mainly due to areas with low or null EGR, where  $T_{egr}$  does not make any sense ( $\dot{m}_{egr}=0$ ).



**Figure 4.15.** Relevant results from the EGR flow model. Left plot:  $EGR$  calculated by the model and from the gas analyser measurements. Right plot:  $T_{egr}$  calculated by the model and measured by a sensor.

**Model robustness.** This section turns out the attention to the model robustness analysing the model sensitivity to errors in the main parameters and inputs. 2% disturbances in the parameters presented in Table 4.2 are applied and the final variation in  $\text{NO}_x$  model output is stored, e.g. a change of 2% in  $n$  produce an error of 4.51% in  $\text{NO}_x$ . Parameters with more influence on model output are  $n$ ,  $m_f$  and  $\dot{m}_a$ . These signals basically represent the engine operating conditions, and define the EGR setting (jointly with  $p_{boost}$  that presents a minor influence) and the trend was already expected.  $p_{boost}$  and  $\eta_v$  have the same influence as expected because both are multiplied for calculating  $\dot{m}_{int}$  and consequently  $\dot{m}_{egr}$ .



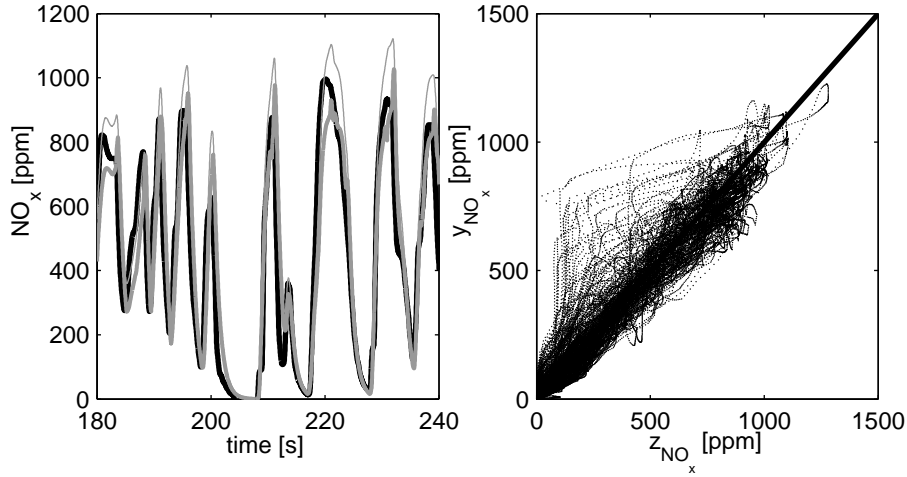
**Table 4.2.** Results over the final  $\text{NO}_x$  model output of applying variations of 2% on different variables.

Variable	Variation [%]	Mean error [%]	STD error [%]	Max. error [g/h]	Min. error [g/h]
$n$	+2	4.51	3.27	15.23	-1.41
$m_f$	+2	3.67	2.53	13.42	-0.36
$m_{boost}$	+2	3.42	3.50	12.60	-1.12
$\text{NO}_{x,0}$ MAP	+2	2	0.00	2.00	2.00
$\dot{m}_{int}$ MAP	-2	1.79	0.74	3.16	0.06
$\eta_v$ MAP	-2	1.03	2.78	8.29	-2.93
$p_{boost}$	-2	1.03	2.78	8.29	-2.93
$T_{cool}$	+2	0.97	0.39	1.81	0.08
$EGR_0 \lambda_0^{-1}$	+2	0.66	0.79	3.22	0.00
$k_{egr}$ MAP	-2	0.16	0.31	1.39	-0.51
$\Delta T_{ie}$ MAP	+2	0.06	0.14	0.97	-0.14
$k_{\text{NO}_x}$ MAP	+2	0.04	0.48	2.46	-2.42

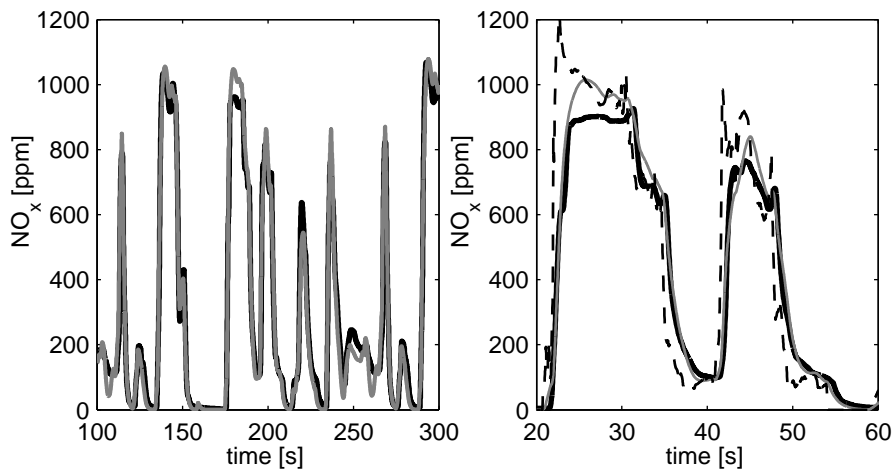
An important conclusion of this study is that variables with more influence on  $\text{NO}_x$  are directly measured, and then having a reliable measurement is crucial for the model accuracy. In the second group,  $\text{NO}_{x,0}$ ,  $\dot{m}_{int}$  and  $\eta_v$  steady maps denote the most relevant maps to calibrate. The rest of variables have lower influence but not negligible.

#### 4.3.4.2 Dynamic results

Figure 4.11 already showed the ability of the  $\text{NO}_x$  model to fit the sensor output and the necessity of utilising the factor  $C_{dyn}$ . Figure 4.16 shows the model results with the SDMP test. This test is a good benchmark for testing the dynamic ability of the model as fast transients are performed. The influence of  $C_{dyn}$  is remarkable and can be seen in the left plot of the figure. Thin gray line corresponds to the model output but cancelling the effect of the thermal dynamic factor, i.e.  $C_{dyn} = 1$ , whilst the thick gray line is  $y_{\text{NO}_x}$ . If using  $C_{dyn} = 1$ ,  $\text{NO}_x$  is overestimated (about 20% higher than expected) due to the model does not consider the rapid load variations and the magnitude of them. In the SDMP, the influence of  $C_{dyn}$  is of crucial importance as the profile is ranging from low to high loads (thermal transients are critical). The right plot compares  $y_{\text{NO}_x}$  with  $z_{\text{NO}_x}$  obtaining a good fitting indeed when using ppm as output. The mean error in the SDMP tests (B5 and B6) is in the order of 50 ppm, which can be minimised by using online observers.



**Figure 4.16.** Left:  $y_{\text{NO}_x}$  calculated normally (thick gray) and cancelling thermal dynamic factor; i.e.  $C_{\text{dyn}} = 1$  (thin gray), and  $z_{\text{NO}_x}$  (black) during a part of the SDMP B5. Right:  $y_{\text{NO}_x}$  against  $z_{\text{NO}_x}$ .



**Figure 4.17.** Left:  $y_{\text{NO}_x}$  (gray) and  $z_{\text{NO}_x}$  (black) in the TRAN A ( $T_{\text{cool}} = 75^\circ\text{C}$ ,  $\Delta u_{\text{egr}} = 5\%$ ,  $\Delta p_{\text{boost}} = 0$ ). Right:  $x_{\text{NO}_x}$  (dashed black),  $y_{\text{NO}_x}$  (thin gray) and  $z_{\text{NO}_x}$  (black) in the same test.  $x_{\text{NO}_x}$  is the actual  $\text{NO}_x$ , no delay nor filtering are expected.

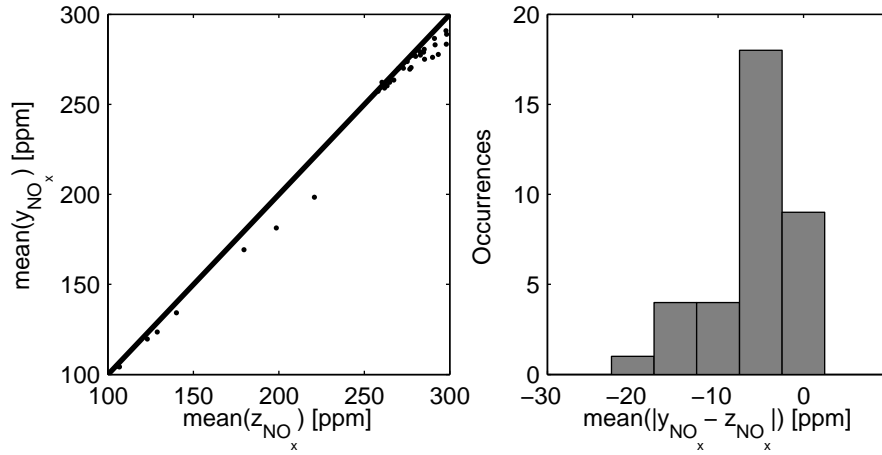
Figure 4.17 shows results on the TRAN A ( $T_{cool} = 75^\circ C$ ,  $\Delta u_{egr} = 5\%$ ,  $\Delta p_{boost} = 0$ ) cycle; the left plot compares  $y_{NO_x}$  with  $z_{NO_x}$  while the right plot includes the  $x_{NO_x}$  estimation. The sensor model fits well enough to have a reliable  $x_{NO_x}$ , that is observed by cancelling sensor filtering and lag. Even though the NO<sub>x</sub> model is based on quasi-static maps, system and physics dynamics are contained in the model inputs, which are directly measured from sensors. As shown in this Figure, the results are also promising.

The absolute mean value for  $z_{NO_x}$  and  $y_{NO_x}$  is calculated for every FTP cycle

$$mean(y_{NO_x}) = \frac{1}{l_{tran}} \sum_{k=1}^{l_{tran}} (abs(y_{NO_x}(k))) \quad (4.43a)$$

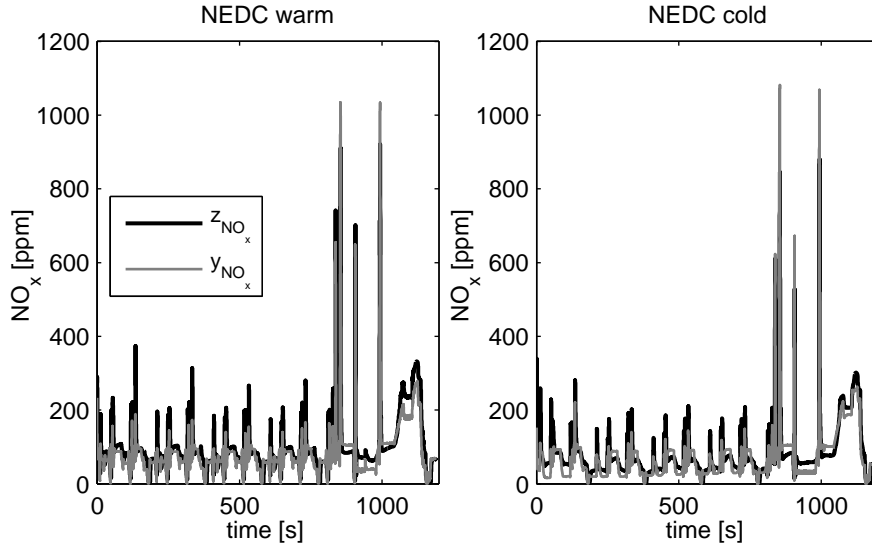
$$mean(z_{NO_x}) = \frac{1}{l_{tran}} \sum_{k=1}^{l_{tran}} (abs(z_{NO_x}(k))) \quad (4.43b)$$

where  $l_{tran}$  is the total number of samples for every TRAN cycle. Results are plotted in Figure 4.18, where the left plot shows  $mean(y_{NO_x})$  against  $mean(z_{NO_x})$  while the right plot shows the error distributions in ppm. The results demonstrate the model accuracy.



**Figure 4.18.** Left: mean value of NO<sub>x</sub> emissions, comparing the sensor  $z_{NO_x}$  against the model  $y_{NO_x}$  by simulating the NO<sub>x</sub> model in all the TRAN cycles. Right: absolute error bar distribution of the model against the sensor mean ( $|y_{NO_x} - z_{NO_x}|$ ). The mean absolute error is less than 10%.

Finally, Figure 4.19 shows the results with the NEDC A1 (warm) and A2 (cold). Especially during the highway part, there exists a bias in both



**Figure 4.19.**  $y_{\text{NO}_x}$  model output and  $z_{\text{NO}_x}$  measured by the sensor in the NEDC. Left: NEDC warm cycle A1. Right: cold cycle A2.

cycles. Since the model is optimised for a set of steady-state tests, dispersion effects or errors due to non-modelled variables ( $\text{NO}_x$  is highly affected by the ambient conditions) can affect the model accuracy. Admittedly, the model dynamics are able to reproduce  $z_{\text{NO}_x}$  and then the drift can be corrected by using observers as discussed in subsequent chapters.

## 4.4 Conclusions

An ECU oriented  $\text{NO}_x$  model has been developed by considering that variations of  $[\text{O}_2]_{\text{int}}$  cause exponential changes in the nominal  $\text{NO}_x$ , which is conveniently mapped. A study on different alternatives for estimating  $[\text{O}_2]_{\text{int}}$  is presented in the chapter, selecting the inert gas rate  $EGR\lambda^{-1}$  as the most convenient. A set of corrections and factors with respect to nominal operating points are included for considering the most sensitive parameters. Furthermore, a thermal dynamic model for simulating thermal loading is proposed as function of the product  $\lambda^{-1}n$ . Two additional models are needed for the  $\text{NO}_x$  model: an EGR gray box model for estimating  $EGR$  and a model for estimating the fuel-to-air ratio  $\lambda^{-1}$ .

A total of 363 steady-state tests are performed for calibrating all parameters and maps. Step-like tests for system and sensors identifications are made and different cycles are tested in order to validate the model. The  $\text{NO}_x$  prediction is successful and the simple programming as well as the fact that all required signals are available in commercial ECUs makes the model suitable for online engine control or diagnosis.

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# Chapter 5

## Adaptive observers for the dynamic estimation of engine variables

### Contents

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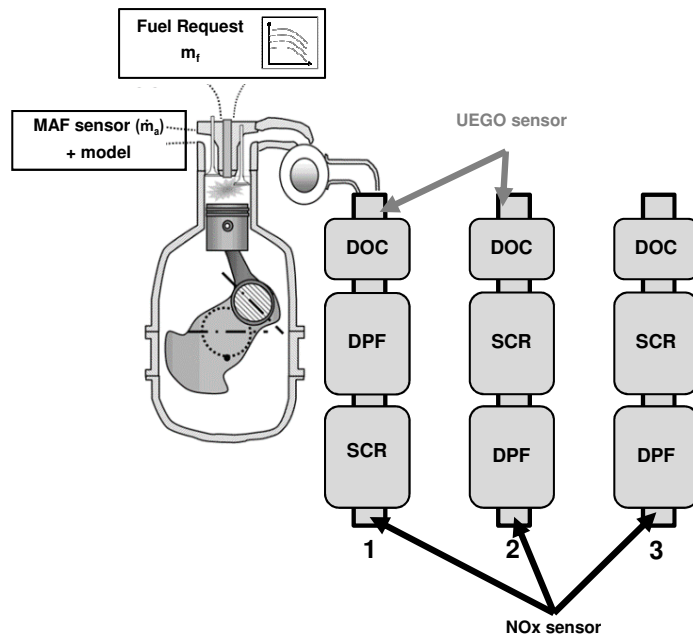
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## 5.1 Introduction

Along the previous chapters, different sensors and models have been explored in order to estimate engine variables, with focus on the exhaust  $\lambda$  and  $\text{NO}_x$ . From all the possible problems in the signals, the dynamic filtering, delay and drift are noticeable from sensor and model outputs. These issues are affected not only by sensors and models themselves, but by the engine layout.

**Sensor layout.** Figure 5.1 shows three possible layouts for on-board gas concentration sensors in order to estimate the engine-out  $\text{NO}_x$  and  $\lambda^{-1}$  for OBD and control of the AT devices. The final implantation depends heavily on the AT configuration and a variety of solutions were already discussed in the Chapter 2. Manufacturers are struggling for finding the configuration of systems and sensors that optimises the trade-off between low emissions, efficient torque production, low cost and robustness.



**Figure 5.1.** Possible UEGO and  $\text{NO}_x$  sensors configuration setup for a turbocharged diesel engine.

The usual position of UEGO sensors is directly downstream of the turbine: this way the sensor is near of the source and avoids the large range of pressure



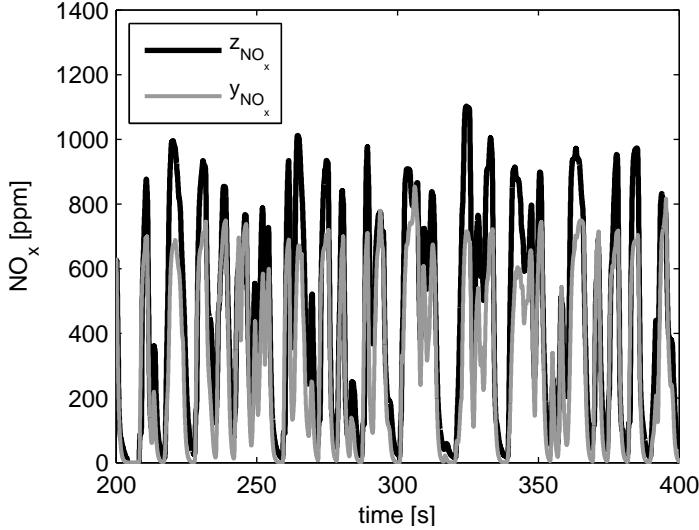
variation that exists in the exhaust manifold. However, if a SCR is installed, a  $\text{NO}_x$  sensor is needed for their control. Such sensor might be placed downstream of the AT devices in order to measure tail-pipe emissions, as sketched in the options 1, 2 and 3 of Figure 5.1. Hence it seems possible to avoid the UEGO sensor and just rely on the  $\text{NO}_x$  sensor information, which measures  $\lambda^{-1}$  and  $\text{NO}_x$ , as depicted in the third exhaust line of Figure 5.1. However, the slow response of the  $\text{ZrO}_2$  sensor, if compared with that of a UEGO sensor, is poor for RT purposes as was shown in Figure 2.5.

**The drift problem on models.** Models can be used instead for estimating variables but despite of the promising results with the models presented in Chapter 4, two problems can be underlined. On one hand, the model accuracy is driven by the collection of the appropriate data and calibration of all the parameters. This is a hard and time consuming task. In fact, the ECU has a big number of maps and parameters for the engine and vehicle management. On the other hand and independently of how well the model has been calibrated, there is inevitably a drift between the system and the model as the surrounding conditions change and the engine ages (this occurs with all kind of models). DDMs are highly sensitive to the calibration data set and will have problems with ageing, manufacturing discrepancies, slowly varying parameters and other non-modelled variables.

Figure 4.19 showed the  $\text{NO}_x$  model (filtered) output  $y_{\text{NO}_x}$  for two NEDC cycles (warm A1 and cold A2) and a clear drift exists, especially during the highway part in the A1 and during the whole cycle in A2. Another example is shown in Figure 5.2, where an old and non-optimised calibration set for the model (maps and sensor dynamics) is used for simulating the SDMP B6 test. The dynamics are well caught but a clear drift also exists. In this case it is not only ageing that affects, but test or ambient conditions.

**The use of adaptive strategies.** The dynamic problems of the sensors and the drift on models motivate the design of adaptive on-board strategies. Concretely, two kind of models and algorithms are proposed in order to solve these issues, which are:

- augmented models for drift correction and dynamic estimation of variables, and
- learning algorithms for updating look-up tables and parameters on the basis of COMs, which can be used for calibration and/or online adaptation.



**Figure 5.2.** Model output  $y_{NO_x}$  and  $NO_x$  sensor signal  $z_{NO_x}$  in the SDMP B6 test with a non-optimized calibration data set. Even though the dynamics are well caught,  $y_{NO_x}$  is biased with respect to  $z_{NO_x}$ .

In the next, the dissertation focuses on the description of the proposed methods, while Chapter 6 will give the results with experimental data.

## 5.2 Augmented models for drift correction

The design of augmented models for bias tracking and the implementation of a KF for observing this bias is discussed. A known case is considering a fast model that keeps high frequency components of the considered variable while a slow but steady-state accurate sensor permits to correct the model drift. This problem is solved by observing the bias  $\theta$  between a reference measurement  $z$  and an initial estimation  $u$ , considering the measurement dynamics and delay.

### 5.2.1 Drift correction algorithm

The model is presented in the discrete state-space form. The sample time is considered constant and the model equation is defined as follows

$$x_k = \mathbf{F}x_{k-1} + Bu_k + \mathbf{W}_k \quad (5.1a)$$

$$z_k = Hx_k + v_k \quad (5.1b)$$

Without loss of generality, the problem can be stated for estimating the drift  $\theta_k \in \mathbb{R}$ , augmenting the standard model with such state.  $u_k \in \mathbb{R}$  is the state-space input, and corresponds to a model output with a certain delay  $\tau$ . The effect of the delay over the augmented model observation is discussed in Chapter 6. For the models presented in Chapter 4, the vector  $u$  is as follows

$$u_k^{\text{NO}_x} = x_{\text{NO}_x}(k - \tau_{\text{NO}_x}/T_s) \quad (5.2a)$$

$$u_k^{\lambda^{-1}} = x_{\lambda^{-1}}(k - \tau_{\lambda^{-1}}/T_s) \quad (5.2b)$$

depending if  $\text{NO}_x$  or  $\lambda^{-1}$  is the key variable. The sensor signal  $z_k \in \mathbb{R}$  the output, while the first order filter presented in Equation (3.3) with response time  $a$  is used for modelling  $z_k$ , giving the signal  $x_{fk} \in \mathbb{R}$ . The state vector  $x_k \in \mathbb{R}^2$  is built as follows

$$x_k = [\theta_k \quad x_{fk}]^T \quad (5.3)$$

$\mathbf{F}$ ,  $B$  and  $H$  are the constant state-space matrices

$$\mathbf{F} = \begin{bmatrix} 1 & 0 \\ 1 - a & a \end{bmatrix} \quad B = \begin{bmatrix} 0 \\ 1 - a \end{bmatrix} \quad H = [0 \quad 1] \quad (5.4)$$

Noises  $W_k \in \mathbb{R}^2$  with  $w_k \in \mathbb{R}$

$$W_k = [w_k \quad 0]^T \quad (5.5)$$

and  $v_k \in \mathbb{R}$ , from process and output respectively, correspond to Gaussian processes with the following stationary constant covariance matrices

$$\mathbf{Q} = \text{cov}[ww^T] = \begin{bmatrix} \sigma_w^2 & 0 \\ 0 & 0 \end{bmatrix} \quad (5.6a)$$

$$r = \text{cov}[vv^T] = \sigma_v^2 \quad (5.6b)$$

The system is fully observable and linear time invariant (LTI). Note that only a noise to the state  $\theta$  is applied as far as this is the key observed variable and the noise is directly transmitted to  $x_f$ . Applying noise to  $x_f$  adds an extra degree of freedom that does not necessarily improve the identification. The system can easily be augmented with other key variables if necessary. In such sense, the state vector  $X$  (4.18) and the input vector  $U$  (4.19) from the RT  $\text{NO}_x$  model could be included in the vectors of the state-space model (5.1) for a compact programming.

**The Kalman filter for drift observation.** The optimal solution of the estimation problem was addressed by Kalman [1], and has been applied by different authors to the automotive domain (see for example [2–5]). In a first step, the state vector is predicted considering the system input and their expected dynamic characteristics and in a second step, this is updated by calculating the estimation error  $e_k$  and weighting it by means of the KF gain  $K_k$

$$\hat{x}_{k|k-1} = \mathbf{F}\hat{x}_{k-1} + Bu_k \quad (5.7a)$$

$$e_k = z_k - H\hat{x}_{k|k-1} \quad (5.7b)$$

$$\hat{x}_k = \hat{x}_{k|k-1} + K_k e_k \quad (5.7c)$$

Kalman gain  $K_k$  was solved in the Equation (2.13).

**The steady-state Kalman filter.** Since current application considers a linear time invariant (LTI) and a fully observable system, the filter is steady-state [6]. A constant Kalman gain  $K_\infty$

$$K_\infty = [k_1 \quad k_2]^T \quad (5.8)$$

is then considered and calculated offline

$$K_\infty(\sigma_w^2/\sigma_v^2, a) = \lim_{k \rightarrow +\infty} K_k \quad (5.9)$$

The use of a constant  $K_\infty$  leads to a sub-optimal filter neglecting the  $K$  updating [7] but alleviating the computational burden, since all the iterative process for obtaining  $K_\infty$  is performed offline. Anyway, if a more complex sensor model (including the effect of the operating conditions on the sensor dynamics) were employed, the Kalman gain would vary with time, and then  $K_k$  might be approximated online by means of the steady-state solution of the Riccati Equation (2.13):

$$k_1 = \sqrt{(1 - k_2) \frac{\sigma_w^2}{\sigma_v^2}} \quad (5.10a)$$

$$k_2 = \frac{s}{s + 1} \quad (5.10b)$$

with

$$s = (1 - a) \sqrt{\frac{\sigma_w^2/\sigma_v^2}{1 - k_2}} (1 + a - ak_2) + a^2 k_2 \quad (5.11)$$

which is demonstrated in the Appendix 5.A.2.  $k_2$  is solved by a fixed point iteration, and  $k_1$  by direct substitution of  $k_2$ .

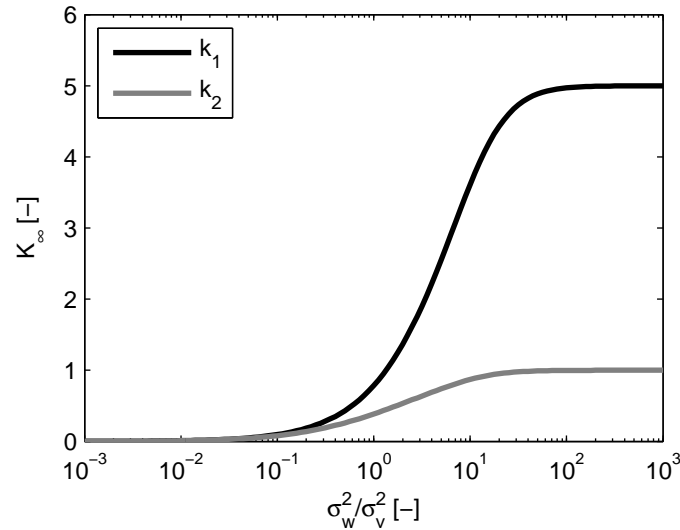
**Final estimation.** The final estimation  $\hat{x}$  on a delayed basis is computed by adding the observed bias to the input  $u$  (see (5.2))

$$\hat{x}(k - \tau/T_s) = u_k + \hat{\theta}_k \quad (5.12)$$

This equation might be included in the state-space model (see Section 6.3.3); the online estimation solving the delay problem is discussed later (see Sections 6.2.3 and 6.3.3).

### 5.2.2 Observer tuning

The KF is tuned with the adequate selection of the quotient  $\sigma_w^2/\sigma_v^2$  that affects the model and sensor respectively, defining an optimal  $K_\infty$ . Figure 5.3 shows the different values of the two elements of  $K_\infty$  for different values of  $\sigma_w^2/\sigma_v^2$ .  $k_1$  is the bias gain, while  $k_2$  corrects the sensor model. Note that the lower the sensor noise is considered (higher  $\sigma_w^2/\sigma_v^2$ ), the faster the states' estimations are updated (because sensor measurement is propagated to the states). The filter behaves as a low pass filter for the sensor and a high pass filter for the model.



**Figure 5.3.** Evolution of the elements of  $K_\infty$  gain vector when varying  $\sigma_w^2/\sigma_v^2$ . Coordinate axis is in logarithmic scale and  $a = 0.8$ .

Hence the filter tuning defines how fast the bias is cancelled, but two issues must be considered:

- The filter also rejects the sensor noise. Using a very high Kalman gain implies do not filtering the measurement noise.
- If the sensor dynamics characteristics are not perfectly known, a high value of the gain causes the system to rely on incorrect information, and then peaks appear in the estimation especially during transients.

For illustrating the latter issue, the augmented model (5.1) is simulated by using synthetic signals. A step transition in the objective variable, named  $x_r \in \mathbb{R}$  is produced by defining the following parameters for the sensor model

$$a' = a \pm \sigma_a \quad \tau' = \pm \sigma_\tau$$

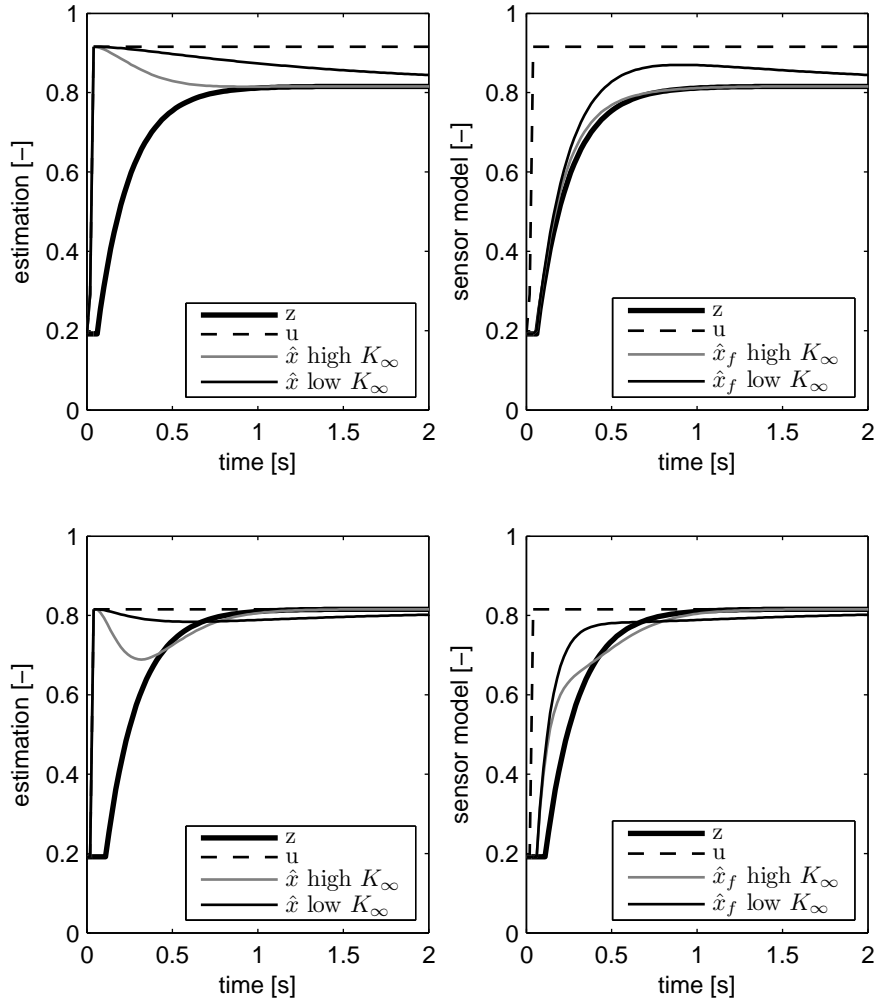
with an initial drift  $\theta^0$  and a final drift after the step  $\theta^1$ . Since synthetic signals are used, filter performance can be evaluated through the comparison of  $\hat{x}$  with actual  $x_r$  (which in real applications it is not usually known). The results can be seen in Figure 5.4.

The basic cases correspond to:

1. Sensor perfectly known ( $a' = a, \tau' = \tau$ ) with an step in the bias ( $\theta^0 \neq \theta^1$ ). In this case, the higher  $K_\infty$  is, the faster the bias is corrected, as depicted in top plots of Figure 5.4.
2. Constant bias and uncertainty on the sensor description ( $a' \neq a, \tau' \neq \tau$ ). In this case, higher  $K_\infty$  values yield to rely excessively on the poorly known sensor dynamics, creating artificial peaks, as shown in bottom plots of Figure 5.4.

The general error case in the sensor modelling is a (non-linear) superposition of the previous cases, and it is clear that a trade-off in the selection of  $K_\infty$  must be considered according to the signals uncertainties and the required convergence speed. For the sensors considered in this work, sometimes they exhibit a significant variation in their dynamic properties and a consistent model cannot be derived for the whole engine operating range. This would force a low value of  $K_\infty$  (and hence a slow correction of the model bias), but several modifications are proposed in the implementation.

Even though data based methods could be used for estimating the appropriate noise trade-off [8], trial-and-error tuning is usual when working with KF. A Monte Carlo based method application presented in [9] is proposed for tuning the filter, where the uncertainties on the sensor knowledge, expected



**Figure 5.4.** Top plots: Observation of the augmented model when the sensor model is perfectly known ( $a' = a, \tau' = \tau$ ) with an step in  $\theta$ . Bottom plots: Observation of the augmented model when the sensor model is not perfectly known ( $a' \neq a, \tau' \neq \tau$ ) and with a constant  $\theta$ . The left plots show the model estimation  $\hat{x}$  when the sensor performance is perfectly known and for two different  $\sigma_w^2/\sigma_v^2$  quotients, i.e. two different values for  $K_\infty$ . The right plots show the sensor model estimation  $\hat{x}_f$  for these two  $K_\infty$  values.

working points ( $n$  and  $m_f$  values) and different expected  $\theta$  are used for evaluating the filter performance under different  $\sigma_w^2/\sigma_v^2$  values. These variations are considered with statistical distributions, and the configuration which minimises the total error is selected.

### 5.3 Learning algorithms for updating look-up tables

Look-up tables are often used in automotive in order to model systems with complex (and highly non-linear) expressions in a grid basis. The table elements are used for interpolating in order to generate the outputs. They can be included in a model, such as the NO<sub>x</sub> model presented in Chapter 4, or can be used directly for storing the bias  $\theta$  calculated in the previous section. The possibility of developing RT algorithms for updating tables makes possible to develop adaptive look-up tables. The methods based on the EKF are well-suited but require a heavy computation burden, topic that is especially treated in this section. In the following, the section describes the problem of updating look-up tables and presents three methods based on the EKF.

**Defining look-up tables.** A look-up table  $\Theta \in \mathbb{R}^{\prod_{i=1}^N n_i}$  is defined as a N-Dimensional mapping  $\{\Theta : \mathbb{R}^N \rightarrow \mathbb{R}\}$  represented by a grid in  $N \in \mathbb{Z}^+$  dimensions, where each one has  $n_i$  grid points. The mapping further relies on a multivariate interpolation  $q(\cdot)$  to calculate the function value from the input using the grid for the interpolation variables and  $\Theta$ . In automotive systems, the multivariate interpolation schemes are often linear in each one of the dimensions, and this is the case that will be considered here. Without loss of generality the presentation will use 2D tables as being the most frequently occurring dimensions, since working with other dimensions is only a matter of reducing or increasing indexes. Then, for N=2,

$$\Theta = [\Theta_{i,j}] \quad (5.13)$$

where  $i = 1, \dots, n_r$  and  $j = 1, \dots, n_c$  ( $r$  stands for *row* and  $c$  for *column*). The multivariate interpolation function for generating the output  $x_\Theta(k)$  from the input  $u_k = [u_1(k) \ u_2(k)]^T$  can be expressed as

$$x_\Theta(k) = \text{vec}(q(u_k))^T \text{vec}(\Theta) \quad (5.14)$$

$\text{vec}(\cdot)$  is the vector transformation and  $q(u_k)$  is the interpolation matrix that both selects the elements to be interpolated and contains the weights. In the 2D case  $q_k(u_k)$  selects the 4 ( $2^N$  in the ND case) active elements  $\Theta_{i,j}$ ,  $\Theta_{i,j+1}$ ,



$\Theta_{i+1,j}$ ,  $\Theta_{i+1,j+1}$  (with  $i, j$  fulfilling  $u_{1,k} \in [r_i, r_{i+1}]$  and  $u_{2,k} \in [c_j, c_{j+1}]$ ) and thus contains the following block with non-zero weights

$$\begin{bmatrix} q(u_k)_{i,j} & q(u_k)_{i,j+1} \\ q(u_k)_{i+1,j} & q(u_k)_{i+1,j+1} \end{bmatrix} = \begin{bmatrix} (1 - \eta_{1,k})(1 - \eta_{2,k}) & (1 - \eta_{1,k})\eta_{2,k} \\ \eta_{1,k}(1 - \eta_{2,k}) & \eta_{1,k}\eta_{2,k} \end{bmatrix} \quad (5.15)$$

where

$$\eta_{1,k} = \frac{u_{1,k} - r_i}{r_{i+1} - r_i}, \quad \eta_{2,k} = \frac{u_{2,k} - c_j}{c_{j+1} - c_j} \quad (5.16)$$

This Section analyses the computational aspects of updating look-up tables with the KF. Two algorithms are designed based on the KF for a computationally efficient table updating: a simplified KF (SKF) which manipulates the covariance matrix  $\mathbf{P}$  and the associated updates efficiently, and a steady-state approach for the KF (SSKF) which directly neglects covariance information. Both methods are inspired by the works presented in [10–12].

**Modelling for learning, drift or ageing in tables.** In the setting,  $\Theta$  models a nonlinear function and the interesting aspect is to allow the model to adapt to the system to either learn the system and/or follow the ageing of the system. This is modelled in the standard way as a random walk process, where the table parameters are collected in a state vector  $x_k \in \mathbb{R}^{n_r \times n_c}$ . The full model can be written as follows

$$x_k = x_{k-1} + w_k \quad (5.17a)$$

$$z_k = \text{vec}(q(u_k))^T x_k + v_k \quad (5.17b)$$

noting that no uncertainties are allowed in the interpolation variable  $u_k$ . For convenience, the non-zero elements in  $\text{vec}(q(u_k))$  are denoted  $q_k^o \in \mathbb{R}^{1 \times 4}$  ( $o$  stands for *observable*) and the corresponding elements in the state vector,  $x_k^o \in \mathbb{R}^4$  for the 2D case. Then, the following expression is obtained for the output

$$z_k = q_k^o x_k^o + v_k \quad (5.18a)$$

$$q_k^o = [(1 - \eta_{1,k})(1 - \eta_{2,k}) \quad \eta_{1,k}(1 - \eta_{2,k}) \quad (1 - \eta_{1,k})\eta_{2,k} \quad \eta_{1,k}\eta_{2,k}] \quad (5.18b)$$

where  $\mathbf{Q}$  and  $\mathbf{R}$  are

$$\mathbf{Q} = \sigma_w^2 \mathbf{I}_{n_r \times n_c} \quad (5.19a)$$

$$\mathbf{R} = \sigma_v^2 \quad (5.19b)$$

and  $\mathbf{I}_{n_r \times n_c} \in \mathbb{R}^{n_r \times n_c}$  is the identity square matrix. If there is application knowledge available,  $\sigma_w^2$  might be selected individually for each table element

building a vector  $\Sigma_w^2 = [\sigma_{w,1}^2 \dots \sigma_{w,n_r \times n_c}^2]$  that contains individual variances in such way  $\mathbf{Q} = \Sigma_w^2 \mathbf{I}_{n_r \times n_c}$ . In automotive, this is useful for considering the order of magnitude of the parameters, i.e. the absolute error will not be the same for low emissions at lower loads than highest peaks at higher loads, and on the other hand, if a driving pattern exists, the noise could be mapped over the table grid. This is linked with the foreseen probability that the engine is running at a certain operating condition, i.e. the ageing of the elements corresponding to more frequent areas will be lower and the opposite.

### 5.3.1 The extended Kalman Filter, KF

The EKF (along the text and figures, the acronym KF refers to both EKF and KF) can now be used to observe  $x_k$ , when measurements  $z_k$  are given. At every iteration, the table is updated by

$$\hat{x}_k = \hat{x}_{k-1} + K_k(z_k - \text{vec}(q(u_k)))^T \hat{x}_k \quad (5.20)$$

where  $K_k$  is calculated using Equation (2.13),  $H_k = \text{vec}(q(u_k))^T$  varies with the operating point conditions. Although only the active elements  $x_k^o$  are updated at every  $k$ , all  $\mathbf{P}$  elements enter in the equation, which leads to huge calculations and big required memory resources. This makes difficult the implementation in commercial ECUs.

There are several publications on computational aspects of Kalman filters. Some authors have studied the filter optimization when different nonlinear functions are handled as for instance [13–15]; where the latter makes an interesting study on the total number of operations required for the updating phase. Chandrasekar *et al.* [16] present an interesting methodology when the system order is extremely large by using the finite-horizon optimization technique for obtaining reduced-order systems. But the key observation here is that the look-up table estimation has structural properties that can be exploited to reduce the computational and memory requirements significantly in a simple way, as suggested by Vogt *et al.* [11].

To understand the problem, the covariance matrix  $\mathbf{P}$  in the KF is studied. For instance, a 2D look-up table with a size  $n_r n_c$  needs system  $\mathbf{F}$  and covariance matrices  $\mathbf{P}$  of  $(n_r n_c) \times (n_r n_c)$ . However, only the active elements are influenced during the update and by defining a local observable system (which corresponds to the active elements), one receives a system which is  $4 \times 4$  against  $(n_r n_c) \times (n_r n_c)$ . That means that the non-active elements do not affect the updating until they become active (zero values for the  $K$  related elements). Furthermore,  $\mathbf{P}$  is a symmetrical and positive-semidefinite matrix, which allows further simplifications in the KF calculations.

**The local observable system.** The *Local Observable System* is defined and analysed and a computationally efficient approximation for the KF is described. Following with the 2D look-up table application, at every iteration a maximum number of 4 elements can be updated and then, if no dynamics are considered, the general  $n_r n_c$  system (5.17) can be simplified to

$$x_k^o = \mathbf{I}_{4 \times 4} x_{k-1}^o + w_k \quad (5.21a)$$

$$z_k = q_k^o x_k^o + v_k \quad (5.21b)$$

where  $\mathbf{I}_4 \in \mathbb{R}^{4 \times 4}$  is the unitary matrix and  $x_k^o \in \mathbb{R}^4$  is the state-vector that stores the active states at every  $k$ . The simplest 2D map is a table of only 4 elements, and where the one existing area is always active, i.e. the system (5.21) is exactly (5.17). For the general case, (5.21) must be rewritten at every  $k$  as elements and matrices change.

**About the system observability.** The system (5.17) is not fully observable in one iteration. But the local observability is ensured if the parameters or states in the local system (5.21) converge with a given data set. As the system is not LTI, the ordinary observability rank condition [17] is not directly applicable.  $q_k^o$  depends on the input data, and it is evident that if a enough level of excitation is given, then the system could be observed, whatever the method chosen. This minimum level of excitation could be proved if 4 elements have been excited. A sufficient observability matrix for local observability may be built with the first 4 independent observations not necessarily consecutive in instants  $i_1, i_2, i_3$  and  $i_4$  of the elements of the involved area:

$$\mathbf{O} = [q_{i_1}^o \quad q_{i_2}^o \quad q_{i_3}^o \quad q_{i_4}^o]^T \quad (5.22)$$

and if the rank of this matrix is 4, then the system (5.21) might be observable. But here the problem is linked with the noise tuning and indeed full rank  $\mathbf{O}$  does not lead to system full convergence (full observability does not lead to full convergence as the method and the model structure affect). The most of the elements are also included in other neighbour areas and because of this, the minimum condition of observability of a given state depends on the number of independent measurements that affects every state. The condition stated in [10] gives a general condition.

### 5.3.2 The steady-state KF approach, SSKF

If the problem is considered stationary (the same input is repeated once and again), then  $q_k^o = q^o$  and the filter is steady-state. In addition to this and

given a noise ratio  $\sigma_v^2/\sigma_w^2$ ,  $K_k$  in (2.13) converges after a certain number of iterations as expressed in (5.9). That occurs although  $\mathbf{P}$  does not converge and can be ill conditioned. Then, the stationary equivalent filter is obtained and an analytical expression is derived for the *local observable system*, resulting a Kalman gain  $K^{SS} \in \mathbb{R}^4$

$$K^{SS} = \begin{bmatrix} k_i(\eta_1, \sigma_v^2/\sigma_w^2) \cdot k_i(\eta_2, \sigma_v^2/\sigma_w^2) \\ k_i(\eta_1, \sigma_v^2/\sigma_w^2) \cdot k_i(1 - \eta_2, \sigma_v^2/\sigma_w^2) \\ k_i(1 - \eta_1, \sigma_v^2/\sigma_w^2) \cdot k_i(\eta_2, \sigma_v^2/\sigma_w^2) \\ k_i(1 - \eta_1, \sigma_v^2/\sigma_w^2) \cdot k_i(1 - \eta_2, \sigma_v^2/\sigma_w^2) \end{bmatrix} \quad (5.23)$$

where  $k_i$  is computed as follows

$$k_i(\eta, \sigma_v^2/\sigma_w^2) = \frac{0.5(1 - \eta)(1 + s)}{0.5(1 + s)(1 - 2\eta + 2\eta^2) + \sigma_v^2/\sigma_w^2} \quad (5.24a)$$

$$s = \sqrt{1 + \frac{4}{(1 - 2\eta + 2\eta^2)} \frac{\sigma_v^2}{\sigma_w^2}} \quad (5.24b)$$

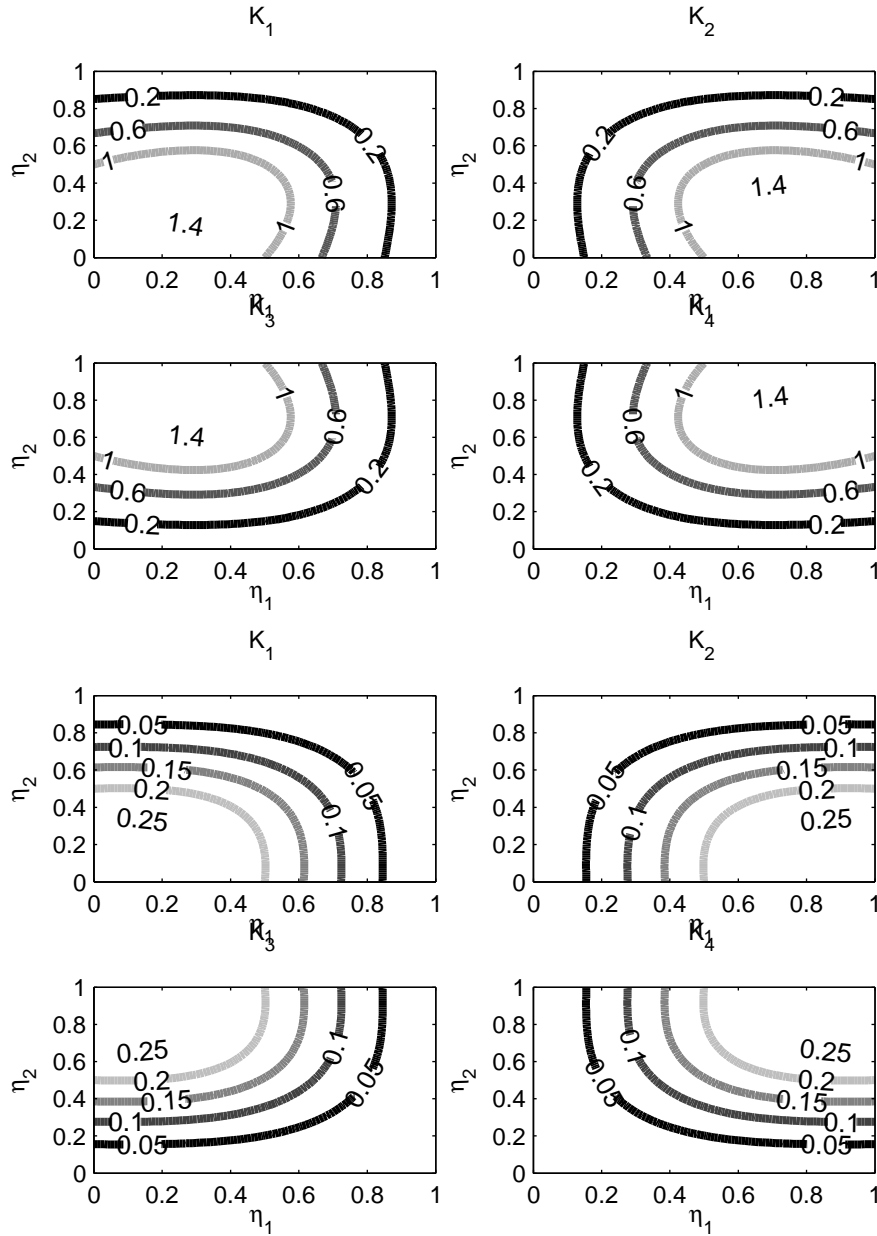
These expressions are solved in Appendix 5.A.2.

The user can evaluate (5.23) for the 2D case for a given  $\sigma_v^2/\sigma_w^2$  or can simulate the functions off-line for mapping the values and reduce the computing time, which is the option suggested by the authors. Figure 5.5 shows the values of the  $K^{SS}$  elements for two different  $\sigma_v^2/\sigma_w^2$  quotients. The symmetry of the functions is clear when comparing the elements of  $K^{SS}$ . Both the noise tuning and the relative position of the operating points affect to the estimation. On one hand, as the operating point approaches to one grid node, the gain is closer to the maximum (it is not linear) and the opposite. On the other hand, a higher  $\sigma_v^2/\sigma_w^2$  leads to a higher correction. Non-active elements have a null correction, i.e. their related gain is zero.

This approach allows to map  $K^{SS}$  and this only depends on the inputs ( $\eta_1$  and  $\eta_2$  from (5.16)) and the noise trade-off  $\sigma_v^2/\sigma_w^2$ . Nevertheless, as the parameters ageing is not considered nor the estimation error computed, the algorithm is quite fast and light but the robustness must be assessed in the applications. The pseudo-code of the algorithms is in the Appendix 5.B. The next subsection presents the SKF algorithm which tracks the estimation error in an efficient way.

### 5.3.3 The simplified Kalman filter, SKF

Despite that only the associated active elements affect  $K_k$ , all  $\mathbf{P}_k$  elements are predicted and updated at every  $k$ .  $\mathbf{P}_k$  can be reorganised in such way that



**Figure 5.5.** Membership functions for Kalman gain correction for the 2D case for  $\sigma_v^2/\sigma_w^2 = 0$  (top) and  $\sigma_v^2/\sigma_w^2 = 0.1$  (bottom).

variances related with observable  $\mathbf{P}_k^o$  and unobservable  $\mathbf{P}_k^u$  elements are split.

$$\mathbf{P}_k = \left[ \begin{array}{c|c} \mathbf{P}_k^u & \begin{matrix} \cdot \\ \cdot \\ \cdot \end{matrix} \\ \hline \begin{matrix} \cdot & \cdot & \cdot \end{matrix} & \mathbf{P}_k^o \end{array} \right] \quad (5.25)$$

$K_k$ ,  $q_k$  and  $\mathbf{Q}$  can also be split in the same way. Iterating one step ahead (2.13c) is written

$$\begin{aligned} \mathbf{P}_{k+1} &= \left[ \begin{array}{c|c} \mathbf{I} & \mathbf{0} \\ \hline \mathbf{0} & \mathbf{I}_{4 \times 4} - K_k^o q_k^o \end{array} \right] \left[ \begin{array}{c|c} \mathbf{P}_k^u + \mathbf{Q}^u & \begin{matrix} \cdot \\ \cdot \\ \cdot \end{matrix} \\ \hline \begin{matrix} \cdot & \cdot & \cdot \end{matrix} & \mathbf{P}_k^o + \mathbf{Q}^o \end{array} \right] \\ &= \left[ \begin{array}{c|c} \mathbf{P}_k^u + \mathbf{Q}^u & \begin{matrix} \cdot \\ \cdot \\ \cdot \end{matrix} \\ \hline \begin{matrix} \cdot & \cdot & \cdot \end{matrix} & (\mathbf{I}_4 - K_k^o q_k^o)(\mathbf{P}_k^o + \mathbf{Q}^o) \end{array} \right] \end{aligned} \quad (5.26)$$

where all observable sub-matrices are 4x4, while the unobservable ones are  $(n_r n_c - 4) \times (n_r n_c - 4)$ . (5.26) shows how the diagonal elements of  $\mathbf{P}_k^u$  are only affected by adding the diagonal matrix  $\mathbf{Q}^u$ . However, both diagonal and non-diagonal elements of  $\mathbf{P}_k^o$  are affected as in a 4x4 KF. The crossed relationships between the observable and the unobservable system, shown as dots in the matrix, are kept constant. This allows simplifying the complete system to the 4x4 active state-space system, whose resolution is highly computationally efficient.

The current work develops an approximation of the KF that requires both less memory and computations in the iterations and is named SKF. The SKF builds and solves the local observable system (5.21) at every  $k$ . Some simplifications are assumed: for the non-active elements, whose covariance matrix is  $\mathbf{P}^u$ , non-diagonal elements are neglected building a vector  $P^u = \text{diag}(\mathbf{P}^u)$ . This is justified for two reasons. First, when the system is running, then leaves one area and later returns to it, related variances reflect the time that the system has been out of this area. This is a desired property since the ageing is captured by the variance increase due to  $\mathbf{Q}^u$ , which indeed for this application it is defined as a diagonal matrix. Older crossed correlations are maintained over the time and whether or not they are valid is not sure. Second, the simulation results show that the full EKF and the SKF have similar performances when running on synthetic data and real data from the engine application.

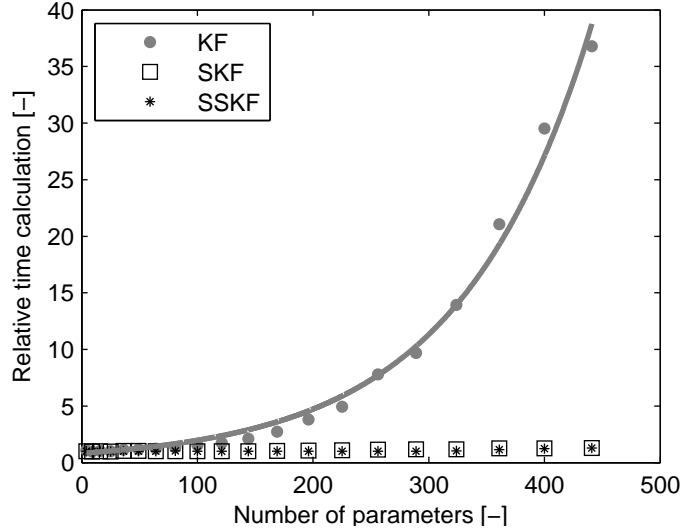
Then, at every iteration, the local system (5.21) is solved with the only additional operation of  $P_k^u + Q^u$  (converted to a vector calculation). The SKF is a suboptimal filter but with a similar behaviour than KF, as demonstrated in Sections 5.3.4.

The pseudo-code for the SKF implementation is in Algorithm 5.B. Note that the code is to show the algorithm where some variables are not strictly needed in the computer programming, e.g.  $P_k^u$  always that  $P_k$  exists, but these steps are kept for the sake of clarity. The noise variance  $\sigma_w^2$  is selected equal and constant for all elements, allowing to pre-define a diagonal matrix  $\mathbf{Q}^o = \sigma_w^2 \mathbf{I}_{4 \times 4}$  for the active system and a vector  $Q^u = \sigma_w^2 \text{diag}(\mathbf{I}_{n_r \times n_c - 4})$  for the non-active one.

In the next, the memory allocation and computational complexity of the three methods, full EKF (here denoted KF), the SKF as well as the SSKF are described.

**Analysis of the memory and computations.** A computation time study is made programming code in the Matlab software on a laptop computer equipping Intel Core 2DUO T9300 2.5 GHz with Windows VISTA 64 bits. Figure 5.6 shows the relative computation time used by the three methods for updating 2D look-up tables of different sizes. Only square  $n \times n$  tables have been considered with  $n$  ranging from 2 to 21, since it is the total number of elements in  $\text{vec}(T)$  that influences memory allocation and computations. The Y-axis shows the relative computation time with respect to the averaged time that the SSKF takes for performing 1000 iterations, i.e., a relative time of 1 means that the method needs exactly the same time than the SSKF. The SKF needs around 1.15 times the SSKF calculation time, whatever the number of parameters. This slight difference is explained by the manipulation of variances and the need of redefining the active vector and matrices  $\mathbf{P}^o$ , and  $x^o$  when the active area changes (see lines 4 and 5 of the Algorithm 2 and compare with Algorithm 1, both included in Appendix 5.B). However, the required time for the KF rapidly grows when the number of parameters increases, i.e. for 256 parameters ( $16 \times 16$ ), the KF needs around 7.8 times the one for the SSKF. An exponential function may be fitted by the least squares method to the KF time as function of the number of parameters  $n_p$ , whose result is the relative time that the KF takes with respect to the one of the SSKF.

Table 5.1 summarises the required memory resources and computation times for the algorithms. The SKF permits to reduce the requirements in terms of memory resources and the general system  $(n_r n_c) \times (n_r n_c)$  is reduced



**Figure 5.6.** Relative computation time required for the methods for updating 2D look-up tables. Times are normalised with the averaged time that the SSKF needs for computing 1000 iterations of the filter.

to a simple 4x4. These results show the benefits when using the SSKF and the SKF compared to the full KF.

**Table 5.1.** Memory storage and computational burden comparison between the Kalman filter KF, the simplified Kalman filter SKF and the steady-state approach SSKF.

	<b>P</b>	<b>K</b>	Rel. Time
KF	$(n_r n_c) \times (n_r n_c)$	$(n_r \times n_c) \times 1$ (by (2.13))	$0.82 \exp(0.0087 n_p)$
SKF	$4 \times 4 + (n_r \times n_c) \times 1$	$4 \times 1$ (by (2.13))	1.15
SSKF	$4 \times 4$	$4 \times 1$ (by (5.23))	1

### 5.3.4 Simulation of the updating algorithms

The convergence and robustness of the approximations need also to be studied and this is first performed by simulations. The objective of this section is comparing the abilities and performance of the algorithms against different synthetic cases where no dynamics are accounted (the real system and the model are statics; Section 5.3.5 later discusses how to consider the dynamics).



For automotive applications, engine speed  $n$  and fuel mass injection quantity  $m_f$  usually define the engine operating point and the look-up table scheduling points selected in the simulation are inspired by these quantities. Hence  $r$  represents a grid for  $u_1(k)$  (the speed in rpm) and  $c$  for  $u_2(k)$  (the injected fuel mass in mg/str)

$$r = [500 \quad 1500 \quad 2500 \quad 3500] \quad (5.27a)$$

$$c = [0 \quad 20 \quad 40 \quad 60] \quad (5.27b)$$

and the true map (the objective map to update)  $\Theta_r$  is defined with the surface

$$\Theta_r = \begin{bmatrix} 1 & 1 & 1 & 1 \\ 1 & 1 & 1 & 1 \\ 1 & 1 & 1 & 1 \\ 1 & 1 & 1 & 1 \end{bmatrix} \quad (5.28)$$

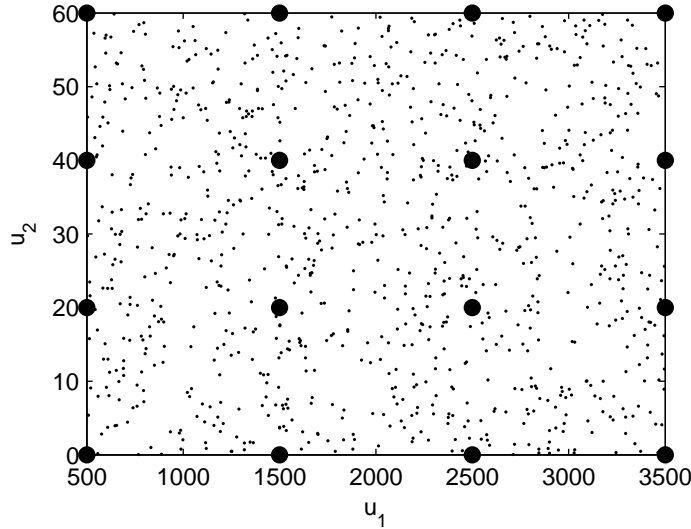
where  $x_r = \text{vec}(\Theta_r)$  is the expanded state vector objective of the estimation process. No initial process knowledge is considered:  $x_0, x_0^o, \mathbf{P}_0$  are initialised with zeros.  $\sigma_v^2/\sigma_w^2 = 1$  for all cases, unless it were pointed.

System measurements are given and the algorithms performance is evaluated. As synthetic signals are used, this allows defining error metrics. Simulations include two worst-cases and a favourable one:

- Input with random variation: a random shot of values following an uniform probability distribution. This is an ideal situation for learning as all areas are excited and the parameters ageing is low (the excitation is high). Measurements are perfect and constant, thus  $z_k = 1 \forall k$ .
- Linear variation: varying  $u_2$  keeping constant  $u_1$ . This variation tests a degenerated case where observability is critical along the  $u_1$ -dimension. Measurements are perfect and constant:  $z_k = 1 \forall k$ .
- Measurement noise rejection: Studying the effects of noise transmission between output and observation when the measurement vector of the previous case is noisy.

#### 5.3.4.1 Simulation 1: Input with random variation

The sequence  $\{u_k\}_{k=1}^{1000}$  following an uniform distribution over the grid defined in (5.27) is used for exciting the system. A complete identification is possible as the grid is completely covered. This situation is not so far of the

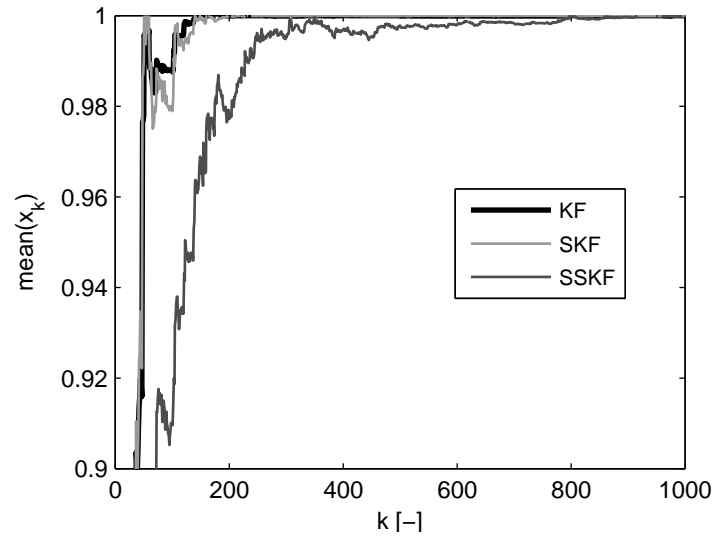


**Figure 5.7.** Random variation of inputs  $u_1$  and  $u_2$  shown with small dots and table grid with thick dots.

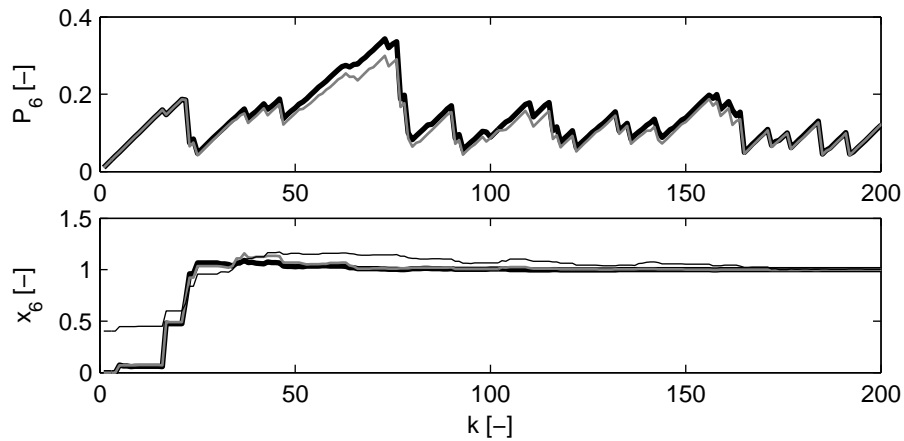
reality for diesel engines, e.g. urban cycles with a lot of speeding/braking actions cause quasi-random variations in partial-low load areas of the engine map. The selected grid and covered points are shown in Figure 5.7.

In this case, the variance information is not as relevant as in other cases, because of the stochastic nature of inputs. The mean value of all states is plotted at each time step in Figure 5.8. All three methods perform well and the rate of convergence can be tuned varying  $\sigma_v^2$  and  $\sigma_w^2$ .

A quick view on the effects of the hypothesis for the SKF is shown for one element  $x_6$  in top plot of Figure 5.9, which shows the variance  $P_6$  of  $x_6$  for the KF and the SKF (not applicable for the SSKF). This value indicates the observability of  $x_6$ : when it increases monotonically,  $x_6$  is not observable. In some parts, an offset between  $P_6$  for the KF and the SKF appears because of the neglected covariances in the SKF, but in the end, this offset is absorbed and the two methods behave similar. The bottom plot of Figure 5.9 shows  $x_6$  evolution for the three methods.  $x_6$  is updated when the element becomes observable. Note that  $x_6$  is already active in the first iteration and  $K_1^{SS}$ , which represents the converged value of the equivalent LTI system, is non-null. Then, the SSKF updates  $x_6$  in  $k = 1$  while  $K_1$  for the KF and the SKF is null. Anyway, this is not an advantage of this method: the KF and the SKF can behave similar if  $P_0$  is non-null, but here for the simulations  $\mathbf{P}_0 = [\mathbf{0}]$ .



**Figure 5.8.** Averaged value of table parameters for the three updating methods when considering an uniform distributed inputs sequence. The correct value is 1.



**Figure 5.9.** Variance analysis for an uniform distributed inputs sequence. Top: variance evolution of the 6th element  $P_6$  where black thick line is the KF and gray line is the SKF ( $P_6$  is normalised by 100 in both cases). Bottom:  $x_6$  value where black thick line is the KF, gray line is the SKF, thin black is the SSKF.

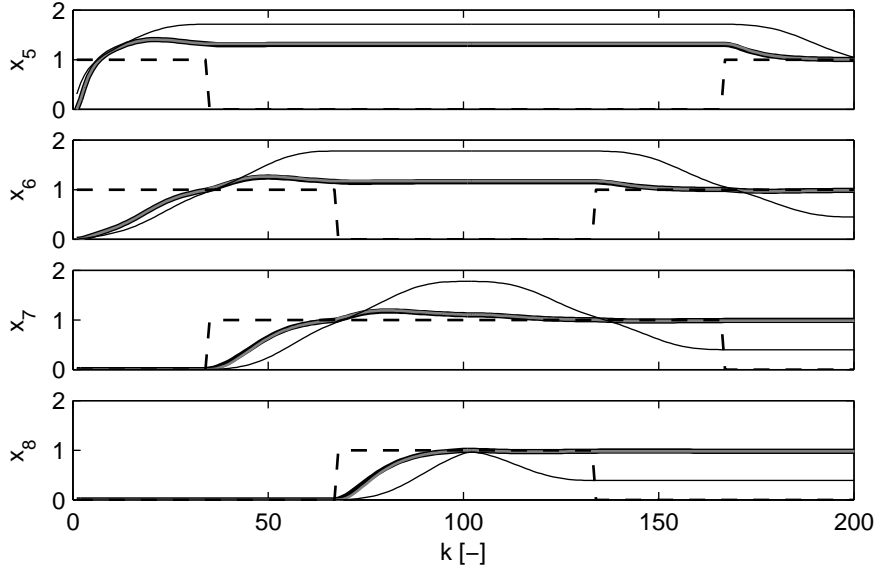
When there exists no previous knowledge of the system to be learnt, it should be advisable to initialize  $P_0$  with a certain value for speeding up the updating during the first iterations.

### 5.3.4.2 Simulation 2: Input with a linear variation

$\Theta_r$  is identified with a sequence  $\{u_k\}_{k=1}^{200}$  where  $u_1$  varies monotonically from 0 to 60 (from  $k = 0$  to 100) and coming back from 60 to 0 (from  $k = 101$  to 200), while  $u_1 = 2000 \forall k$ . This condition is highly restrictive as the excitation level is not high: at each  $k$  only small variations of input  $u_2$  are applied and due to this, the full convergence condition is not fulfilled in the first running of one area, and the interactions between areas and the way back are necessary to ensure the convergence of both the KF and the SKF. Three matrix areas are excited (see (5.29)); area 1: elements 5,6,9,10; area 2: 6,7,10,11; area 3: 7,8,11,12. Moreover, due to the inputs nature, the row 5-8 evolution is equivalent to that of 9-12.

$$\Theta_k = \begin{bmatrix} \cdot & \cdot & \cdot & \cdot \\ \hline x_{5,k} & x_{6,k} & x_{7,k} & x_{8,k} \\ x_{9,k} & x_{10,k} & x_{11,k} & x_{12,k} \\ \hline \cdot & \cdot & \cdot & \cdot \end{bmatrix} \quad (5.29)$$

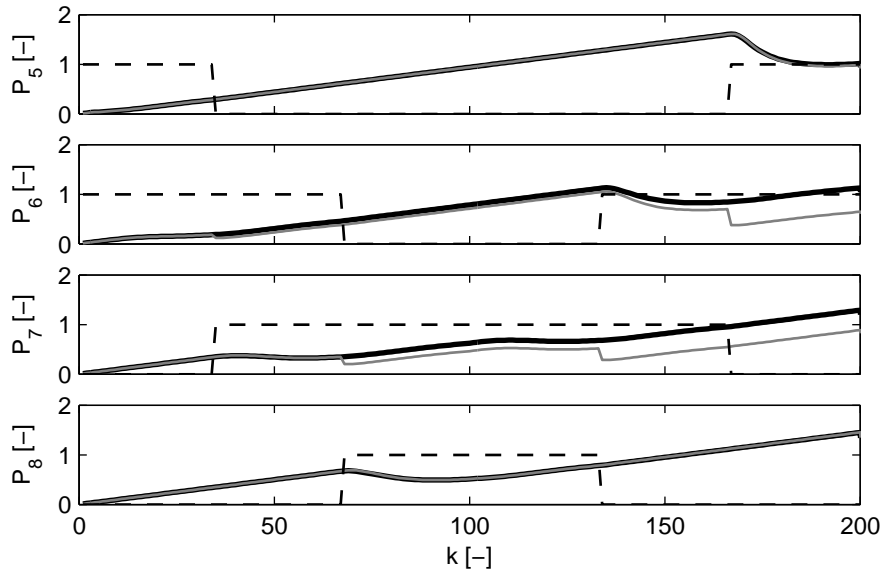
Figure 5.10 shows the evolution of the 4 states  $x_5$  to  $x_8$ . First thing to note is the similar performance of the KF and the SKF and the instabilities of the SSKF. For instance,  $x_5$  is in the area 1, and during the first run it is not capable of converging, but in one complete round the KF and SKF methods converge, and the SSKF seems to do it, at least in  $k = 200$ . The other states have slightly different performances, because the first area has been already covered; e.g.  $x_6$  is closer to the convergence as being a member of area 1 and area 2, but until the way round it does not get the full convergence for the KF and SKF methods. Similar behaviours are observed for  $x_7$  with these methods. In addition,  $x_8$ , as member of the area 3, has the advantage that areas 1 and 2 have been covered first, and the information is already available for getting the full convergence for the KF and SKF in  $k = 100$ . The full observability (in the sense of convergence) of these 8 elements is reached only when  $\mathbf{O}$  in (5.22) copes with the three areas when using the KF and SKF. Nevertheless, the SSKF estimation does not converge for  $x_6$ ,  $x_7$  and  $x_8$  due to no state error information is tracked, and the system and convergence becomes more dependent on data.



**Figure 5.10.** States evolution for a monotonically varied inputs sequence. Black thick line is KF, gray line is SKF, thin black is SSKF and dashed line shows when the element is observable (1 is active; 0 is not active).

In order to understand how KF and SKF behave, it is informative to also study the variance and covariance evolution (remember that this is not present for the SSKF). Figure 5.11 shows the variance evolution for row elements, while Figure 5.12 shows the evolution of a few selected covariances between states (corresponding to non-diagonal elements of  $\mathbf{P}$ ). As for the random case in Figure 5.9, variance evolution for the KF and SKF is similar, although some deviation exists in elements 6 and 7. This is explained for covariances or non diagonal  $P_{ij} \forall i \neq j$  elements that the SKF only accounted for active areas. This is shown in Figure 5.12 where some crossed covariances are plotted. When the pair of elements is not active, covariances of these are not tracked, equivalent to reset them to zero in the global  $\mathbf{P}$ . This is the particular difference with the KF, which always tracks the covariance, although the area would not be active, increasing the computing requirements without a justified improvement in the estimation. However, when an area is active, the SKF covariance evolves as the KF and in the limit is equivalent to the KF. The difference is larger initially but after the observations have converged the difference is negligible, as shown in Figure 5.10. This pattern may slightly be modified linking the covariance values to elements pairs and not to the

involved areas (excepting boundaries, all pair of elements are shared in two different areas), although final results are similar.

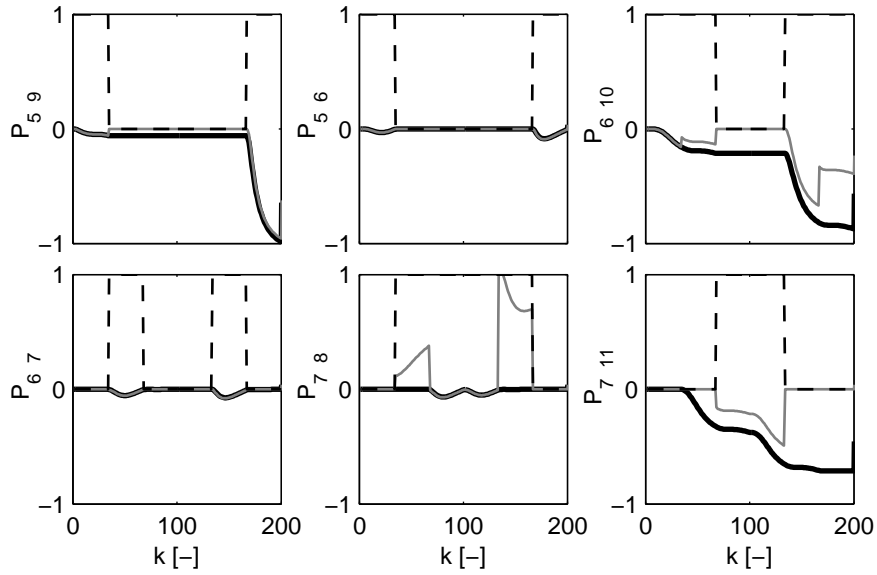


**Figure 5.11.**  $P_{ii}$  analysis for a monotonically varied inputs sequence. Values are normalised by 100. Variances of elements 9-12 are equal to those of 5-9 because of symmetric properties. Black line is KF, gray line is SKF and dashed line shows when the element is active (1 active; 0 is not active).

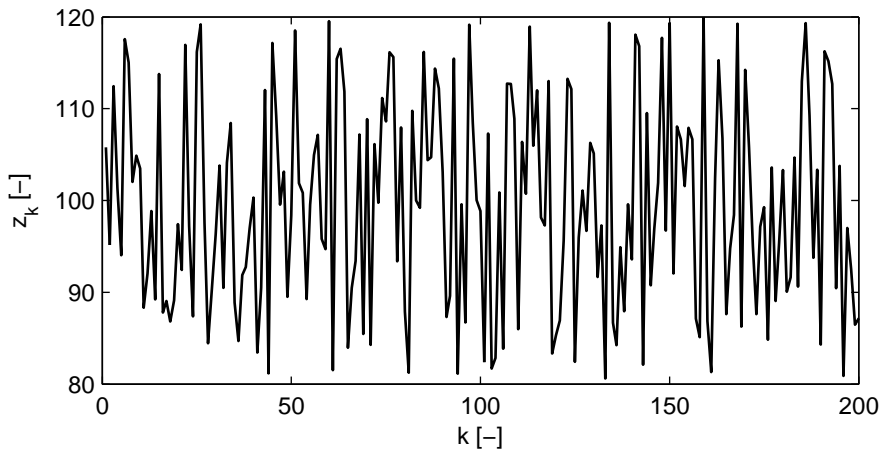
### 5.3.4.3 Simulation 3: Measurement noise rejection

An uniform distributed noise with zero mean and maximum amplitude of 0.2 is applied to the measurement given in simulation 2;  $z_k$  is shown in Figure 5.13. A robust method must filter this noise and converges to the true values.

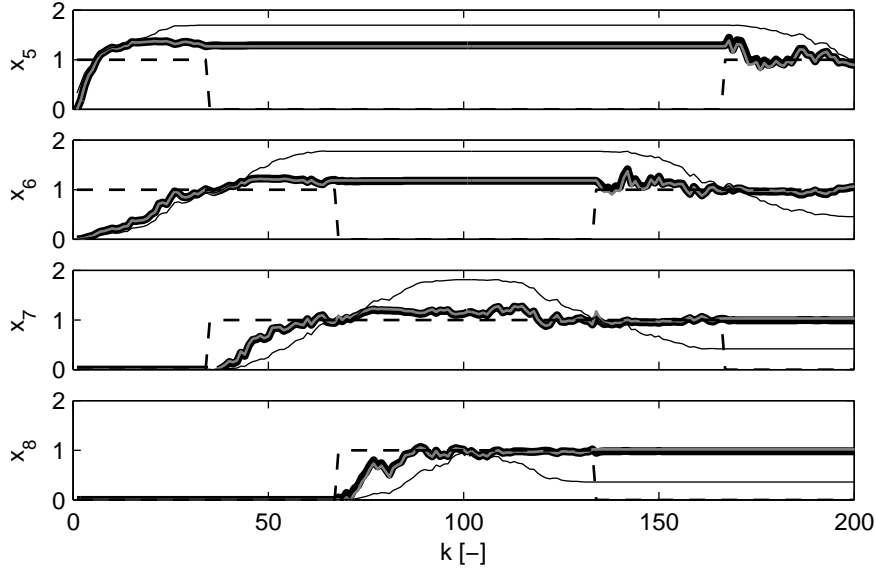
The resulting estimations are shown in Figure 5.14, and the evolutions are quite similar to the ones of Figure 5.10 although with a slight noise transmission. The KF and SKF perform well, being able to also filter the noise, but the SSKF is not capable of converging again having a similar response as in the simulation 2. Variance and covariances are exactly the same of Figures 5.11 and 5.12 as the state-space system and the inputs sequence does not change.



**Figure 5.12.**  $P_{ij}$  analysis for a monotonically varied inputs sequence. Values are normalised by 100. Black line is KF, gray line is SKF and dashed line shows when the element is active (1 active; 0 is not active).



**Figure 5.13.** Measurement  $z_k$  plot with the addition of a uniform distributed noise with zero mean and maximum amplitude of 0.2.



**Figure 5.14.** States evolution for noisy measurements. Black thick line is KF, gray line is SKF, thin black is SSKF and dashed line shows when the element is active (1 is active; 0 is not active).

#### 5.3.4.4 Conclusions from the simulations

The SKF method is demonstrated to have similar accuracy as the KF, at least for the numerical cases proved in this section. Furthermore, the KF and SKF solutions converge for all cases, including the simulations 2 and 3, which are specifically restrictive because of the low level of excitation. The KF, as shown in Figure 5.6 and table 5.1, is computationally heavy. The SSKF is the fastest and lightest method as the variance is not tracked and  $K^{SS}$  formulation is derived analytically. The SKF only needs around 1.15 times the SSKF time for calculation, and the required memory resources are similar, except for the track of the variances. The SSKF behaves well in situations where the variance tracking is not critical, e.g. the simulation 1, but it is less robust when the data is structured and the level of excitation is low, e.g. simulations 2 and 3. The KF and SKF are also capable of filtering the noise, a minimum condition for a correct updating, and SSKF also filters the noise, but its stability depends on the data. In comparison, the SKF method, derived from the KF, is demonstrated to be an efficient and accurate method for updating look-up tables, at least for the presented simulation conditions.



### 5.3.5 The dynamic equations for learning

The learning algorithms presented in the previous subsections have considered a quasi-static model, i.e. the model outputs are directly calculated by a table interpolation. In this subsection, the state-space model (5.17) is slightly modified for coping with sensor dynamics and inputs  $u$  are treated as well.

Usual inputs for the tables in diesel engines are the signals of the ECU: speed  $n$  and injection  $m_f$ , and as far as they are fast signals, no special treatment is necessary in principle, while the delays are still needed to phase the system correctly with the  $\text{NO}_x$  sensor

$$u_{1,k} = z^{-\tau/T_s} n(k) \quad (5.30a)$$

$$u_{2,k} = z^{-\tau/T_s} m_f(k) \quad (5.30b)$$

where  $\tau$  represents the averaged sensor delay obtained with the procedures shown in Chapter 3 ( $T_s$  is the sampling period). This is preferred to apply the delay in the state-space model for avoiding to increase the dimensions of the system.

In this basis, the table output  $x_\Theta(k) \in \mathbb{R}$  is calculated by means of interpolating the adaptive map  $\Theta_k$

$$x_\Theta(k - \tau/T_s) = \Theta_k(u_{1,k}, u_{2,k}) \quad (5.31)$$

defining a quasi-static representation of the table output where  $k$  defines the current time.

Sensor dynamics are considered as in Equation (3.3) for computing  $x_{\Theta_f}(k) \in \mathbb{R}$

$$x_{\Theta_f}(k) = \frac{1-a}{1-az^{-1}} x_\Theta(k - \tau/T_s) \quad (5.32)$$

and the global system (5.17) must be augmented with one extra-dimension considering  $x_{\Theta_f}(k)$

$$x_k^w = \left[ \begin{array}{c|c} \mathbf{I}_{\mathbf{n}_r \mathbf{n}_c \times \mathbf{n}_r \mathbf{n}_c} & \\ \hline (1-a)q_k & a \end{array} \right] x_{k-1}^w + w_k \quad (5.33a)$$

$$z_k^w = H^w x_k^w + v_k \quad (5.33b)$$

with

$$x_k^w = [x \quad x_{\Theta_f}]_k^T \quad (5.34)$$

$$H^w = [0 \cdots 1] \quad (5.35)$$

being  $x_k^w \in \mathbb{R}^{n_r n_c + 1}$  with  $x_k \in \mathbb{R}^{n_r n_c}$  as in (5.17), and  $z_k^w \in \mathbb{R}$ . The Index  $w$  stands for *wide*.

The local observable system (5.21) is now

$$x_k^{ow} = F_k^{ow} x_{k-1}^{ow} + w_k \quad (5.36a)$$

$$z_k^{ow} = H^{ow} x_k^{ow} + v_k \quad (5.36b)$$

$$F_k^{ow} = \left[ \begin{array}{c|c} \mathbf{I}_{4 \times 4} & \begin{matrix} 0 \\ 0 \\ 0 \\ 0 \end{matrix} \\ \hline (1-a)q_k^o & a \end{array} \right] \quad (5.37)$$

$$H^{ow} = [0 \ 0 \ 0 \ 0 \ 1] \quad (5.38)$$

$$x_k^{ow} = [x_1^o \ x_2^o \ x_3^o \ x_4^o \ x_{\Theta f}]_k^T \quad (5.39)$$

where  $H^{ow}$  is constant, because  $q_k^o$  and the filtering parameter  $a$  are included now in the time varying process matrix  $\mathbf{F}_k^{ow}$ .  $\mathbf{Q}^{ow}$  process noise matrix is also augmented, and includes an extra-noise term  $\sigma_f^2$  for the new state  $x_{\Theta f}$

$$\mathbf{Q}^{ow} = \left[ \begin{array}{c|c} \mathbf{Q} & \mathbf{0} \\ \hline \mathbf{0} & \sigma_f^2 \end{array} \right] \quad (5.40)$$

Even though the new local system has one more dimension ( $5 \times 5$ ), the SKF hypotheses are still applicable as proved in Appendix 5.C. Chapter 6 will prove the algorithms under real conditions in a diesel engine considering dynamics (the sensor output will be used there as reference) and completing the study on the methods capabilities.

## 5.4 Conclusions

This Chapter has presented different observers based on the KF in order to estimate engine variables and update look-up tables.

First, a drift correction algorithm is proposed by augmenting a state-space model with an extra-state for tracking the drift which exists between a fast biased model and a delayed and filtered but steady-state accurate sensor signal.

The data fusion is done through a steady-state KF, which may be computed beforehand. The algorithm allows to estimate online the actual considered signal.

Second, the attention turns out to the design of adaptive algorithms for look-up tables fitting and updating. Kalman filtering is used for updating because of its capability for tracking the system and parameters ageing. Computational issues involved when of updating look-up tables online with the KF based methods are then addressed. Two novel methods are designed for the current work:

- A steady-state KF (SSKF) approach using precalculated membership functions based on an off-line solution to the KF main equation. Table elements are updated in a fast way and the computational burden is low. The method relies on neglecting the covariance information, given by the standard KF, and then setting out the stationary problem where the inputs can be considered as constants in order to calculate the steady-state gain.
- A simplified KF (SKF) with similar accuracy as the standard KF, but that requires a much lower memory resources and calculation time. The local observable system serves as inspiration for analysing the variance matrix performance and is used for comparison. The SKF is based on the KF but neglects covariances of locally unobservable states.

The only requirement for the full table adaptation when using online learning algorithms is running the engine in all the matrix areas. The stability and convergence is assessed by means of simulating the algorithms. The memory saving and computational reduction that SKF offer, as well as the robust performance, make this algorithms suitable for being implemented on commercial ECUs. The SSKF has also an interesting performance and a very low computational burden, although the stability depends on the data quality.

In the following Chapter, these algorithms are combined and different results are presented in order to estimate  $\text{NO}_x$  and  $\lambda^{-1}$  in diesel engines.

## 5.A Analytical solutions to the Riccati equations

The drift correction algorithm (Section 5.2.1) and the SSKF approach (Section 5.3.2) for updating look-up tables constitute steady-state versions of the KF. This special property drives to calculate a steady-state Kalman

gain, allowing to precompute it and saving online resources. The Riccati iterative matrix system (2.13) becomes a matrix algebraic system and analytical expressions may be solved.

### 5.A.1 Drift correction model

The dynamic Riccati equation (2.13) is steady-state as  $K$  and  $\mathbf{P}$  converge after a certain number of iterations [7], which depend on the initial values. Then, it is possible to derive an analytical solution as proved by Kalman and Bucy [6]. Some authors have given solutions to particular problems, such as [18–21] or indeed computationally effective solutions as [22]. Here, an analytical solution to the Riccati matrix algebraic system (2.13) considering the model (5.1) is derived.

$\mathbf{P}_{\mathbf{k}-1}$  matrix is

$$\mathbf{P}_{\mathbf{k}-1} = \begin{bmatrix} p_1 & p_2 \\ p_3 & p_4 \end{bmatrix} \quad (5.41)$$

Operating (2.13a),  $P_{k|k-1}$  results

$$\mathbf{P}_{k|k-1} = \begin{bmatrix} p_1 + \sigma_w^2 & (1-a)p_1 + ap_2 \\ (1-a)p_1 + ap_3 & q \end{bmatrix} \quad (5.42)$$

where

$$q = (1-a)((1-a)p_1 + ap_2) + a(1-a)p_3 + a^2p_4 \quad (5.43)$$

Calculating  $K_k$  from (2.13b)

$$K_k = \begin{bmatrix} k_1 \\ k_2 \end{bmatrix} = \begin{bmatrix} \frac{(1-a)p_1 + ap_2}{q + \sigma_v^2} \\ \frac{q}{q + \sigma_v^2} \end{bmatrix} \quad (5.44)$$

From (2.13c)

$$\mathbf{P}_{\mathbf{k}} = \begin{bmatrix} p_1 + \sigma_w^2 - \frac{((1-a)p_1 + ap_2)((1-a)p_1 + ap_3)}{q + \sigma_v^2} & ((1-a)p_1 + ap_2) \frac{\sigma_v^2}{q + \sigma_v^2} \\ ((1-a)p_1 + ap_3) \frac{\sigma_v^2}{q + \sigma_v^2} & \frac{q\sigma_v^2}{q + \sigma_v^2} \end{bmatrix} \quad (5.45)$$

The Kalman filter is steady-state and observable and sequences  $\{\mathbf{P}_{\mathbf{k}}\}_{k=1}^{k=\infty}$  and  $\{K_k\}_{k=1}^{k=\infty}$  are convergent. The system solution can be computed through solving the system that results from setting (5.41) and (5.45) as equal.

$$p_1 = p_1 + \sigma_w^2 - \frac{((1-a)p_1 + ap_2)((1-a)p_1 + ap_3)}{q + \sigma_v^2} \quad (5.46a)$$

$$p_2 = ((1-a)p_1 + ap_2) \frac{\sigma_v^2}{q + \sigma_v^2} \quad (5.46b)$$

$$p_3 = ((1-a)p_1 + ap_3) \frac{\sigma_v^2}{q + \sigma_v^2} \quad (5.46c)$$

$$p_4 = \frac{q\sigma_v^2}{q + \sigma_v^2} \quad (5.46d)$$

$$k_1 = \frac{(1-a)p_1 + ap_2}{q + \sigma_v^2} \quad (5.46e)$$

$$k_2 = \frac{q}{q + \sigma_v^2} \quad (5.46f)$$

From (5.46b) and (5.46c),  $p_2 = p_3$ , which was already known as  $\mathbf{P}$  is symmetric, and system (5.46a) really has 5 equations with 5 unknowns.

From (5.46a) and (5.46e)

$$\sigma_w^2 = k_1^2(q + \sigma_v^2) \quad (5.47)$$

and from (5.46f)

$$q + \sigma_v^2 = \frac{\sigma_v^2}{1 - k_2} \quad (5.48)$$

and combining these two expressions,

$$k_1 = \frac{\sigma_w}{\sigma_v} \sqrt{(1 - k_2)} \quad (5.49)$$

which relates  $k_1$  to  $k_2$ .

$q$  parameter (5.43) depends on  $\mathbf{P}$  elements and must be re-arranged for the correct problem solution. Then, from the symmetry of  $\mathbf{P}$  ( $p_2 = p_3$ ), and considering (5.46b) and (5.46e)

$$p_2 = p_3 = \sigma_w \sigma_v \sqrt{1 - k_2} \quad (5.50)$$

From (5.46d) and (5.46f)

$$p_4 = k_2 \sigma_v^2 \quad (5.51)$$

and finally from (5.46b), (5.48) and (5.50)

$$(1-a)p_1 + ap_2 = \frac{p_2}{1-k_2} = \frac{\sigma_w \sigma_v \sqrt{1-k_2}}{1-k_2} \quad (5.52)$$

Substituting (5.50), (5.51) and (5.52) in (5.43)

$$q = (1-a)(1+a-ak_2) \sqrt{\frac{\sigma_w^2 \sigma_v^2}{1-k_2}} + a^2 k_2 \sigma_v^2 \quad (5.53)$$

For convenience, (5.53) is set as function of the noise covariance trade-off  $\sigma_w^2/\sigma_v^2$

$$q = \left( (1-a)(1+a-ak_2) \sqrt{\frac{\sigma_w^2/\sigma_v^2}{1-k_2}} + a^2 k_2 \right) \sigma_v^2 = s \sigma_v^2 \quad (5.54)$$

defining parameter  $s = f(\sigma_w^2/\sigma_v^2)$

$$s = (1-a) \sqrt{\frac{\sigma_w^2/\sigma_v^2}{1-k_2}} (1+a-ak_2) + a^2 k_2 \quad (5.55)$$

Considering (5.46f), (5.49) with (5.54),  $K$  can be computed

$$k_1 = \sqrt{(1-k_2) \frac{\sigma_w^2}{\sigma_v^2}} \quad (5.56a)$$

$$k_2 = \frac{s}{s+1} \quad (5.56b)$$

The set (5.56) represents an analytical solution of the steady-state Kalman filter that may be solved iteratively. Moreover, (5.56) shows that  $K$  depends exclusively on the noise covariance trade-off  $\sigma_w^2/\sigma_v^2$  given a certain system (5.1).

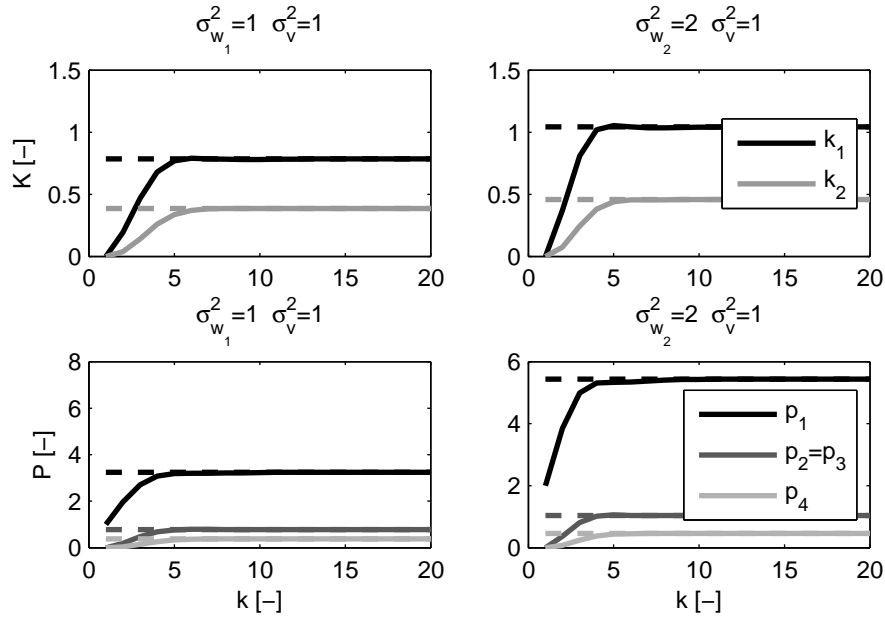
$\mathbf{P}$  is solved from (5.46a)-(5.46d) and (5.56)

$$p_1 = \sqrt{\frac{\sigma_w^2 \sigma_v^2}{1 - k_2}} \left( 1 + \frac{ak_2}{1 - a} \right) \quad (5.57a)$$

$$p_2 = p_3 = \sqrt{\sigma_w^2 (1 - k_2)} \quad (5.57b)$$

$$p_4 = k_2 \sigma_v^2 \quad (5.57c)$$

although is not really necessary for the algorithm implementation.



**Figure 5.15.** Kalman gain  $K$  and covariance matrix  $P$  calculation for two noise sets with  $a = 0.8$ . Solid lines corresponds to the standard Kalman formulation by the iterative Riccati equation (2.13), while dashed lines correspond to the proposed analytical formulation ((5.56) for  $K$  and (5.57) for  $P$ ).

**Numerical example.** A numerical example using (5.56), (5.57) for calculating  $P$  and  $K$  of a steady-state Kalman filter is presented. The result must be

equal to calculate the converged  $\mathbf{P}_k$  and  $K_k$  from (2.13) when the simulation time is high enough:

$$K = \lim_{k \rightarrow +\infty} K_k \quad (5.58a)$$

$$\mathbf{P} = \lim_{k \rightarrow +\infty} \mathbf{P}_k \quad (5.58b)$$

Parameter  $a$  is 0.8 and in all cases  $\sigma_v^2 = 1$ , while two different values are applied to  $\sigma_w^2$ :  $\sigma_{w1}^2 = 1$  and  $\sigma_{w2}^2 = 2$ .

Figure 5.15 shows the results applying the standard Kalman formulation and the ones calculated by the analytical formulation. Note that  $x$  axis shows the number of iterations ( $k$ ), which only makes sense for the standard Kalman formulation as far as the proposed solution is calculated beforehand. Final values after convergence are exactly the same proving the validity of the proposed formulation.

### 5.A.2 The SSKF for updating look-up tables

Given the system (5.17) for updating look-up tables and a filter tuning  $\sigma_v^2/\sigma_w^2$ , the value of  $\mathbf{P}$  does not converge under normal operation, nor does  $K_k$ . This is because  $H_k$  matrix is not constant and depends on the engine operating conditions. But if the operating point condition is constant, system runs in a constant point ( $H_k = H$ ) with a certain  $\sigma_v^2/\sigma_w^2$ , then it is possible to derive a constant  $K^{SS}$ , although  $\mathbf{P}$  does not really converge. This is observed from simulations and then it can be set out for a 2D grid

$$K^{SS}(\eta_1, \eta_2, \sigma_v^2/\sigma_w^2) = \lim_{k \rightarrow +\infty} K_k \quad (5.59)$$

Taking into account the symmetric property of the problem, the solution is set out for a 1D table  $\Theta$ , whose elements are expanded in a state vector and whose 1D local observable has a state vector

$$x_k^o = [\Theta_i \quad \Theta_{i+1}]_k^T \in \mathbb{R}^2$$

where  $i$  denotes the position of the first active element in  $\Theta$ . The state-space equations of the local observable system are

$$x_k^o = \mathbf{I}_{2 \times 2} x_{k-1}^o + w_k \quad (5.60a)$$

$$z_k = q_k^o x_k^o + v_k \quad (5.60b)$$



with

$$q_k^o = [1 - \eta \quad \eta] \quad (5.61)$$

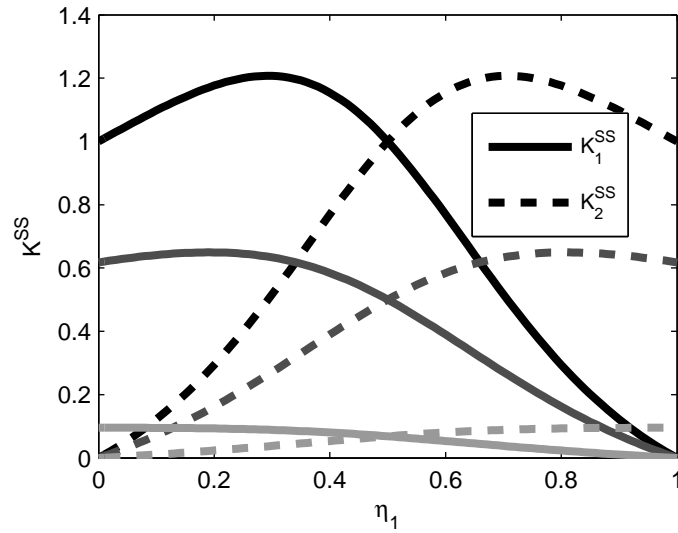
and considering  $r$  as the vector that defines the 1D grid

$$\eta(k) = \frac{u_k - r_i}{r_{i+1} - r_i} \quad (5.62)$$

where  $u_k \in \mathbb{R}$  is now the scheduling point considered as input to the table.

Noise covariances are

$$\mathbf{Q} = \begin{bmatrix} \sigma_w^2 & \\ & \sigma_w^2 \end{bmatrix}; \quad R = \sigma_v^2 \quad (5.63)$$



**Figure 5.16.** Membership functions for the Kalman gain correction for the 1D case for  $\sigma_v^2/\sigma_w^2 = 0$  (black),  $\sigma_v^2/\sigma_w^2 = 0.01$  (medium grey) and  $\sigma_v^2/\sigma_w^2 = 0.1$  (light grey).

By imposing the condition 5.59, the value of Kalman gain  $k_i$  can be derived analytically with a certain effort

$$k_i(\eta, \sigma_v^2/\sigma_w^2) = \frac{0.5(1-\eta)(1+s)}{0.5(1+s)(1-2\eta+2\eta^2) + \sigma_v^2/\sigma_w^2} \quad (5.64)$$

where

$$s = \sqrt{1 + \frac{4}{(1-2\eta+2\eta^2)} \frac{\sigma_v^2}{\sigma_w^2}} \quad (5.65)$$

showing explicitly the dependance of  $k_i$  with the noise trade-off  $\sigma_v^2/\sigma_w^2$  and the interpolation function defined by  $\eta$ . The result is used for calculating gain correction for the node  $i$ , when the observable region is between  $i$  and  $i + 1$  in the 1D case. By using symmetric properties

$$k_{i+1} = k_i(1 - \eta, \sigma_v^2/\sigma_w^2) \quad (5.66)$$

The function is plotted in Figure 5.16. For the 2D case,  $K^{SS}$  (see (5.59)) can be solved by considering at the same time the interpolation properties, and then the gain for a local observable system of a given 2D table with interpolation weightings  $\eta_1$  and  $\eta_2$  defined in (5.16) has the solution

$$K^{SS} = \left[ \begin{array}{l} k_i(\eta_1, \sigma_v^2/\sigma_w^2) \cdot k_i(\eta_2, \sigma_v^2/\sigma_w^2) \\ k_i(\eta_1, \sigma_v^2/\sigma_w^2) \cdot k_i(1 - \eta_2, \sigma_v^2/\sigma_w^2) \\ k_i(1 - \eta_1, \sigma_v^2/\sigma_w^2) \cdot k_i(\eta_2, \sigma_v^2/\sigma_w^2) \\ k_i(1 - \eta_1, \sigma_v^2/\sigma_w^2) \cdot k_i(1 - \eta_2, \sigma_v^2/\sigma_w^2) \end{array} \right] \quad (5.67)$$

The obtained solution can be easily generalised for the nD case by solving

$$\prod_{i=1}^n k_i(\xi_i, \sigma_v^2/\sigma_w^2) \quad (5.68)$$

where  $\xi = \eta_i$  or  $\xi = 1 - \eta_i$  depending on the considered node.

## 5.B Pseudo-codes of the SSKF and SKF methods

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### Algorithm 1: Pseudo-code of the SSKF method

---

**input** :  $\hat{x}_{k-1}, u_k$   
**output**:  $\hat{x}_k$

- 1 **while** *Algorithm is running* **do**
- 2     **if** *Active area changes* **then**
- 3          $x_{k-1}^o$  is stored in the correct positions of  $x_{k-1}$
- 4         Redefinition of new  $x_{k-1}^o$  ;
- 5     **end**
- 6     Computation of  $\eta_{1,k}$  and  $\eta_{2,k}$  as in (5.16)
- 7     Definition of  $q_k(\eta_{1,k}, \eta_{2,k})$  as in (5.15)
- 8     Computing  $K^{SS}$  as in (5.23)
- 9      $x_k^o = x_{k-1}^o + K_k^{SS} (z_k - q_k^o x_{k-1}^o)$
- 10 **end**
- 11 Updating of  $x_k$  considering  $x_k^o$

---



---

### Algorithm 2: Pseudo-code of the SKF method

---

**input** :  $\hat{x}_{k-1}, u_k, P_{k-1}, R, \mathbf{Q}^o, Q^u$   
**output**:  $\hat{x}_k, P_k$

- 1 **while** *Algorithm is running* **do**
- 2     **if** *Active area changes* **then**
- 3          $x_{k-1}^o$  is stored in the correct positions of  $x_{k-1}$
- 4          $diag(\mathbf{P}_{k-1}^o)$  and  $P_{k-1}^u$  are stored in a global vector  $P_{k-1}$
- 5         Redefinition of  $\mathbf{P}_{k-1}^o, P_{k-1}^u, x_{k-1}^o$
- 6     **end**
- 7     Computation of  $\eta_{1,k}$  and  $\eta_{2,k}$  as in (5.16)
- 8     Definition of  $q_k^o(\eta_{1,k}, \eta_{2,k})$  as in (5.15)
- 9      $\mathbf{P}_{k|k-1}^o = \mathbf{P}_{k-1}^o + \mathbf{Q}^o$
- 10      $K_k^o = \mathbf{P}_{k|k-1}^o \mathbf{q}_k^{oT} \left( \mathbf{q}_k^o \mathbf{P}_{k|k-1}^o \mathbf{q}_k^{oT} + \mathbf{R} \right)^{-1}$
- 11      $x_k^o = x_{k-1}^o + K_k^o (z_k - q_k^o x_{k-1}^o)$
- 12      $\mathbf{P}_{k|k-1}^o = (\mathbf{I}_{4 \times 4} - K_k^o q_k^o) (\mathbf{P}_{k-1}^o + \mathbf{Q}^o)$
- 13      $P_k^u = P_{k-1}^u + Q^u$
- 14 **end**
- 15 Updating of  $x_k$  and  $P_k$  considering  $x_k^o, diag(\mathbf{P}_k^o)$  and  $P_k^u$

---

### 5.C The dynamic system for the SKF method

The system (5.36) is the local dynamic observable system. For this system, the split in (5.25) is still valid. The only modification is that the new observable part  $\mathbf{P}_k^{\text{wo}}$  has  $n$  extra dimensions corresponding to the  $n$ :th order discrete filter. In the paper, a first order discrete filter is used, and then  $\mathbf{P}_k^{\text{wo}}$  is  $5 \times 5$ .

$$\mathbf{P}_k^{\text{w}} = \left[ \begin{array}{c|c} \mathbf{P}_k^{\text{u}} & \\ \hline & \mathbf{P}_k^{\text{wo}} \end{array} \right] \quad (5.69)$$

$$\mathbf{P}_k^{\text{wo}} = \left[ \begin{array}{c|c} \mathbf{P}_k^{\text{o}} & P_k^{\text{ndo}T} \\ \hline P_k^{\text{ndo}} & P_k^{\text{do}} \end{array} \right] \quad (5.70)$$

where  $P_k^{\text{do}}$  is the scalar variance coupled to sensor dynamics and  $P_k^{\text{ndo}}$  is the vector  $1 \times 4$  with the covariances between sensor estimation and observable parameters.  $\mathbf{P}_k^{\text{w}}$  is a positive-semidefinite matrix, so is  $\mathbf{P}_k^{\text{wo}}$ .  $K_k$  is reordered to contain the unobservable part  $K_k^{\text{u}} = 0$  and the observable part  $K_k^{\text{o}}$ , and the latter in the parameters related  $K_k^{\theta}$  and dynamics related part  $K_k^{\text{s}}$

$$K_k^{\text{w}} = \left[ \begin{array}{c} K_k^{\text{u}} \\ \hline K_k^{\text{o}} \end{array} \right] = \left[ \begin{array}{c} 0 \\ \hline K_k^{\theta} \\ \hline K_k^{\text{s}} \end{array} \right] \quad (5.71)$$

and (5.26) is now

$$\begin{aligned} \mathbf{P}_{k+1} &= \left[ \begin{array}{c|c} \mathbf{I} & \mathbf{0} \\ \hline \mathbf{0} & \mathbf{I}_{5 \times 5} - K_k^{\text{o}} H^{\text{w}} \end{array} \right] \left[ \begin{array}{c|c} \mathbf{P}_k^{\text{u}} + \mathbf{Q}^{\text{u}} & \vdots \\ \hline \vdots & \mathbf{F}_k^{\text{w}} \mathbf{P}_k^{\text{wo}} (\mathbf{F}_k^{\text{w}})^{\text{T}} + \mathbf{Q}^{\text{wo}} \end{array} \right] \\ &= \left[ \begin{array}{c|c} \mathbf{P}_k^{\text{u}} + \mathbf{Q}^{\text{u}} & \vdots \\ \hline \vdots & (\mathbf{I}_{5 \times 5} - K_k^{\text{o}} H^{\text{w}}) (\mathbf{F}_k^{\text{w}} \mathbf{P}_k^{\text{wo}} (\mathbf{F}_k^{\text{w}})^{\text{T}} + \mathbf{Q}^{\text{wo}}) \end{array} \right] \end{aligned} \quad (5.72)$$

which shows that the SKF hypotheses are also applicable for the widespread dynamic system.  $\mathbf{P}_k$  is a positive-semidefinite matrix so are  $\mathbf{P}_k^w$  and  $\mathbf{P}_k^{wo}$ .

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# Chapter 6

## Adaptive estimation of $\text{NO}_x$ and $\lambda^{-1}$

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### 6.1 Introduction

This Chapter uses the adaptive algorithms developed in Chapter 5 in order to estimate  $\lambda^{-1}$  and  $\text{NO}_x$ , when a  $\text{NO}_x$  sensor is available. The use of adaptive algorithms also allows to tune model parameters with and without an initial knowledge [1], i.e. the algorithms can be used for online model fitting. For that purpose, three different algorithms are designed and the filter gains are calculated on the basis of the methods presented in Chapter 5:

- Algorithm A: The drift correction algorithm corrects the bias in the  $\lambda^{-1}$  and  $\text{NO}_x$  models. The bias is tracked online but models are not re-calibrated.
- Algorithm B: An adaptive table is used for learning and prediction without requiring the RT  $\text{NO}_x$  model. This approach is valid for slow cycles, such as the NEDC. The table is updated by using the SKF or the SSKF methods, as advisable.
- Algorithm C: Model is adapted and re-tuned online by tracking the nominal  $\text{NO}_x$ . With this strategy, the nominal  $\text{NO}_x$  map can be fitted by using the dynamic tests, indeed avoiding the steady-state campaign usually needed for its calibration. The state-space system (5.17) built for updating look-up tables should be slightly modified in order to consider the model structure.

Figure 6.1 schematises the three options, particularising for the  $\text{NO}_x$  estimation. The objective is on one hand, obtaining a fast and reliable estimation of the considered variable, i.e. without sensor filtering and delay and, on the other hand, introducing adaptive capabilities to the models.

In the following, these algorithms are selected for the fast estimation of  $\lambda^{-1}$  and  $\text{NO}_x$ .

## 6.2 Fast estimation of $\lambda^{-1}$

The estimation of  $\lambda^{-1}$  is the core of this section, where a complete methodology based on the drift correction model developed in Section 5.2.1 is proposed. The original problem is modified in order to consider the model and sensor uncertainties as well as the actual variation of the bias  $\theta_{\lambda^{-1}}$  with respect to the operating point and time as remarked in (4.3). In addition, the variable  $\lambda^{-1}$  is a relevant input for the  $\text{NO}_x$  model, and its observation by KF methods can be integrated in an adaptive  $\text{NO}_x$  model. First, the problem set-up and the methodology are explained.

### 6.2.1 Problem set-up and methodology

The goal is providing an estimation  $\hat{x}_{\lambda^{-1}}$  of the actual value of  $\lambda^{-1}$  combining the information provided by the sensor  $z_{\lambda^{-1}}$  and the model  $x_{\lambda^{-1}}$ . The basic idea is keeping the model dynamics while correcting the model bias (and its drift) profiting the sensor steady performance.



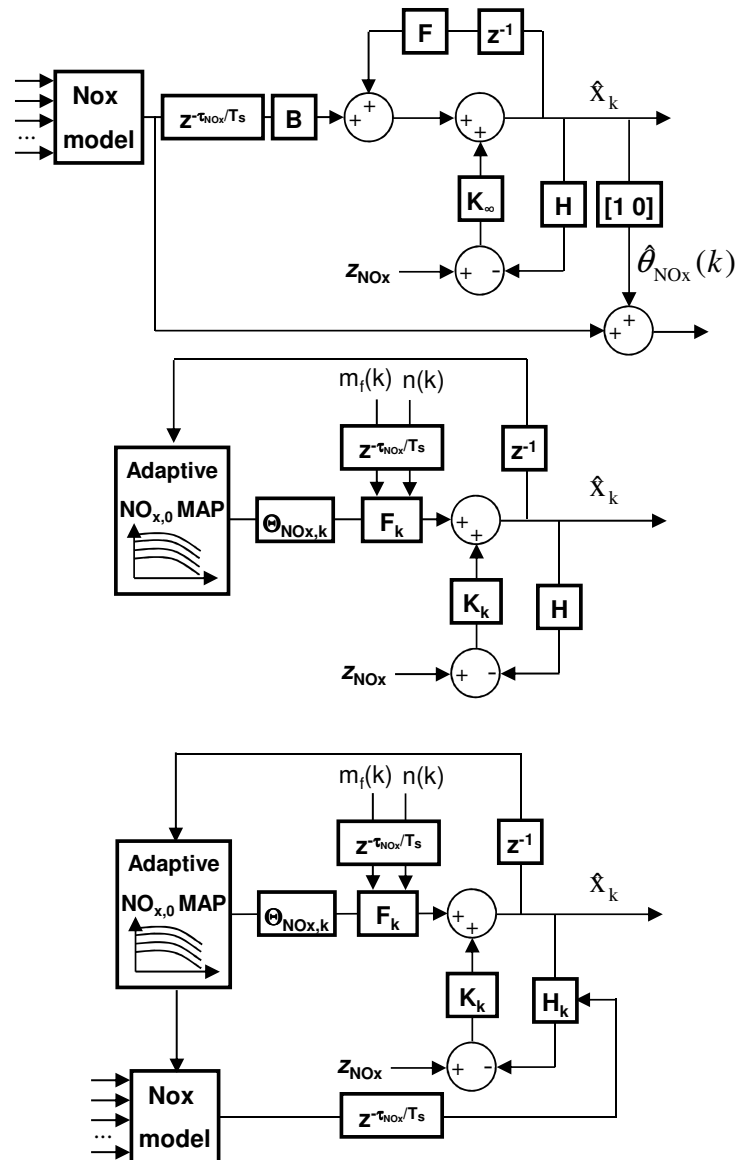


Figure 6.1. From top to bottom, algorithms 1 to 3: drift correction model, adaptive table for  $NO_x$  estimation and adaptive  $NO_x$  model.

Figure 6.2 shows an schematic representation of the proposed structure. All calculations and equations are set-up in discrete form where  $T_s$  represents the sampling period, which corresponds to 20 ms for the results shown in this work. Main inputs are the injected fuel mass  $m_f$ , the air mass flow  $m_a$ , the speed  $n$  and the sensor signal  $z_{\lambda^{-1}}$ .

Because of the bias strong dependence with the operating point conditions, an adaptive 2D look-up table  $\Theta_{\lambda^{-1},\mathbf{k}} \in \mathbb{R}^{n_r \times n_c}$  updated at instant  $k$  is used for accounting with this variation depending on the engine speed and load (see discussion in Section 4.3.1 about engine operating point representation). Hence the current bias at the instant  $k$  may be represented as

$$\theta_{\lambda^{-1}}(k) = \Theta_{\lambda^{-1},\mathbf{k}}(n(k), m_f(k)) + w_k \quad (6.1)$$

where a noise  $w_k \in \mathbb{R}$  is added to the tabulated value coming from the table  $\Theta_{\lambda^{-1},\mathbf{k}}$  at instant  $k$  for considering modelling errors, and for dealing with the system drift.

Through the combination of the sensor and the model information (conveniently delayed for being comparable), a KF is used for tracking the value of the bias and of the filtered value of  $\lambda^{-1}$ , thus providing the estimates  $\hat{\theta}_{\lambda^{-1}}$  and  $\hat{x}_{\lambda^{-1},f}$ . A set of 'freezing' conditions adds robustness to the algorithm stopping the integration when the sensor signals or the model are not reliable.

In order to cope with the system drift, the look-up table  $\Theta_{\lambda^{-1}}^{-1}$  is updated on the basis of the adaptive SSKF algorithm that uses the estimate  $\hat{\theta}_{\lambda^{-1}}$

$$\Theta_{\lambda^{-1},\mathbf{k}} = f(\Theta_{\lambda^{-1},\mathbf{k}}, \hat{\theta}_{\lambda^{-1}}(k)) \quad (6.2)$$

Finally, the estimation  $\hat{x}_{\lambda^{-1}}$  is built up from the current model output  $x_{\lambda^{-1}}$  and the tabulated value of the bias (see (6.5)).

### 6.2.2 Robustness against signals uncertainties

Three main circumstances affect the correct algorithm performance:

- Sensor saturation:  $\text{NO}_x$  sensors exhibit saturation problems for high  $\text{O}_2$  partial pressures (low  $\lambda^{-1}$ ). Figure 6.3 shows the saturation point for the tested  $\text{NO}_x$  sensor, where two units of this model are located downstream of the turbo and downstream of the engine here equipped with an AT line. If the KF is active, the integration would be incorrect in these situations as far as  $z_{\lambda^{-1}}$  is not reliable.

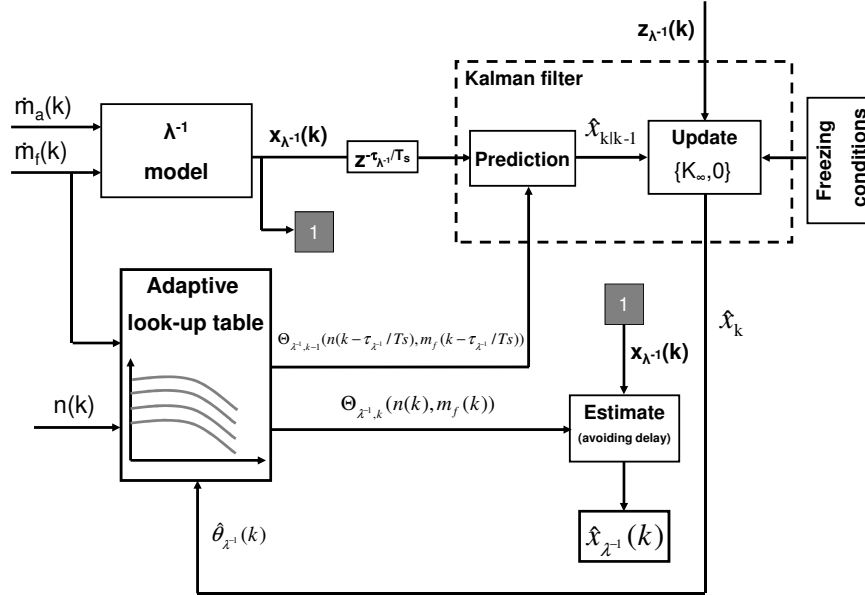
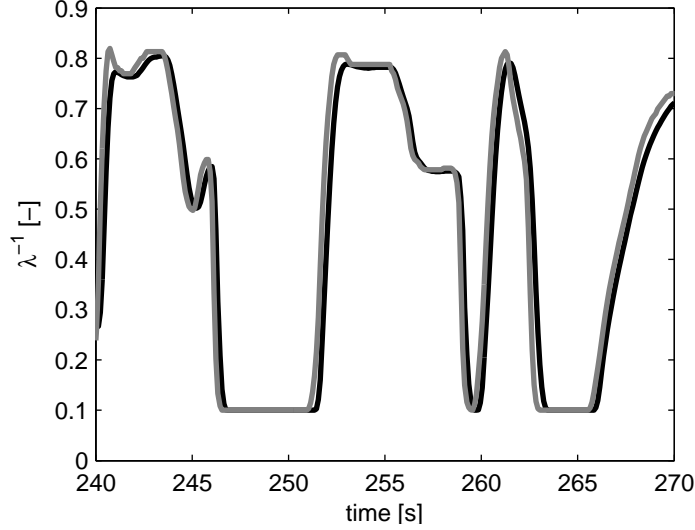


Figure 6.2. Schematic view of the proposed procedure for  $\lambda^{-1}$  estimation.

- Sensor and model uncertainties during sharp dynamic transients: The sensor behaviour, defined by  $a_{\lambda^{-1}}$  and  $\tau_{\lambda^{-1}}$ , varies with time, boundary conditions, system conditions and because of the ageing and unit-to-unit dispersion. For avoiding wrong integrations, these must be considered. Furthermore, the sensor model considered in (3.3) could not be complex enough for accounting with the actual sensor performance. An incorrect sensor model leads to peak errors, especially when a sharp transient occurs.
- Spurious measurements: Signals involved can present outliers or errors that could be fatal for the algorithm.

When these occur, the integration must be stopped. A set of deactivation IF-THEN or 'freezing' rules have been programmed for that. If any of the



**Figure 6.3.**  $z_{\lambda^{-1}}$  measurements from two  $NO_x$  sensors, the first located downstream of the turbine (gray) and the second downstream of the AT line (black) in an engine equipped with DOC, SCR and DPF (see the third configuration of Figure 5.1). Both sensors are the same commercial model and present a saturation in approximately  $\lambda^{-1} = 0.1$ .

following conditions is true, the updating of the KF is null; i.e.  $e_k = 0$ :

$$e_k \frac{1 - z^{-1}}{T_s} > F_1 \quad (6.3a)$$

$$x_{\lambda^{-1}} \frac{1 - z^{-1}}{T_s} > F_2 \quad (6.3b)$$

$$z_{\lambda^{-1}} \frac{1 - z^{-1}}{T_s} > F_3 \quad (6.3c)$$

$$z_{\lambda^{-1}} < F_4 \quad (6.3d)$$

$$z_{\lambda^{-1}} > F_5 \quad (6.3e)$$

$$\hat{x}_{\lambda^{-1}} < F_6 \quad (6.3f)$$

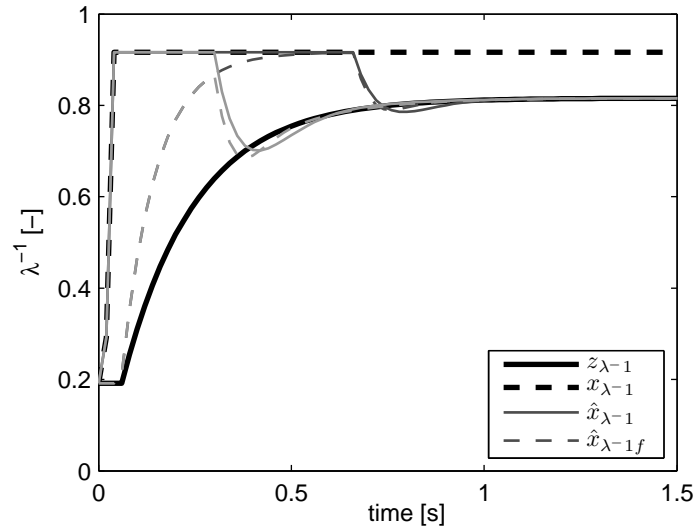
$$\hat{x}_{\lambda^{-1}} > F_7 \quad (6.3g)$$

where (6.3a) to (6.3c) deactivate the filter in sharp dynamic transients or in the case of spurious measurements; (6.3d) if sensor saturation or outliers appear; (6.3e) if measurement outliers are detected; finally (6.3f) and (6.3g) are included for providing robustness to the estimation. These conditions

make the table update highly insensitive to sensor model errors, sensor output errors and model output errors.

The freezing conditions allows to set a faster integration, i.e. a higher Kalman gain, as integrations at points with high expected errors on  $\hat{\theta}_{\lambda^{-1}}$  are avoided. The definition of thresholds needs to be done accordingly with the level of uncertainty in the sensor knowledge, signals noise level, model reliability and assumptions concerning the dynamic characteristics of the  $\lambda^{-1}$  evolution.

Figure 6.4 compares  $\hat{x}_{\lambda^{-1}}$  obtained with two different tunings of the thresholds and for one sharp injection transient, where  $\lambda^{-1}$  value is drastically varied and the bias is affected by the operating conditions (it varies from 0 at the beginning of the test to 0.1 after the step). In all cases, a significant error in the sensor model is assumed ( $a'_{\lambda^{-1}} \neq a_{\lambda^{-1}}$  and  $\tau'_{\lambda^{-1}} \neq \tau_{\lambda^{-1}}$ ). It can also be noticed that including freezing mitigates the overshoot in the correction; although a conservative tuning will result to completely deactivating the integration of the filter, and then no bias integration would be made during the engine operation. The Monte Carlo calibration method presented in [2] might be used for tuning the thresholds  $F_1$  to  $F_7$ .



**Figure 6.4.**  $\hat{x}_{\lambda^{-1}}$  for two different conditions of freezing for an injection step from 15 to 30 mg/str at 2250 rpm. Thin solid lines represents  $\hat{x}_{\lambda^{-1}}$ , lighter to darker gray is less to more severe freezing condition.  $\hat{x}_{\lambda^{-1}f}$  models linked with the estimate  $\hat{x}_{\lambda^{-1}}$  are also included with dashed lines.

### 6.2.3 An adaptive look-up table for modelling the drift

An adaptive look-up table (see Figure 6.2) is used for coping with the bias dependence with the operative conditions. The look-up table interpolation principle is based on a 2D bilinear interpolation. At every iteration, the initial estimation of the bias is interpolated from the table  $\Theta_\lambda^{-1}$  conveniently updated

$$\hat{\theta}_{\lambda^{-1}}(k|k-1) = \Theta_{\lambda^{-1}, \mathbf{k}-1}(n(k-\tau/T_s), m_f(k-\tau/T_s)) \quad (6.4)$$

where the table scheduling inputs  $n$  and  $m_f$  are conveniently delayed in order to consider the sensor delay. The KF provides an updated value of the bias (at the delayed input conditions) that is used for updating the table with a given learning method. Chapter 5 have made a review and developed two methods based on the EKF for updating look-up tables. For this formulation, the SSKF method presents the advantage that the table can be considered as an independent block, and then the converged  $K_\infty$  for the drift correction is still applicable. In this section, the results by using the SSKF are presented (although for a robust implementation, the SKF is advised).

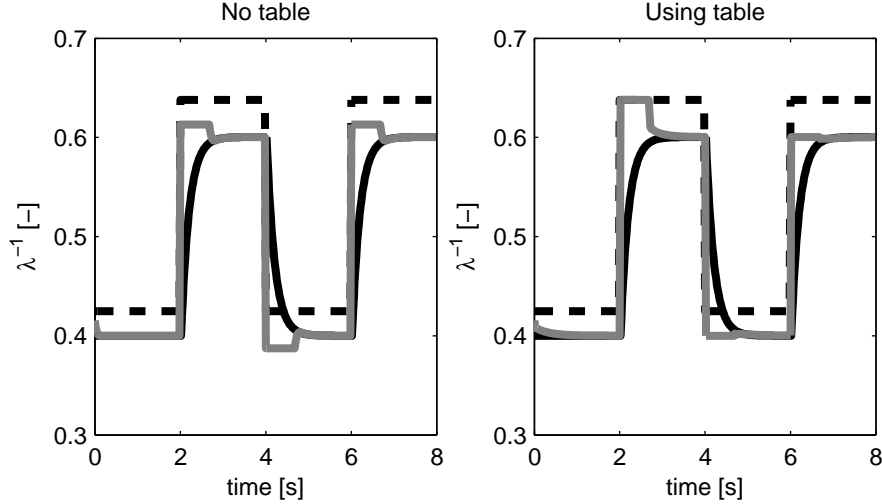
Once the table has been adapted, an updated value of the bias is obtained for the current operative conditions. This is used for building the final estimation

$$\hat{x}_{\lambda^{-1}}(k) = x_{\lambda^{-1}}(k) + \Theta_{\lambda^{-1}, \mathbf{k}}(n(k), m_f(k)) \quad (6.5)$$

where the current values of the inputs and the model are used at the instant  $k$ , in order to provide the more updated information and avoiding the causality problem related with the sensor delay when the table is not used.

For illustrating the look-up table performance, Figure 6.5 shows the simulation of  $\lambda^{-1}$  for a repetition of the same profile.  $\tau_{\lambda^{-1}}$  is known and null and  $a_{\lambda^{-1}}$  presents a certain error. The top plot represents the evolution if the adaptive table is not used. In the case of using the adaptive table (right plot), first part of the cycle serves for the bias identification (initially all elements of the table are set to zero), and in the second part of the cycle the estimation is significantly improved because of the stored values. Note that during transients, the error integration starts slightly after the step because of the freezing rules.

Different tables may also be used in case of using different combustion modes (split injection, exhaust system regeneration) for coping with the different trend of the bias on the  $m_f$  estimate when the injection settings change. Each table is updated only when the mode is activated and can add more accuracy to the estimate although incrementing the programming complexity.



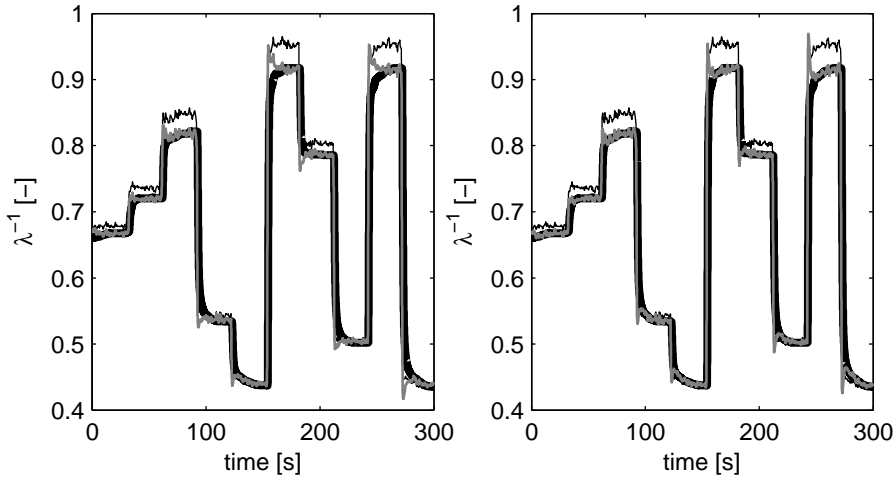
**Figure 6.5.**  $\hat{x}_{\lambda^{-1}}$  without table (left) and using an adaptive table (right) changing the operating point condition of the engine repeating two times a cycle by simulating. Black line is  $z_{\lambda^{-1}}$ , dashed black line is  $x_{\lambda^{-1}}$  and gray line is  $\hat{x}_{\lambda^{-1}}$ . Deactivation freezing rules are applied.

#### 6.2.4 Experimental results

In order to provide a reference signal for assessing the algorithm, injection steps are performed in the engine and  $\hat{x}_{\lambda^{-1}}$  calculated. Since the injection is a fast acting control variable, the actual  $\lambda^{-1}$  response is judged to be instantaneous and no more than one engine cycle delay is expected; being able to generate a reference for comparison. Furthermore, this signal is available online, making that the algorithm may be tuned and proved without a specific test rig beyond the engine and the ECU itself.

The Sensor model identification was discussed in Chapter 3, hence a delayed first order filter (see (3.3)) with parameters  $a_{\lambda^{-1}} = 0.92$  and  $\tau_{\lambda^{-1}} = 0.5s$  with a sampling frequency of 50 Hz ( $T_s=0.02s$ ) is used.  $K_{\infty}$  is precalculated offline for all cases as the KF is steady-state. Figure 6.6 shows the results for a set of different injection steps for a constant engine speed  $n = 1500$  rpm. For that Figure, the  $ZrO_2$  sensor has been moved downstream of the AT systems and a significant bias, 5% of its own value, has been added for exaggerating the correction; i.e  $x_{\lambda^{-1}}^{drift} = 1.05x_{\lambda^{-1}}$ . Furthermore,  $x_{\lambda^{-1}}$  presents a significant variable bias depending on the operating point with respect to  $z_{\lambda^{-1}}$ . The drift correction algorithm is applied to the estimation of  $\lambda^{-1}$ . Freezing conditions are relaxed because the sensor is well modelled. The left plot shows an ex-

ample where the integration is quite slow; the right plot shows the opposite situation where the integration is quite fast, provoking some peaks but with a lower settling time. Here, no model or table is used for modelling  $\theta_{\lambda^{-1}}$ . This makes that after every step,  $\theta_{\lambda^{-1}}$  shall be integrated for coping with its variation.



**Figure 6.6.**  $\hat{x}_{\lambda^{-1}}$  performing injection steps at 1500 rpm without using a look-up table. Left plot: Slow correction, with  $\sigma_w^2/\sigma_v^2 = 10^{-5}$ . Right plot: Fast correction, with  $\sigma_w^2/\sigma_v^2 = 10^{-4}$ . Legend: Thick black is  $z_{\lambda^{-1}}$ , thin black is  $x_{\lambda^{-1}}$  and gray is  $\hat{x}_{\lambda^{-1}}$ .

The use of an adaptive table  $\Theta_{\lambda^{-1}}$  smoothes the bias integration when the operating point changes once the table has been learnt. Figure 6.7 shows the same situation of Figure 6.6 but now using a look-up table with the grid

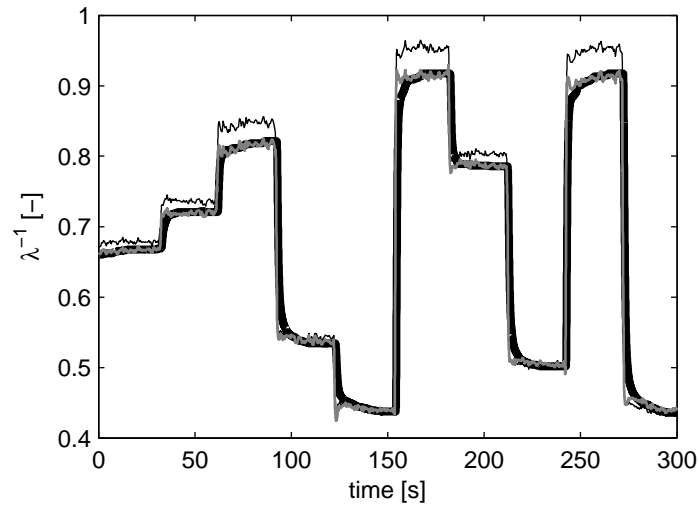
$$r = [750 \quad 1000 : 500 : 5000] \quad (6.6a)$$

$$c = [0 : 5 : 30 \quad 34 : 4 : 50 \quad 55 : 5 : 80] \quad (6.6b)$$

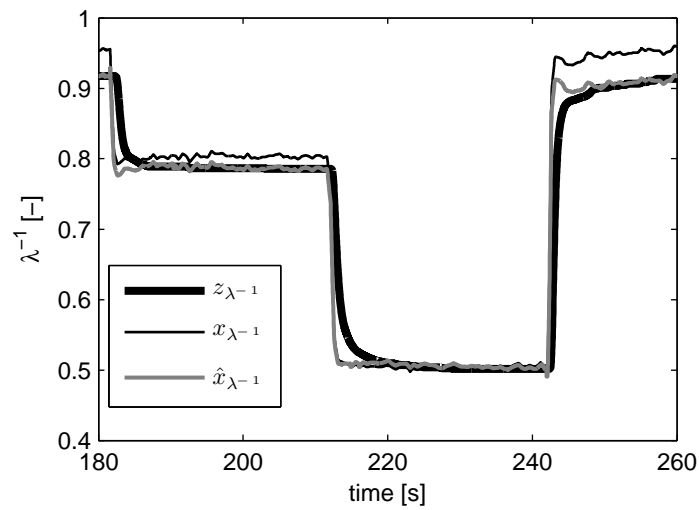
Figure 6.8 shows final results for three different injection steps of the previous cycle once that the bias has been perfectly learnt.  $\hat{x}_{\lambda^{-1}}$  keeps fast dynamics but converges to the steady-state value of  $x_{\lambda^{-1}}$ , in contrast to the biased model and the slow sensor.

Finally, the algorithm is proven with the NEDC B1 cycle and is shown in Figure 6.9 for a window of the cycle. The described conditions and data for previous tests are still valid here.  $\hat{x}_{\lambda^{-1}}$  keeps the model dynamics but evolves correcting the drift and relying on the steady-state value of the sensor. This



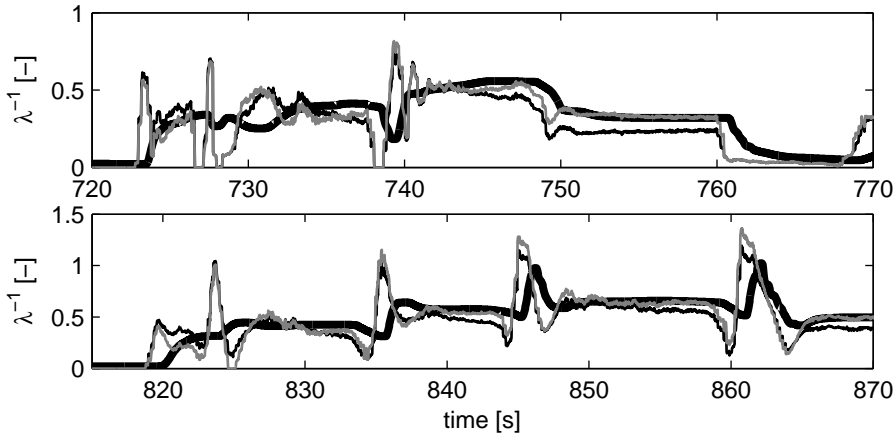


**Figure 6.7.**  $\hat{x}_{\lambda^{-1}}$  performing injection steps at 1500 rpm and using an adaptive look-up table for modelling  $\theta_{\lambda^{-1}}$ . Legend: Thick black is the sensor  $z_{\lambda^{-1}}$ , thin black is the model  $x_{\lambda^{-1}}$  and gray is  $\hat{x}_{\lambda^{-1}}$ . Values for tuning the KF are  $\sigma_w^2/\sigma_v^2 = 1/80$ .



**Figure 6.8.**  $\hat{x}_{\lambda^{-1}}$  using table and freezing conditions performing different injection steps at 1500 rpm.

test is representative of real driving conditions and the procedure demonstrates its feasibility for being used in commercial vehicles.



**Figure 6.9.** Top and bottom:  $\hat{x}_{\lambda^{-1}}$  using the table and applying freezing conditions to the NEDC B01 cycle.  $z_{\lambda^{-1}}$  is the sensor output (thick black),  $x_{\lambda^{-1}}$  is the model output (thin black) and  $\hat{x}_{\lambda^{-1}}$  is the observation (gray). The drift is corrected while the dynamics of the observation  $\hat{x}_{\lambda^{-1}}$  are maintained.

### 6.3 Fast estimation of $\text{NO}_x$

The attention is turned out to the problem of the online  $\text{NO}_x$  estimation. With an algorithm like B, an adaptive  $\text{NO}_x$  map can be used in order to infer  $\text{NO}_x$  as function of only  $m_f$  and  $n$  as in (4.8), which is justified if the transitions are slow enough for utilising a quasi-static approach (plus a filter for the sensor dynamics). An algorithm C permits to update  $\text{NO}_x$  model parameters and look-up tables. Finally, and for both designs, the algorithm A and the drift correction algorithm could be used in order to track the model bias.

This section also completes the comparison of the different methods for updating look-up tables already presented in Chapter 5 by testing the cycle SDMP B5 and the algorithm B. Afterwards, the adaptive tuning of the RT  $\text{NO}_x$  model is proposed. On the basis of the presented results, the estimation of the actual  $\text{NO}_x$  is proposed.

### 6.3.1 Online updating of look-up tables for modelling $\text{NO}_x$

In a first approach,  $\text{NO}_x$  is modelled by using a 2D adaptive look-up table function of  $n$  and  $m_f$ , whose parameters are estimated with the updating methods. The state-space model (5.33) is used for designing an observer like the algorithm B presented in Figure 6.1.

The model parameters are estimated online while the engine is running without any special calibration procedure or test rig, beyond the sensor measurements. The first time that the engine is running, parameters evolve, and when the engine switches off, the stored parameters can be used for predicting  $\text{NO}_x$ . When the engine is running again, the parameters keep evolving for correcting drift and slowly varying effects. Furthermore, the observer built for updating the map can be profited for having an actual  $\text{NO}_x$  estimation, i.e. avoiding the filtering and the delay of the sensor. The SKF updating method is appropriate for online usage because of the light computational burden and the estimation capabilities.

Moreover, this configuration might be applied to estimate local models for defining the  $\text{NO}_x$  emissions as function of  $u_{soi}$  and  $u_{egr}$ , and for different engine operating points [3,4]. These maps are useful for the joint air and fuel path control [5].

With respect to the dimensions of the maps used for the adaptive estimation of  $\text{NO}_x$  and according to the author's experience, usual dimensions for look-up tables in production engines are between 200 and 400 parameters. The selection of the appropriate dimensions is a trade-off between the accuracy and the computational resources. An excessive density can compromise the table convergence, but a higher density of scheduling points can improve the accuracy of the solution.

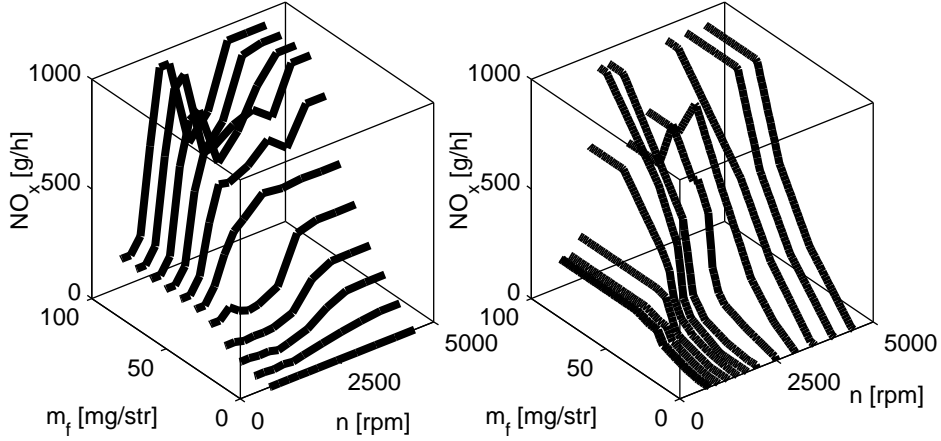
Figure 6.10 shows the nominal  $\text{NO}_{x,0}$  map used for the  $\text{NO}_x$  model in Chapter 4. This map was filled with an intensive DOE comprising 363 steady-state tests. The  $\text{NO}_x$  sensitivity depends on the engine area, with a clear difference between the EGR area, transition and the areas without EGR. Note that all elements of  $\text{NO}_{x,0}$  are filled in order to avoid numerical problems, indeed areas that are not reached for the engine (out of the full load line).

The grid used in this work for the observation is defined as follows

$$r = [750 \quad 1000 \quad 1250 \quad 1500 \quad 1750 \quad 2000 : 500 : 4500] \quad (6.7a)$$

$$c = [0 : 5 : 30 \quad 34 : 4 : 50 \quad 55 : 5 : 80] \quad (6.7b)$$

where  $r$  represents the grid for  $n$  [rpm] and  $c$  stands for  $m_f$  [mg/str].



**Figure 6.10.**  $NO_x$  variation as function of  $n$  and  $m_f$  in the directions. Left plot: along  $n$  direction. Right plot: along  $m_f$  direction.

### 6.3.1.1 Comparison of the updating methods

The SDMP B5 cycle is used for comparing the algorithms with real engine data.  $\Theta_{NO_x,0}$  is the null matrix and is updated with the cycle. The state-space model (5.33) is built with the following numerical values

$$a_{NO_x} = 0.96; \quad \tau_{NO_x} = 0.75s \quad (6.8)$$

while the sampling frequency is 50 Hz ( $T_s = 20ms$ ). Sensor parameters come from the methods presented in Chapter 3 by performing SOI steps. The state-space model output  $z_{NO_x}$  is given by the  $NO_x$  sensor, while the inputs are

$$u_{1,k} = z^{-\tau_{NO_x}/T_s} n(k) \quad (6.9a)$$

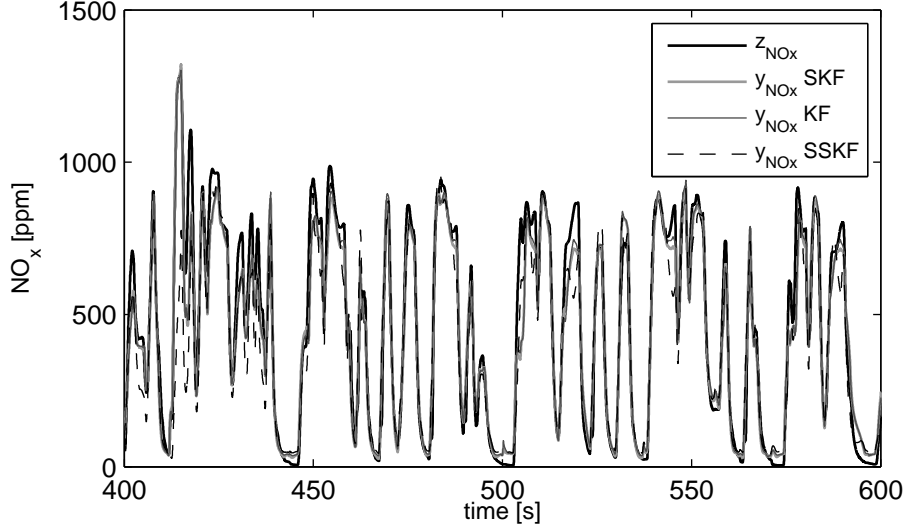
$$u_{2,k} = z^{-\tau_{NO_x}/T_s} m_f(k) \quad (6.9b)$$

and the state vector  $x$  is

$$x_k = [vec(\Theta_{NO_x}) \quad x_{NO_x f}]_k^T \quad (6.10)$$

The filter is tuned with the values

$$\sigma_v^2 = 25^2; \quad \sigma_w^2 = 15^2; \quad \sigma_f^2 = 50^2 \quad (6.11)$$



**Figure 6.11.**  $\text{NO}_x$  prediction for the three methods by performing the SDMP B5 cycle. The maps are updated during 400 s, i.e. a subset of 20000 measurements of  $z_{\text{NO}_x}$ .  $y_{\text{NO}_x}$  is simulated with the updated tables at 400 s  $\Theta_{\text{NO}_x,400}$ . The sensor measurement  $z_{\text{NO}_x}$  is provided for comparison.

and a test subset of  $t = 400\text{s}$  is used for updating the adaptive map and the look-up table  $\Theta_{\text{NO}_x,400}$  is used for predicting  $\text{NO}_x$  in the whole cycle plus considering the first order filter that models the sensor response.

$$y_{\text{NO}_x} = \frac{1 - a_{\text{NO}_x}}{1 - a_{\text{NO}_x} z^{-1}} \Theta_{\text{NO}_x,400}(\mathbf{u}_{1,k}, \mathbf{u}_{2,k}) \quad (6.12)$$

The KF, SKF and SSKF methods are used for updating the model and the results can be seen in Figure 6.11. All three methods behave well and are capable of predicting  $\text{NO}_x$  emissions. Dynamics assumptions are also good enough for having a good fitting in the SDMP. Note that the SSKF results are also acceptable as the covariance tracking is not relevant in this cycle as it is in slow varying tests. The fitting should be optimized by independently tuning the filter for each method.

It is worth comparing the convergence of the methods and for this, the SDMP cycle is also used as benchmark for testing the absolute mean error when the updating time is varied monotonically around the cycle

$$e_M(k) = \frac{\sum_{i=1}^{n_y} (z_{\text{NO}_x}(i) - y_{\text{NO}_x}(i)|_{t=kT_s})}{kT_s} \quad (6.13a)$$

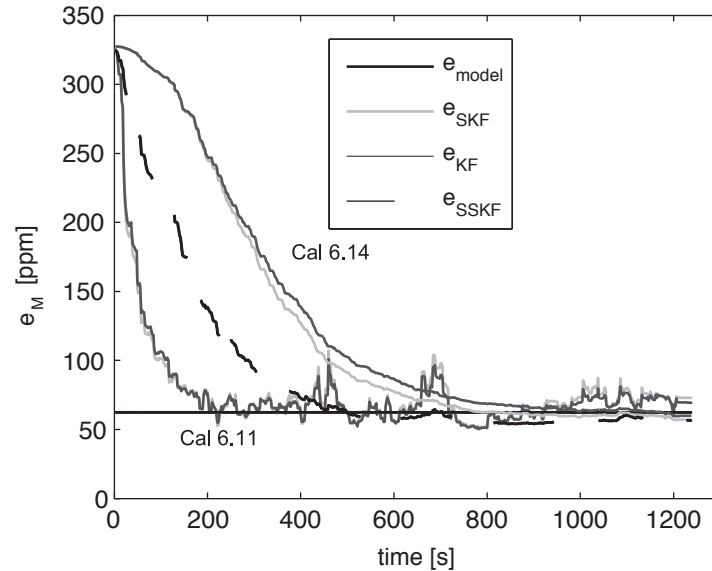
where  $t$  is the time that the table has been updated and  $y_{\text{NO}_x}$  is calculated by (6.12). The results are shown in Figure 6.12. Focusing the attention in the lines generated with the filter calibration Cal (6.11),  $e_M$  tends to be lower as  $t$  grows at least during first iterations when the system knowledge is poor. The horizontal line shows the benchmark model error, which is constant as the model calibration is fixed. It is clear how all the three methods converge to an error similar to that of the benchmark model, but with the advantage that the drift and ageing are considered by the online versions but not by the model. The KF and SKF behaviour is quite similar as expected. On the other hand, Figure 6.12 also shows that the KF and the SKF are faster than the SSKF and this is due to the filter tuning. Anyway, there exists a trade-off between the convergence speed and the estimation robustness, e.g. the KF and SKF have a significant oscillation with respect to the SSKF around the benchmark model line. This is a matter of the noise tuning and the results are also generated by using a lower gain observer with the calibration set

$$\sigma_v^2 = 25^2; \sigma_w^2 = 7^2; \sigma_f^2 = 50^2 \quad (6.14)$$

for the KF and the SKF. The results are shown in lines generated by Cal (6.14) in Figure 6.12. With this configuration, the convergence speed is lower but the noise transmission and overfitting is avoided when  $t$  is large. This goes in the direction of the robustness and the optimised calibration data set must solve this trade-off, considering uncertainties in the sensor behavior knowledge or in the model quality and data-set quality, among others. Strategies such as the freezing applied to the  $\lambda^{-1}$  estimation may also be applied to this problem.

With respect to the global observability, variances of the 4 table parameters are compared for both the KF and SKF using the calibration (6.14). Figure 6.13 shows the results. The top plot shows an element that is never observed during first 200 s and due to this, the variance increases monotonically and  $x_{126}$  is still null in  $t = 200$ . Second and third elements represent two elements that are active at some instants. This is clearly seen when the variance decreases, while when the elements are not active, the variance increases again monotonically. Note that in a production system application, that will run for the life time of a vehicle, the elements of the estimation error covariance matrix need to be limited so that they do not grow too much and cause numerical problems, see e.g. Höckerdal *et al.* [6] who propose a saturation for avoiding too large variances. The bottom plot shows the variance related to  $x_{\text{NO}_x f}$  that is fairly constant and is a proof of the system global observability.

The variances by using the KF and the SKF behave similar and differences can only be found when zooming in the Figure 6.13. Therefore, the

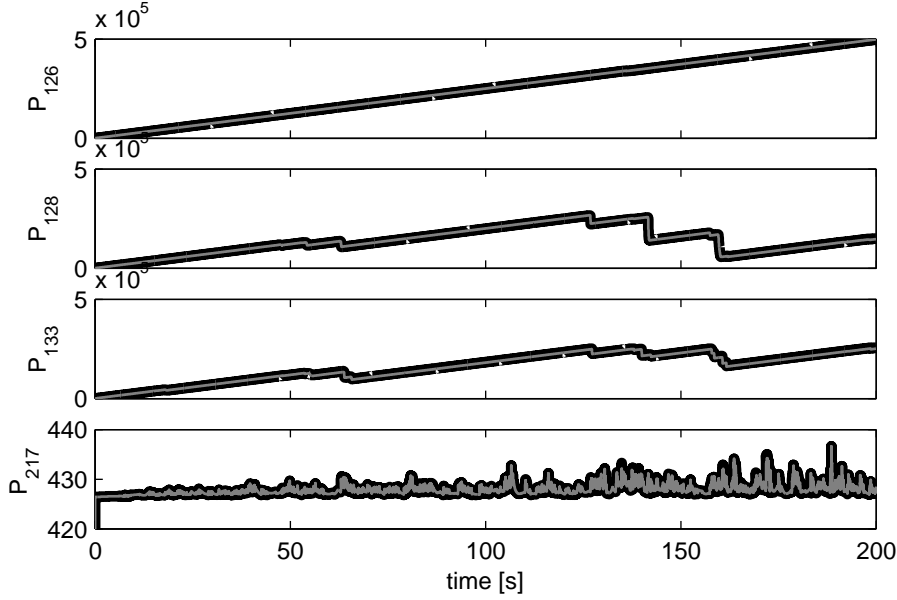


**Figure 6.12.** Evolution of the mean of the absolute error  $e_M$  for the different methods in the SDMP B5 cycle and for two different filter calibrations. SSKF dashed line is calculated by using the filter calibration (6.11). Lines left to the SSKF one: using calibration (6.11). During first iterations, the error is high, as far as the table has no initial knowledge, but once that the time is around 400 seconds, the error is minimum, although the noise transmission is evident. Lines right to the SSKF one: using calibration (6.14), noise transmission is reduced but the convergence speed as well. The chosen filter tuning must be a trade-off between the robustness and convergence speed.

SKF method performs quite similar to the standard KF but with a much lower computational burden. Due to the SDMP presents sharp transients, the SSKF method also performs quite well, and proves that this method can be useful when the considered data set fulfils some conditions in order to ensure robustness. To the nature of the application considered here, the SKF is the best solution for online updating of maps.

### 6.3.1.2 Adaptive maps for predicting $\text{NO}_x$

The SDMP is a test with sharp variations on the operating point conditions but the homologation cycles, such as the NEDC, are much slower. This could compromise the global observability since higher levels of excitation are beneficial for updating. As shown in Figure 4.3, the air path dynamics are fast enough to follow the load variations in the NEDC and then  $\dot{m}_a$  and  $p_{boost}$  are able to track their references. In such case, the model (4.8) is valid and



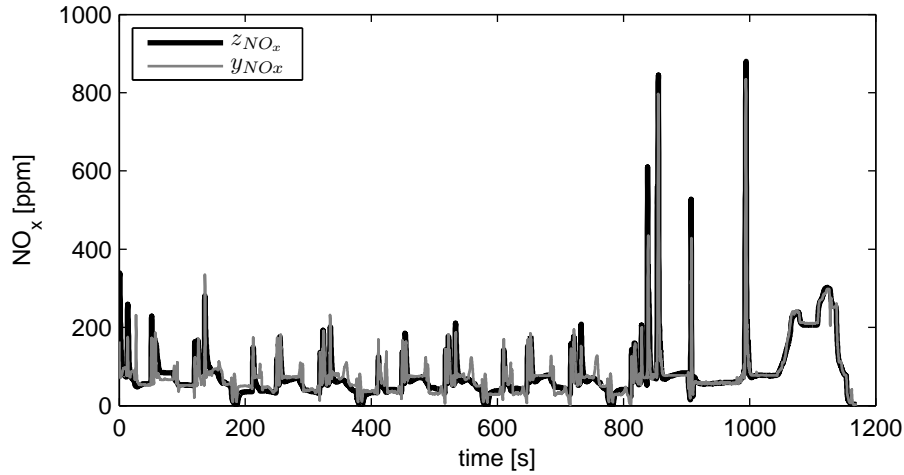
**Figure 6.13.** Variance analysis by updating a  $\text{NO}_x$  map in the cycle SDMP B5. From top to bottom, variances of elements 126, 126, 133 and 217 are plotted. Black line is the KF and gray line is the SKF.

the table  $\Theta_{\text{NO}_x}$  might directly replace the nominal  $\text{NO}_{x,0}$  table in the RT  $\text{NO}_x$  model (see (4.22)). Despite  $\text{NO}_{x,0}$  could be used for initializing  $\Theta_{\text{NO}_x}$ , here the null matrix is the initial map, which is a worst-case condition.

The NEDC A1 cycle is run for updating the map. Figure 4.19 already showed the RT  $\text{NO}_x$  model output when comparing with  $z_{\text{NO}_x}$  in that cycle, where a clear drift exists. After using an adaptive map and the SKF method with the calibration (6.14), the results are shown in Figure 6.14. There,  $y_{\text{NO}_x}$  is calculated on the basis of a table  $\Theta_{\text{NO}_x}$  which has been updated during the whole cycle. After comparing with  $z_{\text{NO}_x}$  and  $x_{\text{NO}_x}$  coming from the  $\text{NO}_x$  model (see Figure 4.19), the adaptive map approach gets a good  $\text{NO}_x$  estimation which is slightly better than the one of the  $\text{NO}_x$  benchmark model.

The SKF could be used for the online adaptation and/or calibration of complex models, where a large number of maps and parameters must be updated. Anyway, a deep study is required for ensuring the observability, convergence and robustness properties and for getting a computationally efficient learning structure. A big number of parameters and maps should be updated and the





**Figure 6.14.** Adaptive  $NO_x$  estimation in the NEDC A01, where it is estimated with a table that is updated during the whole cycle and then used for simulating  $y_{NO_x}$ .  $z_{NO_x}$  is provided for comparison. The estimation keeps the dynamic properties of  $z_{NO_x}$  but after updating, the drift in Figure 4.19 is cancelled.

problem is non-convex. Admittedly the computational issue is critical and there the SKF can be an effective solution.

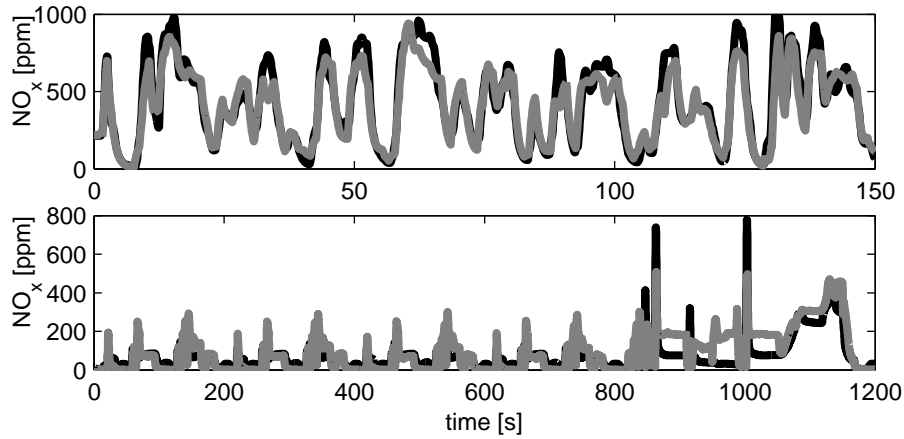
In the case of a fast cycle, i.e. the SDMP, the EGR valve is closed the most of the time. Therefore, the states are function of  $n$  and  $m_f$  for the engine settings corresponding to a closed EGR valve

$$x = f(n, m_f, m_a^*(n, m_f), p_b^*(n, m_f))_{u_{egr}=100} = f(n, m_f)_{u_{egr}=100} \quad (6.15)$$

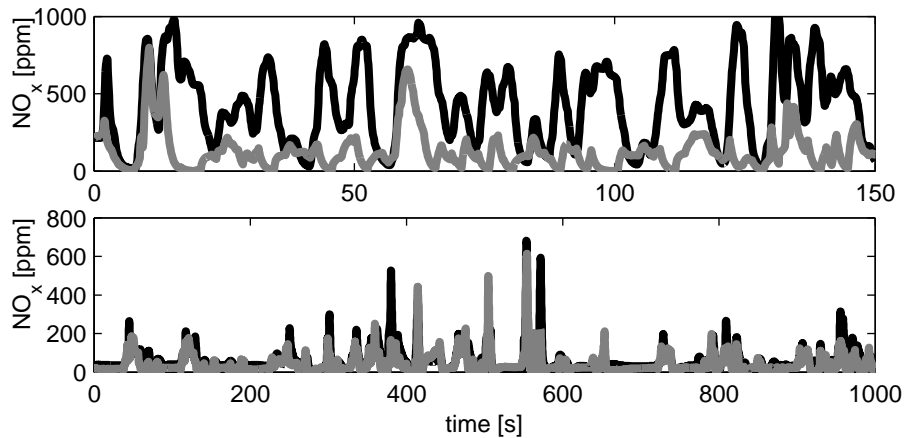
In the case of the NEDC or CADC cycles, the transients are much slower and the engine works closer to the nominal settings. In such case, the states are also function of  $n$  and  $m_f$  but for the nominal operation of the EGR-VGT controller. Possible states can now be approximated with

$$x = f(n, m_f, m_a^*(n, m_f), p_b^*(n, m_f))_0 = f(n, m_f)_0 \quad (6.16)$$

This point links with the discussion of Section 4.3.1 about how to represent the engine conditions. Figure 6.15 shows the predicted  $NO_x$  emissions in the cycles SDMP B6 (top plot) and NEDC B1 (bottom plot) after using a map  $\Theta_{NO_x}$  that has been updated in the cycle SDMP B5. The table output is filtered and delayed by a sensor model with the parameters  $a_{NO_x} = 0.96$  and  $\tau_{NO_x} = 0.75s$ . The ability to reproduce  $NO_x$  on the SDMP B6 is clear while a



**Figure 6.15.** Offline  $\text{NO}_x$  prediction  $y_{\text{NO}_x}$  (gray) against  $\text{NO}_x$  sensor (black) by using a table  $\Theta_{\text{NO}_x}$  updated online in the SDMP B5 cycle. Top plot: results in the SDMP B6 cycle. Bottom plot: results in the NEDC B1 cycle.



**Figure 6.16.** Offline  $\text{NO}_x$  prediction  $y_{\text{NO}_x}$  (gray) against  $\text{NO}_x$  sensor (black) by using a table  $\Theta_{\text{NO}_x}$  updated online in the NEDC B1 cycle. Top plot: results in the SDMP B6 cycle. Bottom plot: results in the CADX B2 cycle.

slight drift exists in the NEDC B01. Anyway, the NO<sub>x</sub> estimation during the urban part is not as bad as in the highway part, where NO<sub>x</sub> is overestimated as expected ( $\Theta_{\text{NO}_x}$  has been fitted with nearly zero EGR conditions).

However, if updating  $\Theta_{\text{NO}_x}$  with the NEDC B1 cycle and then using the results for predicting the SDMP B6, shown in top plot of Figure 6.16, the results are quite bad as a big area of the engine map is not excited and the covered areas present lower NO<sub>x</sub> emissions because of the EGR actuation. The prediction using  $\Theta_{\text{NO}_x}$  in the CADC B2 urban cycle is quite better, as the covered areas are similar.

### 6.3.2 Online updating of the NO<sub>x</sub> model

In Chapter 4, a NO<sub>x</sub> model was developed based on a set of look-up tables for representing nominal engine conditions with several corrections based on the estimated intake oxygen, thermal loading and others as in (4.22). In this section, adaptive capabilities are proved in the NO<sub>x</sub> model [7], and the new state-space vector  $X^w$  might be augmented as follows

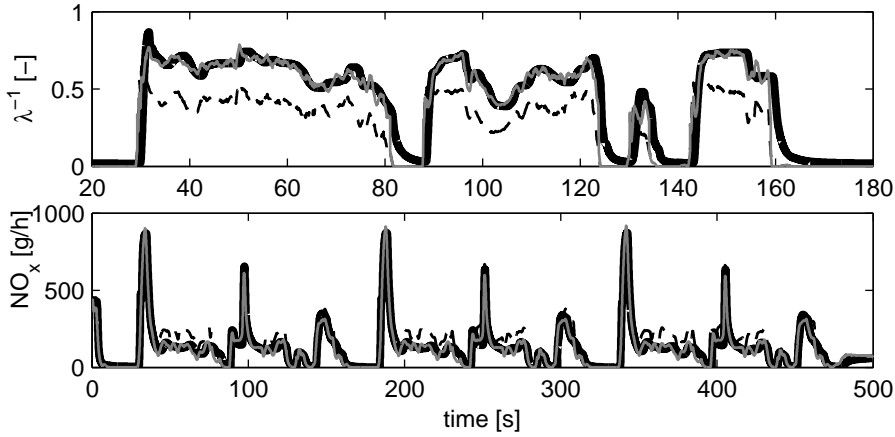
$$X^w = [x_{\text{NO}_x} \quad y_{\text{NO}_x} \quad x_\lambda^{-1} \quad y_\lambda^{-1} \quad \text{vec}(\Theta_{\text{NO}_x}) \quad \theta]^T \quad (6.17)$$

where besides  $x_{\text{NO}_x}$  and  $y_{\text{NO}_x}$ , the following states are included:  $x_\lambda^{-1}$  and  $y_\lambda^{-1}$  representing the actual  $\lambda^{-1}$  and the filtered  $\lambda^{-1}$  (for comparing with the  $\lambda^{-1}$  output from the sensor as was made in (6.12) for NO<sub>x</sub>) respectively and used as inputs for the RT NO<sub>x</sub> model, the nominal NO<sub>x</sub> map  $\Theta_{\text{NO}_x}$  and the model bias  $\theta_{\text{NO}_x}$ .

On one side, the use of an observer for  $\lambda^{-1}$  if the NO<sub>x</sub> sensor is available is completely justified due to the NO<sub>x</sub> sensitivity to errors on mistaking the intake oxygen [8] (see Table 4.2). Furthermore, the error sources in the  $\lambda^{-1}$  model were already explained in Section 4.2 and an example of  $\lambda^{-1}$  observation is shown for the NEDC in Figure 6.9 where the model drift is not negligible. Figure 6.17 precisely compares the model output  $y_{\text{NO}_x}$  when a drift on  $x_\lambda^{-1}$  estimation occurs. The top plot shows  $z_{\lambda^{-1}}$  by a ZrO<sub>2</sub> sensor,  $x_\lambda^{-1}$  from (4.1) and the observation  $\hat{x}_{\lambda^{-1}}$  by a KF, when a clear drift on  $x_\lambda^{-1}$  exists<sup>1</sup>. If the drifted signal is used for inferring NO<sub>x</sub> by the model, the output is also drifted from the sensor signal as shown in the bottom plot. However, if observing  $\hat{x}_{\lambda^{-1}}$

<sup>1</sup>The drift on  $x_\lambda^{-1}$  is caused due to failures in the air mass flow or the injected fuel mass signals. This drift will affect to the EGR flow model and the look-up tables interpolation. In those cases, the sensor signal  $z_{\lambda^{-1}}$  can be used for correcting errors on the air mass flow and/or injected fuel mass signals. In this section and for the simulation, the error is only considered in the  $x_\lambda^{-1}$  model.

by a KF, the  $NO_x$  output keeps the model accuracy. Therefore, the benefits of observing  $\lambda^{-1}$  are clear.



**Figure 6.17.**  $\lambda^{-1}$  and  $NO_x$  estimations in a TRAN B ( $T_{cool} = nom$ ,  $\Delta u_{egr} = 0\%$ ,  $\Delta p_{boost} = 0$ ). Top plot:  $z_{\lambda^{-1}}$  sensor signal (black),  $x_{\lambda^{-1}}$  calculated by the model (dashed thin black) and  $\hat{x}_{\lambda^{-1}}$  observed by a drift augmented model (gray). Bottom plot:  $NO_x$  estimation by three ways,  $z_{NO_x}$  sensor signal (black),  $y_{NO_x}$  by the  $NO_x$  model using  $x_{\lambda^{-1}}$  as input (dashed thin black) and  $y_{NO_x}$  by using an adaptive model with  $\hat{x}_{\lambda^{-1}}$  as  $\lambda^{-1}$  input.

On the other side, model errors and ageing can be corrected online by tracking the bias in the same way as explained in Section 6.2 for  $\lambda^{-1}$ , or by observing the nominal  $NO_{x,0}$  table for also improving the prediction capabilities of the model. The former is useful for actual  $NO_x$  prediction, while the latter is advisable for ensuring the model accuracy.

**A learning algorithm for updating the  $NO_x$  model.** The state-space vector

$$X^w = [x_{NO_x} \quad y_{NO_x} \quad x_{\lambda^{-1}}^{-1} \quad y_{\lambda^{-1}}^{-1} \quad vec(\Theta_{NO_x})]^T \quad (6.18)$$

is built for designing an algorithm like C (see Figure 6.1) in order to adapt the nominal  $NO_x$  table ( $NO_{x,0}$ ) to possible changes in the  $NO_x$  emissions, mainly due to the ageing. Nevertheless, this structure is useful not only for learning  $NO_{x,0}$  where

$$\Theta_{NO_x} = NO_{x,0} \quad (6.19)$$

but for fitting a table without initial knowledge. Depending on the model and sensor uncertainties, the strategy of the filter tuning can be conservative (slow corrections) or aggressive (fast corrections).

The new state-space model remains similar to that of (5.33), used for the observer B, but including NO<sub>x</sub> model variations and corrections around the nominal values (see (4.22)) in the output matrix  $H_k$

$$H_k^w = [\text{zeros}(\text{length}(\text{vec}(\Theta_{\text{NO}_x}))) \quad C_{\text{mod}}] \quad (6.20)$$

where

$$C_{\text{mod}} = e^{-k_{\text{NO}_x} \cdot (EGR\lambda^{-1} - EGR_0\lambda_0^{-1})} \cdot C_{\text{NO}_x} \cdot C_{\text{dyn}} \quad (6.21)$$

Other model maps and parameters, such as the  $\eta_v$ , might be updated by just applying an EKF similar to that applied to NO<sub>x,0</sub> (with the correspondent linearisation). If a full adaptive model is proposed, in a way that all model parameters are observed, then the learning structure would lack of robustness and it would be difficult to discriminate the observed error between all tables, curves and coefficients. An alternative strategy could be to observe a drift  $\theta_{\text{NO}_x}$  and afterwards weighting the error between model the parameters by LS or other strategies. Other limitations are the computational and memory resources or the lack of awareness in the error sources, making inadvisable to design an strategy based on the together updating of all model parameters and tables [9].

$\Theta_{\text{NO}_x}$  is designed with the grid defined by (6.7) and is updated with the SKF method. The RT NO<sub>x</sub> model is simulated with the tests of the campaign B and the errors are computed. Due to the system ageing, the model presents a drift. The SKF method is also used for updating NO<sub>x,0</sub> coming from the NO<sub>x</sub> model and thus cancelling the drift. The updated model is simulated in those cycles and the new errors are computed.

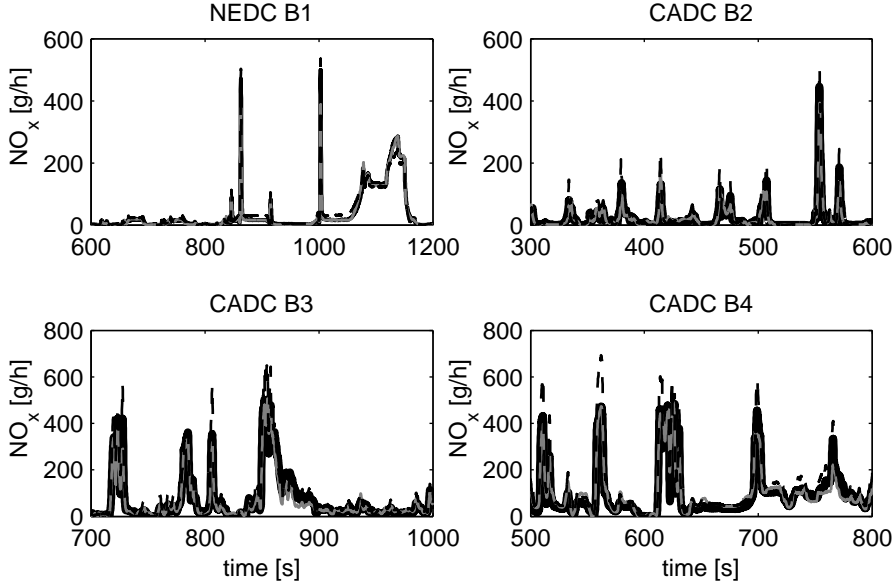
For the application, the NO<sub>x</sub> sensor output  $z_{\text{NO}_x}$  is converted from ppm to g/h by using (4.15) and the updating method is run with the output in g/h in order to directly modify NO<sub>x,0</sub> table.

The noise calibration is adjusted by the trial and error method and is not guaranteed to be the optimised one

$$\sigma_v^2 = 1; \sigma_w^2 = 0.01; \sigma_f^2 = 1 \quad (6.22)$$

The tests in the test campaign B are run by maintaining the calibration (6.22), and the offline NO<sub>x</sub> prediction is calculated just after finishing the cycle by using the updated NO<sub>x,0</sub> in the NO<sub>x</sub> model. The computed errors are compared with the ones of the model. The cycles used for the updating are the B1 to B6, totalling 6 tests. In these simulations, the initial matrix is NO<sub>x,0</sub> fitted for the NO<sub>x</sub> model in Chapter 4. Figure 6.18 shows the results

on 4 different cycles, where the original and updated  $y_{NO_x}$  are compared with the sensor signal  $z_{NO_x}$ .  $y_{NO_x}$  with the original calibration presents a bias with respect to  $z_{NO_x}$  that is corrected after running the cycles with the SKF.



**Figure 6.18.** Comparison of the  $NO_x$  estimation by three different sources in 4 different cycles: sensor signal  $z_{NO_x}$  (black),  $y_{NO_x}$  by using the RT  $NO_x$  model (dashed thin black) and  $y_{NO_x}$  computed by updating the  $NO_{x,0}$  table with the SKF method (gray). Left to right and top to bottom: NEDC B1, CADC B2, CADC B3 and CADC B4 cycles.

Table 6.1 compares the absolute mean errors by using the original model output and the predictions of the updated model

$$e = \text{mean}\left(\sum_{k=1}^{k=l_{end}} z_{NO_x}(k) - y_{NO_x}(k)\right) \quad (6.23)$$

where  $l_{end}$  defines the number of samples taken for the considered cycle. The errors can be minimised achieving quite better results by using the learning structure. These results validate the SKF method and the capabilities for updating models.

With respect to the filter tuning, it has not been optimised for obtaining the results, and a trial-and-error method has been followed. Further improvements are in the line of optimising the tuning by improving the sensor model, e.g.

**Table 6.1.** Mean absolute errors (g/h) in the  $\text{NO}_x$  model output for different tests, and comparing the error of the original model (Model column) with the ones of the updated model (Updated column). The cycles are run one time and the matrix  $\text{NO}_{x,0}$  is updated online.

Cycle	Model	Updated
NEDC B1	6.14	3.08
CADC Urban B2	6.40	3.75
CADC Rural B3	14.81	10.46
CADC Highway B4	36.23	22.51
SDMP B5	49.46	49.13
SDMP B6	46.34	43.64

including a variable delay, or augmenting the state vector with other model parameters. Anyway, the obtained results show the capabilities of the method for its online implementation in order to re-tune models and fit tables and parameters.

### 6.3.3 Online observation of the actual $\text{NO}_x$

The actual  $\text{NO}_x$  can be predicted by utilizing the state-space models and tables properly. For that, the state vector is augmented with an extra-state for representing the actual  $\text{NO}_x$ , i.e.  $\text{NO}_x$  emissions before sensor filtering and correcting the delay. Estimating the actual  $\text{NO}_x$  is advisable for RT purposes, such as on-board  $\text{NO}_x$  control or diagnosis.

One of the important problems when of working on RT systems is the causality of the solution, as far as sensors usually measure in a delayed basis, mainly due to the transport and hardware delays. In this work, two possibilities are described for solving this issue: tracking the bias and the use of a model learning structure. Both possibilities are similar, and are based on the ageing cancellation by means of adaptive filtering.

**Actual  $\text{NO}_x$  by drift correction models.** The following state-space vector is built

$$X^w = [\theta_{\text{NO}_x} \quad x_{\text{NO}_x} \quad x_{\text{NO}_x f}]^T \quad (6.24)$$

where now  $x_{\text{NO}_x}(k) \in \mathbb{R}$  is added to track the actual  $\text{NO}_x$ , the model input  $u_k = x_{\text{NO}_x}(k - \tau_{\text{NO}_x}/T_s)$  is conveniently delayed and  $z_k = z_{\text{NO}_x}(k)$ . Matrices

$\mathbf{F}$ ,  $B$  and  $H$  are modified accordingly

$$\mathbf{F} = \begin{bmatrix} 1 & 0 & 0 \\ 1 & 0 & 0 \\ 0 & 1 - a_{\text{NO}_x} & a_{\text{NO}_x} \end{bmatrix} \quad B = \begin{bmatrix} 0 \\ 1 \\ 1 - a_{\text{NO}_x} \end{bmatrix} \quad H = [0 \quad 0 \quad 1] \quad (6.25)$$

Due to the sensor delay, the bias  $\hat{\theta}_{\text{NO}_x}(k - \tau_{\text{NO}_x}/T_s)$  at the instant  $k$  is observed with a certain delay  $\tau_{\text{NO}_x}$ , and that means that at a given time, the actual drift will be calculated with the delay  $\tau_{\text{NO}_x}$

$$\hat{x}_{\text{NO}_x}(k) = x_{\text{NO}_x}(k) + \hat{\theta}_{\text{NO}_x}(k - \tau_{\text{NO}_x}/T_s) \quad (6.26)$$

However, the actual  $\text{NO}_x$  estimation should be

$$\hat{x}_{\text{NO}_x}(k) = x_{\text{NO}_x}(k) + \hat{\theta}_{\text{NO}_x}(k) \quad (6.27)$$

which corresponds to a non-causal operation.

Using the delayed  $\hat{\theta}_{\text{NO}_x}(k - \tau_{\text{NO}_x}/T_s)$  could lead to important errors when sharp load transients. Alternatively, an adaptive map  $\Theta_{\text{NO}_x} \in \mathbb{R}^2 : [n, m_f] \rightarrow \theta_{\text{NO}_x}$  could be introduced for modelling the drift as a function of the operating point conditions, as was made for  $\lambda^{-1}$ . This option is well described in Chapter 5: a map  $\Theta_{\text{NO}_x}$  is adapted by means of a steady-state version of a KF (similar to a RLS filtering) and is interpolated in a 2D basis for giving an initial estimate in the prediction of the state, in such way that the KF is modified as follows

$$\hat{x}_k = \mathbf{F}\Theta_{\text{NO}_x, \mathbf{k}}(n(k), m_f(k)) + Bu_k + K_{\infty} (z_{\text{NO}_x}(k) - H(\mathbf{F}\hat{x}_{k-1} + Bu_k)) \quad (6.28)$$

and the table  $\Theta_{\text{NO}_x}$  is independently updated from the KF in order to store  $\theta_{\text{NO}_x}$

$$\Theta_{\text{NO}_x, \mathbf{k}} = f(\Theta_{\text{NO}_x, \mathbf{k}-1}, \hat{x}_k, z_{\text{NO}_x}(k)) \quad (6.29)$$

The map itself is interpolated for inferring  $\theta_{\text{NO}_x}$  avoiding the causal problem and huge integrations when the drift varies with the operating point conditions and time. The final estimation is then calculated as follows

$$\hat{x}_{\text{NO}_x}(k) = x_{\text{NO}_x}(k) + \Theta_{\text{NO}_x, \mathbf{k}}(n(k), m_f(k)) \quad (6.30)$$

**Actual  $\text{NO}_x$  by updating look-up tables.** The state-space model for learning maps is useful not only for calibrating the map but also for the RT



observation of the NO<sub>x</sub> signal. Therefore, the matrices and the state vector in (5.33) are also modified and augmented for introducing  $x_{\text{NO}_x}$

$$X^w = [\text{vec}(\Theta_{\text{NO}_x}) \quad x_{\text{NO}_x} \quad x_{\text{NO}_x f}]^T \quad (6.31)$$

with the local observable matrix  $\mathbf{F}_k^o$

$$\mathbf{F}_k^o = \left[ \begin{array}{c|cc} \mathbf{I}_{4 \times 4} & 0 & 0 \\ \hline C_{mod} q_k & 0 & 0 \\ 0 & (1 - a_{\text{NO}_x}) & a_{\text{NO}_x} \end{array} \right] \quad (6.32)$$

For the case of the algorithm B for updating only look-up tables,  $C_{mod} = 1$  (model (4.8)). For the case of an algorithm like C and for the NO<sub>x</sub> model,

$$C_{mod} = e^{-k_{\text{NO}_x} \cdot (EGR\lambda^{-1} - EGR_0\lambda_0^{-1})} \cdot C_{\text{NO}_x} \cdot C_{dyn} \quad (6.33)$$

with the parameters calibrated offline. Finally,  $H^w$  is now constant for both algorithms B and C

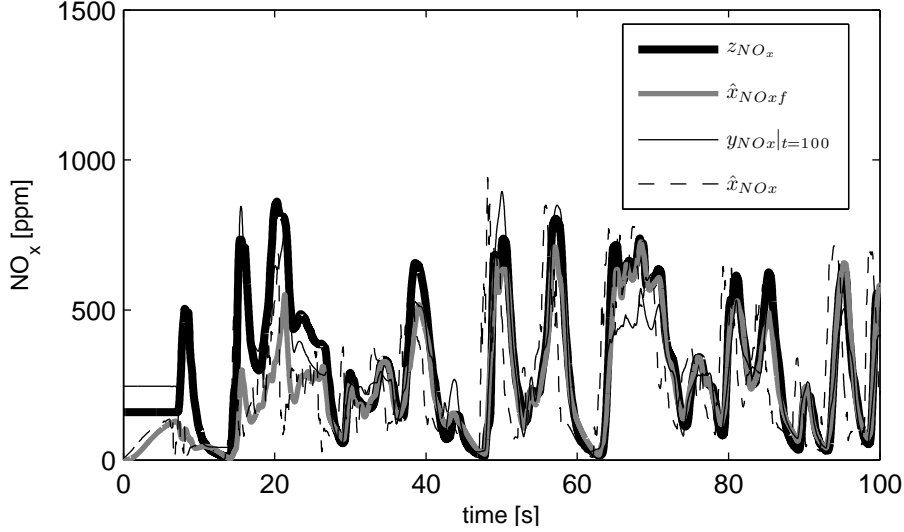
$$H^w = [0 \quad 0 \cdots 1] \quad (6.34)$$

The use of the table allows to solve the non-causalities related with the RT estimation in a smart way: note that the observer is built on the basis of delayed inputs.

The use of a drift correction model or adaptive look-up tables NO<sub>x</sub> leads to similar results, but the look-up tables structure is beneficial as far as can be used also for updating the model. In order to decouple the filter tuning for updating the look-up table and the output drift correction, the two methods could be combined by observing both the table parameters, the bias and the filtered output, designing a state-space vector

$$X^w = [\text{vec}(\Theta_{\text{NO}_x}) \quad \theta_{\text{NO}_x} \quad \mathbf{x}_{\text{NO}_x} \quad \mathbf{x}_{\text{NO}_x f}]^T \quad (6.35)$$

**Results.** The model with the state-space vector (6.31) is observed in order to show an application example of the fast estimation of the actual NO<sub>x</sub>, taking into account that this algorithm can be adapted to other applications. Therefore, the look-up tables learning structure can be utilised for updating models and also estimating actual signals. Figure 6.19 shows results in the cycle SDMP B5, by fitting the NO<sub>x</sub> model NO<sub>x,0</sub> map with initial null values



**Figure 6.19.**  $NO_x$  estimation by three ways:  $z_{NO_x}$  sensor measurement, online observation of the sensor signal by the SKF method  $\hat{x}_{NO_xf}$ , offline prediction  $y_{NO_x}$  by using the map  $\Theta_{NO_x,100}$  and online actual estimation  $\hat{x}_{NO_x}$ .

and keeping the calibration set (6.14).  $\hat{x}_{NO_xf}$  represents the observation of  $z_{NO_x}$  by using the adaptive map at every iteration. The adaptive map provides a perfect fitting in about 30 seconds (it might be accelerated varying the filter calibration but it is not advisable if the sensor model is not perfectly identified). Alternatively,  $y_{NO_x}|_{t=100}$  shows the offline  $NO_x$  sensor prediction using the model  $\Theta_{NO_x,100}$  with acceptable results.

The table itself might predict the actual  $NO_x$  emissions if interpolating the table directly applying no delay nor filtering. Furthermore, the actual  $NO_x$  is directly observed with the state-space system for learning having a compact programming and providing an adaptive estimation. For that,  $x_{NO_x}$  is included in the state vector. This does not compromise the assumptions made for the SKF. Coming back to Figure 6.19,  $\hat{x}_{NO_x}$  is the online observation of the actual  $NO_x$ . This signal is advised for using it as a valid reference of the actual  $NO_x$ .

## 6.4 Conclusions

This section has presented different applications for the methods developed in Chapter 5 applied to the fast estimation of  $\lambda^{-1}$  and  $NO_x$ .

For the case of  $\lambda^{-1}$ , a complete procedure for the drift correction of a  $\lambda^{-1}$  model is presented by programming an augmented model. The robustness against signal uncertainties, coming from sensors or models, is ensured by introducing the so-called freezing rules that deactivate the updating when some conditions are fulfilled. A map for modelling the bias is proposed in order to solve the drift problem and to overcome the temporal and operating point variability of the bias.

The second part of the chapter emphasises the fast estimation of  $\text{NO}_x$  by following different methods.

First, the only use of look-up tables and a model sensor for estimating  $\text{NO}_x$  is presented. This section completes the experimental validation of the different methods for updating look-up tables, presented in Chapter 5. The results with all the methods are acceptable, but due to the robust properties of the SKF, this method is selected for the rest of the work.

Second, different observers are proposed in order to the online adaptation of the  $\text{NO}_x$  model presented in Chapter 4: the  $\lambda^{-1}$  observation, the online updating of the nominal  $\text{NO}_{x,0}$  table and the model bias correction. The results show that the model bias can be tracked by the online updating of the model parameters and inputs.

Finally, the estimation of the actual  $\text{NO}_x$  by avoiding the delay and sensor filtering is proposed by augmenting the state-space models in order to observe the actual signal. Concretely, two alternatives are proposed: a drift correction model and a look-up tables learning model. The results are presented and validated by utilising the second option.

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# Chapter 7

## Conclusions and future works

### Contents

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### 7.1 Main contributions and conclusions

The present dissertation covers the topic of the online dynamic estimation of  $\lambda^{-1}$  and  $\text{NO}_x$  in diesel engines. For that purpose, different sources of information are utilised:

- physical sensors for measuring exhaust gas concentrations (and intake oxygen when necessary), and
- control oriented models using ECU signals as fast inputs for estimating gas concentrations.

By combining the sensor and models, data fusion strategies are designed by means of observers based on the Kalman filter (KF) for the adaptive estimation of engine variables.

Specific tools, covering the online characterisation of sensors, the design of fast models and adaptive observers are developed in this work. Even though the methods of this work are applied to the estimation of  $\lambda^{-1}$  and  $\text{NO}_x$ , these can be applied to other engine variables, such as other gas concentrations or the volumetric efficiency, if a set-up with sensors and models is available. In the following, the main contributions and conclusions are presented and organised according to the thematic.

### 7.1.1 Online characterisation of gas concentration sensors

The measurement of engine-out gas concentration variables of a diesel engine is subjected to gas dilution, transport delays and is affected by the sensor hardware. In this work, a  $\text{NO}_x\text{-ZrO}_2$  sensor is used for measuring  $\lambda^{-1}$  and  $\text{NO}_x$  at the exhaust line of a diesel engine (while a gas analyser and a UEGO sensor are used for calibration and comparison when necessary).  $\text{NO}_x$  sensors are useful for control and diagnosis of AT systems, but the specific characteristics of the dynamic response must be assessed. The methods for the dynamic characterisation used in this work are based on the application of step-like transitions of the objective variables. For that purpose, ECU actuators are used for generating these profiles, allowing the on-board implementation of these methods.

**Static characterisation of  $\lambda^{-1}$  and  $\text{NO}_x$ .** Both  $\lambda^{-1}$  and  $\text{NO}_x$  outputs of the  $\text{ZrO}_2$  sensor are compared with those of a gas analyser, used as standard reference for the signals. The static accuracy, after measuring a representative set of the engine operating conditions, is enough for engine control purposes. A linear correlation is calculated with the static results.

**Dynamic characterisation of the  $\text{NO}_x$  output.** A simple on-engine experiment is proposed for the  $\text{NO}_x$  output characterisation. The method uses the engine as a gas generator and consists of applying steps in the start of injection (SOI) producing sharp variations ( $\Delta\text{NO}_x$ ) in the  $\text{NO}_x$  concentration. The use of SOI is justified by the high sensitivity of  $\text{NO}_x$  to SOI variations; the fact that SOI variations produce low variations in temperatures and pressures; and that SOI application is nearly instantaneous and no more than a

cycle-to-cycle response time is expected. Furthermore, the SOI command can easily be varied in the diesel engine.

The dynamic response of the sensor is identified with a delayed first order filter. The main parameters, namely response time and delay, are identified by utilising least squares (LS) tools. The response time of the sensor is highly constant whatever the conditions, even though some variations appear at low  $\text{NO}_x$  and/or low  $\Delta\text{NO}_x$  conditions. The identified delay has a bigger dispersion, especially when also approaching to idling. However the supposition of a constant delay for the designed algorithms gives satisfactory results.

**Dynamic characterisation of the  $\lambda^{-1}$  output.** A similar procedure to that of the  $\text{NO}_x$  output is made by performing injection step transitions provoking fast responses of  $\lambda$ , and thus  $\lambda^{-1}$ . LS methods are also used for identifying the delay and response time of a first order filter that models the sensor response.

The conclusions reached for the  $\text{NO}_x$  output are also applicable to the  $\lambda^{-1}$  output. The difference here is that the responses are a little bit faster, both in terms of delay and response time, because the output generated by an oxygen ion pump is located in a first cavity closer to the engine runner; while the  $\text{NO}_x$  output is proportional to a second oxygen ion pump located in a second chamber, thus increasing the delay and filtering.

In this work, a constant response time and delay have been utilised for modelling both  $\text{NO}_x$  and  $\lambda^{-1}$  outputs from the  $\text{NO}_x$  sensor. The results with the applied algorithms justify the selection of these simplistic sensor models.

### 7.1.2 Control oriented models for $\lambda^{-1}$ and $\text{NO}_x$

The present dissertation designs virtual sensors for the on-board prediction of  $\lambda^{-1}$  and  $\text{NO}_x$  by means of fast models which take profit of signals coming from the ECU.

**Fast modelling of  $\lambda^{-1}$ .** The relative fuel-to-air ratio ( $\lambda^{-1}$ ) is estimated on the basis of the ECU signals for the injected fuel mass flow ( $\dot{m}_f$ ) and the air mass flow ( $\dot{m}_a$ ). Both signals are fast, which in principle guarantees the dynamic response of the model. For simplicity, this model neglects the mass accumulation effects that affect the intake and exhaust manifolds. Such effects could be considered for correcting the value  $\dot{m}_a$  during the engine transients, e.g. when the boost pressure ( $p_{boost}$ ) is increased.

The model presents significant errors in the steady-state determination of  $\lambda^{-1}$ . This is due to the deviations that exhibit the hot wire sensor used in the series engine for measuring  $\dot{m}_a$ , and because the fuel mass is not measured, but determined through the rail pressure and the injection duration.

The bias on  $\dot{m}_a$  is expected to vary with the speed (because of the fluctuation patterns in the manifolds) and load (different set-points for  $p_{boost}$  and  $\dot{m}_a$ ). Possible leaks in the manifold and blow-by also depend on the engine load, thus affecting the total signal error. On the other hand, the error in  $\dot{m}_f$  estimation is affected by the injection settings (and hence load and engine speed).

When comparing the dynamic performance of the model with that of the sensor, the model exhibits almost instantaneous response:  $\dot{m}_f$  estimate, although biased, is instantaneous; and  $\dot{m}_a$  characteristic response time is in the order of milliseconds.

**Fast modelling of NO<sub>x</sub>.** An ECU oriented model for estimating NO<sub>x</sub> is designed based on an exponential variation around the nominal behaviour as a function of the available intake oxygen and a set of corrections for catching possible deviations from the nominal. First principle equations function of different ECU signals, such as  $n, m_f, \dot{m}_a, p_{boost}$  and the engine coolant temperature ( $T_{cool}$ ), are combined with different maps used for modelling the steady-state behaviour of the engine.

The available intake oxygen is approximated by means of the product of the EGR rate and  $\lambda^{-1}$ . A calculation based on the volumetric efficiency map, and the  $\dot{m}_a$  and  $p_{boost}$  signals, is utilised for estimating the EGR rate, after computing the intake temperature by a mixing model. The exhaust temperature ( $T_{exh}$ ) is modelled by tables and used for inferring the EGR manifold temperature ( $T_{egr}$ ). A set of corrections ( $C_{NO_x}$ ) for coping with different deviations in the intake mass flow ( $\dot{m}_{int}$ ), the intercooler temperature ( $T_{wic}$ ), ambient humidity ( $H$ ) and the intake temperature ( $T_{int}$ ) is proposed. The effect of the thermal transients is considered by means of a factor  $C_{dyn}$  that considers the temperature in the cylinder walls by means of a dynamic equation depending on the thermal loading, characterised by the product  $\lambda^{-1}n$ .

The model is tuned by means of steady-state tests for filling the maps and parameters, while it is also validated by means of dynamic engine cycles. The identified sensor model for the NO<sub>x</sub> output of the ZrO<sub>2</sub> sensor is used for comparing the model with the sensor output.



The results are satisfactory with both static and dynamic cycles, showing the capabilities of the model. The dynamic assumptions are sufficient for getting acceptable dynamic responses. The model tuning (mainly maps) is dependent on the calibration data set, and the ageing and changing conditions affect the model output, producing a biased response.

### 7.1.3 Observers for the fast estimation of engine variables

Two different observers are proposed for being used as adaptive estimators of engine variables:

- augmented models for drift correction, and
- on-board methods for updating look-up tables.

It is worth noting that these methods, presented in Chapter 5, are suited for a general problem, and can be applied for other engine variables if required inputs are available.

**Augmented models for drift correction.** The data fusion is done through a steady-state KF for observing the bias between a fast model affected by drift and a slow but accurate sensor. The filter works as a pass filter that keeps the high frequency components of the model and the low frequencies of the sensor. The model states are augmented for tracking the model bias. The steady-state analytical solution of the Kalman gain is solved and computed beforehand as function of the response time of the sensor and a noise quotient between the process and sensor noises. Guidelines for the observer tuning are proposed by a sensitivity analysis of the estimations covering possible errors or variations in the sensor modelling.

**Methods for updating look-up tables.** The problem of updating look-up tables by methods based on the extended Kalman filter (EKF) is treated. The table parameters are treated as states that can be observed if an output measurement is given. The table output is computed with a first order linear interpolation, which defines the output matrix of the learning state-space system. In a first approach, the problem has been focused on the use of a static table whose output is directly the aimed variable, i.e. the sensor dynamics are not considered in a first approach.

The EKF, with respect to other recursive methods for identification such as the recursive least squares (RLS), presents the advantage of tracking the

estimation error by means of a covariance matrix. Nevertheless, when coping with large systems, the computational burden is high, in terms of memory, fact that makes difficult its implementation for on-board strategies.

Beyond the EKF and based on it, two methods are developed:

- A steady-state KF (SSKF) calculated by means of building independent state-space models for every relative position of the matrix, which corresponds to a given operating point, e.g. scheduled by  $n$  and  $m_f$  in a general case. Therefore, if the inputs are considered stationary, the KF is steady-state under these hypotheses and a constant Kalman gain can be derived (the filter is linear time invariant). If covering the whole grid, the Kalman gain can be mapped for the operating range, depending on the inputs and noise trade-off.

The analytical solution of the steady-state filters is demonstrated and a proportional correction method is developed, resulting in a fast filter, whose gains can also be mapped beforehand if the noise variances are constants. Even though the method does not track the covariance error, the magnitude of the corrections could be modified online just varying the noise trade-off and recalculating the Kalman gain (by the analytical functions of by interpolating the map). The method behaves well for data when variations are rapid, closer to random data. However, it presents some convergence problems when the system knowledge is poor and the excitation level is not sufficient.

- A simplified KF (SKF) that lights the computation of the covariance matrix of the KF. This method exploits the fact that a N-D look-up table has a structure where at every iteration a maximum number of  $2^N$  elements are active, allowing to skip the updating of the elements out of this area. Nevertheless, a KF tracks the ageing of all elements, indeed the inactive ones, by means of storing large covariance matrices and compromising the memory and processing. The SKF method neglects the non-diagonal relationships of the covariance matrix (crossed covariances between table elements), but keeps the updating of the diagonals (variances).

The SKF method requires similar resources to those of the SSKF, but the performance is more similar to the one of the KF. The solution of the SKF is robust and the method behaves well under different data settings. The SKF is well-suited for a robust on-board implementation of look-up table estimators.

The static look-up tables can be linked with first order filters for representing dynamic equations, related with the sensor measurement or physical processes. For including the filters in the learning structure, an augmentation of the state-space model is proposed. After validating the methods by learning an adaptive table that models the NO<sub>x</sub> emissions, the SKF method is advised instead of the KF (heavier) or the SSKF (less robust), due to its performance is similar to that of the KF but with a comparable computational cost to that of the SSKF.

#### 7.1.4 Application of adaptive estimators to infer $\lambda^{-1}$ and NO<sub>x</sub> in diesel engines

Considering the methods developed in Chapter 5, the following three algorithms present useful data fusion strategies:

- algorithm A for tracking bias by an augmented model and a steady-state KF estimator,
- algorithm B for directly updating a static look-up table, and
- algorithm C for including the algorithm B structure inside a model for updating a table or parameter of a given model.

The results with the proposed structures are shown by applying to the engine-out  $\lambda^{-1}$  and NO<sub>x</sub> signals in a diesel engine. The uncertainties on the sensors and the models are considered as well.

##### 7.1.4.1 Fast observation of $\lambda^{-1}$

An observer is initially proposed for tracking the bias on the  $\lambda^{-1}$  signal computed on the basis of the fuel-to-air ratio. A set of rules is established in order to stop the bias integration when certain conditions do not ensure the problem robustness, which are:

- uncertainties on the sensor model when sharp transients,
- saturation problems of the sensor, and
- unexpected behaviours due to inaccuracies in the model or involved sensor signals, i.e.  $m_f$ ,  $\dot{m}_a$ ,  $\lambda^{-1}$ .

The estimation is accurate, even though two problems are underlined. First, the causality of the solution since the sensor response is delayed and therefore the actual bias is not computed on real time; second, the bias varies with the operating point conditions which involve high variations of the state.

These problems are solved by introducing an adaptive look-up table for modelling the bias. The table elements and the bias can be included directly in the observer structure. However, the table is kept as an independent block of the drift correction, being able to program independent filters for the bias and the table. Both the SSKF and the SKF methods are suitable for updating the table, although the results are shown by applying the former. The table output gives the prediction for the bias, while a steady-state KF is applied to update this bias and at the same time for giving the reference for also updating the table, and so on.

#### 7.1.4.2 Fast observation of NO<sub>x</sub>

The three different algorithms (A, B and C), and some modifications based on them, are applied to the estimation of NO<sub>x</sub> emission in diesel engines.

**Adaptive look-up tables for modelling NO<sub>x</sub>.** NO<sub>x</sub> emissions are modelled first by introducing a static 2D look-up table, as function of  $n$  and  $m_f$  in order to catch the nominal NO<sub>x</sub>, and a first order filter for representing the sensor dynamics.

The three learning methods (EKF, SKF and SSKF) are implemented and the results compared in order to fill the table parameters without any previous knowledge. The results are satisfactory for the three methods, but due to the special properties of the SKF, this method is advisable.

These tables can model with acceptable accuracy NO<sub>x</sub> emissions when the cycles are slow enough, allowing to simplify the closed loop engine response to a function that depends uniquely on  $n$  and  $m_f$ . Under this condition, the learning on a given cycle, i.e. NEDC, can be used for representing NO<sub>x</sub> emissions in other similar cycles such as the CADC, provided that the table elements have already been updated.

**Online updating of the NO<sub>x</sub> model.** Different modifications are introduced to improve the online and offline capabilities of the NO<sub>x</sub> model: the observation of  $\lambda^{-1}$  when a sensor signal is available, the adaptation of the main table of the NO<sub>x</sub> model (NO<sub>x,0</sub>), and the possible state augmentation

by tracking the model bias (discussed in the next paragraph). These modifications are proposed for correcting the ageing while rejecting spurious signals.

The benefits of the  $\lambda^{-1}$  observation are clear from the  $\lambda^{-1}$  estimation, avoiding that incorrect model estimations affect the  $\text{NO}_x$  model output. A drift correction algorithm is proposed when the  $\text{NO}_x$  sensor (or a UEGO sensor) is available.

The use of a learning model for updating the  $\text{NO}_{x,0}$  table in the  $\text{NO}_x$  model allows to avoid the ageing, minimising the online estimation error and the offline prediction error (model is adaptive). The SKF method is used for the application. The updating method is useful for both fitting a table when no initial knowledge exists or updating the model for cancelling the drift.

**Online observation of the actual  $\text{NO}_x$ .** The actual  $\text{NO}_x$  is inferred on the basis of the adaptive filtering models. Two alternatives are designed in this work.

First, a drift correction model is proposed for tracking the bias, and thus correcting the filtered  $\text{NO}_x$  model output when a  $\text{NO}_x$  sensor is available. This model is augmented with a state for tracking the actual  $\text{NO}_x$  (before filtering and delaying), resulting from adding the bias to the  $\text{NO}_x$  model output. The algorithm and causality problems are similar to those of the  $\lambda^{-1}$  estimation.

Second, the learning model for look-up tables can also be augmented with the actual  $\text{NO}_x$  state. In this case, the structure avoids the causality problem when the RT estimation is required. The direct interpolation of the table is the actual  $\text{NO}_x$  output, when the only delays are associated with the ECU signals, mainly  $n$ ,  $m_f$ ,  $\dot{m}_a$  and  $p_{boost}$ .

A further augmentation of the model with a bias state makes possible to decouple the table updating with respect to the fast actual  $\text{NO}_x$  estimation, in such a way that the related noises for the bias and table elements are tuned separately:

- Table parameters may be updated slowly by selecting a low noise related to those parameters,
- while the bias can be integrated faster if the sensor model is reliable and hence selecting a bigger noise for this state.

Other different trade-off might be selected in order to speed up the table convergence, but this structure allows to decouple in some way the table updating and the bias tracking. Of course that the lower the model error (table output) the lower the bias.

## 7.2 Future works

Even though the results presented in this work prove the validity of the proposed methods, the on-board implementation of them still deserves a deeper study. Particularly, the works to be performed as a continuation of this dissertation can be divided into those related with possible improvements of the sensor models, the design of an adaptive MVEM of the engine air path, the design of more complex emission models (and their possible adaptation), the validation of the methods and the different applications of the algorithms.

**Sensor knowledge.** A deeper study on the  $\text{NO}_x$  sensor response is mandatory in order to identify a physical model that can predict both the delay and response time with higher accuracy. The sensor delay and response times are variable, especially at low loads and low  $\text{NO}_x$  emissions (partly due to the loss of accuracy of the sensor for low  $\text{NO}_x$  concentrations [1]). Different phenomena and effects deserve an attention: the protective layer distortion, the internal diffusion speed between chambers and the effect of flux velocities, pressures, temperatures and concentrations over the final sensor output. Even though the pressure effect over the  $\text{NO}_x$  output from the  $\text{NO}_x$  sensor has been identified and it is not high for the pressures that can appear at the diesel engine exhaust, a deeper study should be conducted in order to characterise an appropriate correction. Furthermore, the influences of the pressure and other operation variables over the sensor should also be analysed.

To sum up, the use of a first order filter model with constant response times and delay could be improved by considering possible variations in these parameters as well as other influences that can be modelled for instance with transfer functions, or alternatively by considering a more representative non-linear model. In order to fulfil this task, the use of fast response measurements is required for calibration and validation.

**Design of an adaptive MVEM for the air path.** The air path model used for computing the EGR mass flow entering the engine is based on the volumetric efficiency and the measurements of the air mass flow and the boost pressure. The influence of the storage dynamics at the intake and the exhaust junction are not considered nor the dynamic effects over the intake and exhaust temperatures. Furthermore, the exhaust temperature is also modelled by a map with a set of corrections. The model could also consider the turbocharger (TC), the air filter, the actuators or the intercooler as well as the after-treatment (AT) systems.

The design of a MVEM considering these effects is not a challenge (this topic is well covered in the bibliography and the state of the art) but the adaptation of a MVEM is an interesting problem. Different strategies, similar to those presented in this work, can be implemented in order to estimate model states and inputs (see e.g. [2]).

With regards to the observation of the model inputs and states, drift correction models can be set for tracking the bias between the model states and sensor measurements. Good examples are the drifts on  $\dot{m}_a$  (by considering a model for the TC) and the injector (by using  $\lambda^{-1}$  and the corrected  $\dot{m}_a$ ).

With respect to the observation of model parameters as states, the fusion strategies can be utilised for updating the model while also for fitting the ECU parameters. Hence the drift correction and the updating algorithms are useful tools. However, there exists a problem with respect to the observability of the states and the filter tuning, i.e. it is difficult to weight the errors between all the parameters and the robustness must be assessed. The problem could be split up by designing local observers for the different sub-models, e.g. observing compressor maps by comparing with the MAF sensor output. Anyway, the closed-loop interaction should be assessed and a global cost function optimised.

This problem deserves a specific study and the use of synthetic signals should be a first step for evaluating the possibilities of the observers for solving a global optimisation problem. When involving look-up tables (frequent on MVEMs and ECUs), methods like the SKF permit the online adaptation with an affordable computational effort.

If the updating problem is designed globally and considering the closed loop interaction on real time, the discussion made for the SKF in order to simplify the KF when the system is large and only a few parameters are observable is also valid. Here, the use of parameters for defining how observable elements are at every iteration (e.g. grammians) allows to simplify the observation problem.

**Emission models.** The designed  $\text{NO}_x$  and  $\lambda^{-1}$  models are based on static maps and equations with a set of corrections and ECU signals as inputs, plus filters and delays for modelling physical and sensor dynamics. If other sensors are available, such as a cylinder pressure sensor, the prediction of the exhaust conditions can be benefited from a more accurate combustion model. The model states could also be augmented by including the soot prediction and considering further states from a MVEM of the air path.

In addition, an interesting improvement is modelling the AT devices. For instance, the monitoring of soot upstream and downstream of the DPF and the tight control of the exhaust  $\lambda^{-1}$  is required for an optimal control of the DPF. For a correct estimation, the combination of sensors and models with similar methods to those developed in this dissertation is advisable. The same example can be applied to SCR.

The physical interactions should also be included in control-oriented models of the AT line [3]. This is of particular importance for tracking the  $\text{NO}_x$  emissions at different points, i.e. although the total  $\text{NO}_x$  could be highly constant upstream of the de $\text{NO}_x$  system (SCR or LNT), the concentration of NO and  $\text{NO}_2$  species vary due to the oxidation in the DOC. And of course, the total  $\text{NO}_x$  will be reduced after de $\text{NO}_x$  systems. The crossed effects over other concentrations (e.g. soot) and over the sensor signals, as well as other factors such as the ammonia cross sensitivity should be taken into account. For such cases, the data fusion of sensors and models by observers is also interesting, as shown in Hsieh and Wang [4].

**Validation of the models and adaptive methods.** The assessment of the estimations and sensors should be validated with fast response measurement systems that can offer a dynamic and accurate reference of the considered variables. They are necessary for improving the sensor model identification and the validation of the methods.

Fast gas analysers based on CLD for measuring  $\text{NO}_x$  present response times around 2 ms, those based on NDIR for measuring CO/ $\text{CO}_2$  (thus oxygen or  $\lambda^{-1}$ ) reach 8 ms, while for soot these can be lower than 100 ms. These devices are suitable in order to identify the sensor dynamics and validate the results.

**Applications.** Once these methods and models are validated, the RT estimation of engine states ( $\text{NO}_x$  and  $\lambda^{-1}$  in the present work) could be used for the following functions:

- Control of the air path, fuel path and AT devices with feedback on exhaust emissions and other engine states,
- OBD of the AT systems,
- serving as inputs for other models or functions, as for instance, soot models (which depends on the exhaust  $\lambda^{-1}$ ), and
- tuning of controller parameters by model based optimisation [5].



To sum up, if a proper sensor set is available, engine variables and parameters can be observed. In addition to  $\lambda^{-1}$  and  $\text{NO}_x$ , interesting examples could be the volumetric efficiency, the intake oxygen concentration, the mass flows, temperatures and pressures in the manifolds, and other relevant exhaust concentrations. Due to the EURO VI and future legislations, this topic is of high interest, and recent advancements in sensor technology, as the case of soot sensors, may be benefited by the presented methods. Anyway, the final quality on the estimation also falls on the accuracy of the models used for the data fusion and the knowledge of the sensor behaviour.

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