Identification of Nonlinear Block-Oriented Systems starting from Linear Approximations: A Survey

Maarten Schoukens, Koen Tiels

Vrije Universiteit Brussel, Dept. ELEC, Pleinlaan 2, B-1050 Brussels, Belgium

Abstract

Block-oriented models are popular in nonlinear modeling because of their advantages to be quite simple to understand and easy to use. Many different identification approaches were developed over the years to estimate the parameters of a wide range of block-oriented models. One class of these approaches uses linear approximations to initialize the identification algorithm. The best linear approximation framework and the $\epsilon$-approximation framework, or equivalent frameworks, allow the user to extract important information about the system, guide the user in selecting good candidate model structures and orders, and they prove to be a good starting point for nonlinear system identification algorithms. This paper gives an overview of the different block-oriented models that can be modeled using linear approximations, and of the identification algorithms that have been developed in the past. A non-exhaustive overview of the most important other block-oriented system identification approaches is also provided throughout this paper.

Key words: System Identification, Maximum Likelihood, Nonlinear Systems, Hammerstein, Wiener, Wiener-Hammerstein, Hammerstein-Wiener, Parallel Cascade, Feedback, Best Linear Approximation, Linear Fractional Representation

1 Introduction

Nonlinear models are often used these days to obtain a better insight in the behavior of the system under test, to compensate for a potential nonlinear behavior using predistortion techniques, or to improve plant control performance. Popular nonlinear model structures are, amongst others, nonlinear state-space models (Suykens et al., 1995; Paduart, 2008; Schön et al., 2011), NARX and NARMAX models (Leontaritis & Billings, 1985; Billings, 2013), and block-oriented models (Giri & Bai, 2010). Due to the separation of the nonlinear dynamic behavior into linear time invariant (LTI) dynamics and the static nonlinearities (SNL), block-oriented nonlinear models are quite simple to understand and easy to use.

This paper provides a survey of the identification of block-oriented models using input-output data only. The identification algorithms that are discussed in this paper are all initialized using linear approximation frameworks for nonlinear systems: the Best Linear Approximation (Enqvist, 2005; Enqvist & Ljung, 2005; Pintelon & Schoukens, 2012) and the $\epsilon$-approximation (Schoukens et al., 2015a) framework. Many other block-oriented identification approaches, different from the linear approximation approaches, are described in the literature (e.g. overparametrization, iterative, or kernel-based methods). A non-exhaustive overview of the most important approaches for each model structure is given at the beginning of each section discussing that particular model structure. The presented identification approaches and the considered block-oriented structures go beyond the methods and structures that are presented in previous surveys, such as (Billings, 1980; Hasiewicz et al., 2012) and (Giri & Bai, 2010).

Block-oriented identification approaches starting from linear approximations have the advantage that part of the model selection problem is done in an LTI-framework, which is better understood by many practitioners. These approaches prove to obtain good results in various benchmark problems (such as the Wiener-Hammerstein (Schoukens et al., 2009b) and Silverbox benchmarks (Wigren & Schoukens, 2013), as illustrated in Section 11 of this paper. Linear models of nonlinear systems can also be used to detect the underlying block-oriented structure of the system under test. This structure detection problem using linear approximations is discussed in (Haber & Unbehauen, 1990; Lauwers et al., 2006; Schoukens et al., 2015a), however, no identification approaches are presented there.

* The corresponding author is M. Schoukens (maarten.schoukens@vub.ac.be).
Figure 1. The block-oriented structures that are considered in this paper. The different structures have been obtained by using series, parallel and feedback connections of LTI blocks ($G(q)$ and $S(q)$) and static nonlinear blocks ($f(\cdot)$ and $g(\cdot)$). There are three types of structure classes: single branch structures (Wiener, Hammerstein, Wiener-Hammerstein and Hammerstein-Wiener), parallel branch structures (parallel Wiener, parallel Hammerstein, parallel Wiener-Hammerstein) and feedback structures (simple feedback structure, Wiener-Hammerstein feedback and LFR).
This paper introduces first the different block-oriented structures that are considered throughout this paper in Section 2. Bussgang’s theorem is very useful to determine the BLA of many of the considered block-oriented structures. This theorem is discussed in Section 3. The considered linear approximation frameworks are discussed next in Section 4. Section 5 discusses the maximum likelihood nature of the considered cost function. The identification algorithms for the single branch models (Hammerstein, Wiener, Hammerstein-Wiener and Wiener-Hammerstein) are discussed in Section 6. The parallel block-oriented modeling approaches are discussed in Section 7. Section 8 considers the identification of different nonlinear feedback models. Section 9 provides an overview of the pros and the cons of the model structures and their identification based on the BLA, as explained in Sections 6–8. Some guidelines for the user are presented in Section 10. Finally, the good performance of the identification methods that are discussed in this paper is illustrated by some benchmark results in Section 11.

2 Block-Oriented Models: Structures, Representation and Noise Framework

Model Structures

Block-oriented models are constructed starting from two basic building blocks: a linear time-invariant (LTI) block and a static nonlinear block. They can be combined in many different ways. Series, parallel and feedback connections are considered in this paper, resulting in a wide variety of block-oriented structures as is depicted in Figure 1. These block-oriented models are only a selection of the many different possibilities that one could think of. For instance the generalized Hammerstein-Wiener structure that is discussed in [Wills & Ninness, 2012] is not considered in this paper.

Model Representation

The LTI blocks and the static nonlinear blocks can be represented in many different ways. The LTI block could for instance be modeled as a nonparametric FRF [Giri et al., 2013] or an impulse response [Lacy & Bernstein, 2003], or they can be parametrized using a state-space [Verhaegen & Westwick, 1996], rational transfer function [Narendra & Gallman, 1966], or basis function expansion [Tiels & Schoukens, 2014]. The static nonlinear block can again be represented in a nonparametric way using, for instance, kernel-based methods [Mzyk, 2014], or in a parametric way using, for instance, a linear-in-the-parameters basis function expansion (polynomial, piecewise linear, radial basis function network, ..., [Schoukens et al., 2014b]), neural networks [Schoukens et al., 2015b], or other dedicated parametrizations for static nonlinear functions.

The methods that are presented in this paper will typically use, but are not limited to, a rational transfer function representation for the LTI blocks and a linear-in-the-parameters basis function expansion for the static nonlinearity.

Another issue of block-oriented models is the uniqueness of the model parametrization. Gain exchanges, delay exchanges and equivalence transformations are present in many block-oriented structures [Schoukens et al., 2015b, Schoukens, 2015]. This results in many different models with the same input-output behavior, but with a different parametrization. The Jacobian of the cost function can be rank deficient due to these indistinguishability issues that are common for block-oriented identification problems. The related degenerations in the Jacobian need to be taken into account during the optimization.

The indistinguishability issues do not only impact the parametrization of the chosen model structure, it is also present between model structures. The simple feedback structure (see Figure 1) can both be represented using an LTI block in the forward path and a static nonlinearity in the feedback path or the other way around, a static nonlinearity in the forward path and an LTI block in the feedback path as is depicted in Figure 2. Both representations lead to the same input-output representation, but are structurally different [Schoukens et al., 2005b]. Indeed we have that the outputs of the two possible simple feedback structures are given by:

\[
\begin{align*}
\{y(t) &= G(q)[u(t) - g(y(t))] , \\
\tilde{y}(t) &= \tilde{g}(u(t) - \tilde{G}(q)\tilde{g}(t)) \}.
\end{align*}
\]

Which can be rewritten as:

\[
\begin{align*}
\{u(t) &= G^{-1}(q)y(t) + g(y(t)) , \\
u(t) &= \tilde{g}^{-1}(\tilde{g}(t)) + \tilde{G}(q)\tilde{g}(t) \}.
\end{align*}
\]

Both representations should have the same input-output relation, which results in the following system of equations:

\[
\begin{align*}
\{y(t) &= \tilde{y}(t) , \\
G^{-1}(q)y(t) + g(y(t)) &= \tilde{g}^{-1}(y(t)) + \tilde{G}(q)y(t) \}.
\end{align*}
\]

It can be easily observed that this system of equations holds for \(\tilde{G}(q) = G^{-1}(q)\) and \(\tilde{g}(.) = g^{-1}(.)\).

Noise Framework

Assumption 1. Noise framework: A Gaussian additive, colored zero-mean noise source \(n_p(t)\) with a finite variance \(\sigma^2\) is present at the output of the system only:

\[
y(t) = y_0(t) + n_p(t).
\]
Simple Feedback (a) 

\[ G(q) \]

\[ \gamma(t) \]

Simple Feedback (b) 

\[ \hat{G}(q) \]

Figure 2. Two simple feedback block-oriented model structures: (a) LTI block (\( G(q) \)) in the forward path and a static nonlinearity (\( \gamma(.) \)) in the feedback path, (b) a static nonlinearity (\( \hat{\gamma}(.) \)) in the forward path and an LTI block (\( \hat{G}(q) \)) in the feedback path. Both structures share the same input-output behavior when \( \hat{\gamma}(.) = g^{-1}(.) \) and \( G(q) = G^{-1}(q) \).

This noise \( n_y(t) \) is assumed to be independent of the known input \( u(t) \). The signal \( y(t) \) is the actual output signal and a subscript 0 denotes the exact (unknown) value.

The choice for the noise framework that is described in Assumption 1 is a simplified representation of reality. This simplification can lead to biased estimates when there are other noise sources present, located at other positions inside the system, e.g. process noise passing through a nonlinear subsystem [Hagenblad et al., 2008].

A more realistic noise framework can be obtained by introducing multiple noise sources, or by placing the noise source at a different location in the considered system structure. Multiple noise sources are, for instance, considered in [Hagenblad et al., 2008; Wills et al., 2013; Lindsten et al., 2013; Wahlberg et al., 2014]. This more realistic noise framework comes often at the cost of a more complex identification algorithm.

3 Bussgang’s Theorem and Separable Processes

3.1 Bussgang’s Theorem

Bussgang’s theorem [Bussgang, 1952] is stated in [Papoulis, 1991] as follows:

**Theorem 1.** If the input to a memoryless system \( y = g(u) \) is a zero-mean stationary Gaussian input \( u(t) \), the cross-correlation of \( u(t) \) with the resulting output \( y(t) = g(u(t)) \) is proportional to \( R_{u y}(\tau) \):

\[
R_{u y}(\tau) = \alpha R_{u u}(\tau) \quad \text{where} \quad \alpha = E\{g'(u(t))\} \quad (5)
\]

**Proof.** See [Bussgang, 1952; Papoulis, 1991].

In the frequency domain this becomes:

\[
S_{YU}(e^{j\omega T_r}) = \alpha S_{Uu}(e^{j\omega T_r}). \quad (6)
\]

Bussgang’s theorem proves to be very valuable for the analysis of block-oriented structures using linear approximations in the remainder of this paper.

3.2 Riemann Equivalence Class of Asymptotically Normally Distributed Excitation Signals

Here we consider the Riemann equivalence class of asymptotically normally distributed excitation signals [Schoukens et al., 2009a; Pintelon & Schoukens, 2012]. This signal class \( U \) mostly contains Gaussian noise sequences, but considering the Riemann class allows one to extend \( S_U \) to also contain periodic signal sets, while Bussgang’s Theorem still applies [Pintelon & Schoukens, 2012].

**Definition 1.** Riemann equivalence class of asymptotically normally distributed excitation signals \( S_U \). Consider a signal \( u(t) \) with a piecewise continuous power spectrum \( S_{UU}(e^{j\omega T_r}) \), with a finite number of discontinuities. A random signal \( u(t) \) belongs to the Riemann equivalence class if it obeys any of the following statements:

1. \( u(t) \) is a Gaussian noise signal with power spectrum \( S_{UU}(e^{j\omega T_r}) \).
2. \( u(t) \) is a random multisine or random phase multisine [Pintelon & Schoukens, 2012] such that:

\[
1 \over N \sum_{k=k_1}^{k_2} E\{|U(k)|^2\} = {1 \over 2\pi} \int_{\omega_{k_1}}^{\omega_{k_2}} S_{UU} \left( e^{j\nu T_r} \right) d\nu + O \left( N^{-1} \right),
\]

with \( \omega_k = k 2\pi f_s / N \), \( k \in \mathbb{N} \), \( 0 < \omega_{k_1} < \omega_{k_2} < \pi f_s \), \( f_s \) is the sampling frequency, and \( U(k) \) is the discrete Fourier spectrum of \( u(t) \).

A random phase multisine \( u(t) \) is a periodic signal with period length \( N f_s \) defined in [Pintelon & Schoukens, 2012] as:

\[
u(t) = N^{-1/2} \sum_{k=-N/2+1}^{N/2-1} U_k e^{j(2\pi k + \phi_k) f_s t}, \quad k \neq 0
\]

\[
= N^{-1/2} \sum_{k=1}^{N/2-1} 2U_k \cos(2\pi k t / N f_s + \phi_k), \quad (7)
\]

where \( U_k = U_k \) and \( \phi_k = -\phi_{-k} \) to obtain a real signal \( u(t) \). The phases \( \phi_k \) are random variables that are independent over the frequency and are a realization of a random process on \([0,2\pi] \), such that \( E\{e^{j\phi_k}\} = 0 \). For instance, the random phases can be uniformly distributed over the interval \([0,2\pi] \). The (real) amplitude \( U_k \) is set in a deterministic way by the user and is uniformly bounded by \( M_U / 0 < U_k < M_U < \infty \). Random phase multisines have the advantage of being periodic signals. This avoids the adverse effects of spectral leakage for a proper choice of the period length \( N f_s \). They also offer full control over the applied amplitude spectrum to the user. The Riemann equivalence ensures that
a random phase multisine is asymptotically \((N \to \infty)\) Gaussian distributed \cite{Pintelon2012}.

### 3.3 Separable Processes

Bussgang’s theorem has been extended to other classes of signals, besides the Gaussian class, using the concept of separable processes that is introduced in \cite{Nuttall1958}. Gaussian processes, sine wave and phase modulated signals are shown to be separable processes in \cite{Nuttall1958}. Furthermore, \cite{McGraw1968} have shown that the signals belonging to the class of elliptically symmetric distributions are also separable. A more in depth discussion of separable processes for nonlinear system identification can be found in \cite{Enqvist2005, Enqvist2010}.

### 4 Linear Approximations of Nonlinear Systems

Most real-world systems do not behave completely linearly. Nevertheless, a linear model often explains a significant part of the behavior of a (weakly) nonlinear system. This approximative linear model also provides the user with a better insight into the behavior of the system under test. It motivates the use of a framework that approximates the behavior of a nonlinear system by a linear time invariant model under well-defined system-specific boundary conditions. This paper uses the BLA framework \cite{Ljung2001, Enqvist2005, Enqvist2005, Pintelon2012, Schoukens2005a} and the \(\epsilon\)-approximation \cite{Schoukens2005a} to estimate a linear approximation of a nonlinear system given a fixed set of such boundary conditions. This section gives a brief introduction to the theoretical framework of the BLA and the \(\epsilon\)-approximation.

#### 4.1 Best Linear Approximation

The theoretical framework of the BLA results in a model whose structure is shown in Figure 3. It includes four components: one boundary condition, an input excitation class \(U\), and three model constituents, namely a Linear Time Invariant (LTI) model labeled \(G_{bla}(q)\), a perturbation noise source labeled \(n_y(t)\), and a nonlinear distortion source labeled \(\tilde{y}_s(t)\).

As mentioned above, the BLA of a system depends on the chosen signal class \(U\). Most identification methods that are discussed in this work consider \(U\) to be the Riemann equivalence class of asymptotically normally distributed excitation signals \(S_U\). When this class of signals is considered to provide an excitation signal, the BLA of many block-oriented systems becomes, due to Bussgang’s Theorem (see Section 3), a simple function of the linear dynamics that are present in that system. All signals are assumed to be stationary for the remainder of this work.

Figure 3. The BLA of a nonlinear system for a given class of excitation signals \(U\) consists of the resulting LTI model \(G_{bla}(q)\), the unmodeled nonlinear contributions \(\tilde{y}_s(t)\), and the additive noise source \(n_y(t)\) that is assumed to be present at the output of the nonlinear system. The zero-mean input excitation \(\tilde{u}(t)\) belongs to the signal class \(U\).

The BLA model of a nonlinear system is an LTI approximation of the behavior of that system. It is best in mean square sense for a fixed class of input signals \(U\) only. The BLA is defined in \cite{Enqvist2005, Enqvist2005, Pintelon2012} as:

\[
G_{bla}(q) = \arg \min_{G(q)} E \left\{ \| \tilde{y}(t) - G(q)\tilde{u}(t) \|^2 \right\},
\]  

(8)

where \(E\{\cdot\}\) denotes the expected value operator. The expected value \(E\{\cdot\}\) is taken w.r.t. the random input \(\tilde{u}(t)\). The zero-mean signals \(\tilde{u}(t)\) and \(\tilde{y}(t)\) are defined as:

\[
\tilde{u}(t) = u(t) - E\{u(t)\},
\]

(9)

\[
\tilde{y}(t) = y(t) - E\{y(t)\}.
\]

(10)

This definition of the BLA is equal to the definition of the linear time invariant second order equivalent model defined in \cite{Ljung2001, Enqvist2005, Enqvist2005, Pintelon2012} when the stability and causality restrictions imposed there are omitted.

It is shown in \cite{Enqvist2005, Enqvist2005, Pintelon2012} that eq. (8) is minimized by:

\[
G_{bla}(e^{j\omega T_s}) = \frac{S_{\tilde{y}G}(e^{j\omega T_s})}{S_{U\tilde{u}}(e^{j\omega T_s})},
\]  

(11)

where \(S_{\tilde{y}G}\) and \(S_{U\tilde{u}}\) denote the crosspower spectrum of \(\tilde{y}(t)\) and \(\tilde{u}(t)\) and autopower spectrum of \(\tilde{u}(t)\) respectively.

The output residuals \(n_i(t)\) of the model are given by:

\[
n_i(t) = \tilde{y}(t) - G_{bla}(q)\tilde{u}(t)
\]

(12)

These residuals represent the total distortion that is present at the output of the system. The total distortion can be split in two contributions based on their nature as is depicted in Figure 3. The nonlinear distortion \(\tilde{y}_s(t)\) represents the unmodeled nonlinear contributions, while the noise distortion \(n_y(t)\) is the additive noise that is assumed to be present at the system output. The nonlinear distortion and the noise distortion can be calculated...
separately as:

\[
\tilde{y}_n(t) = y_0(t) - G_{bla}(y) \tilde{u}(t), \\
\tilde{n}_n(t) = y(t) - \tilde{y}_n(t),
\]

where \(\tilde{y}_n(t)\) is the unknown noiseless output. The nonlinear distortion \(\tilde{y}_n(t)\) is uncorrelated with the input \(\tilde{u}(t)\), however, \(\tilde{y}_n(t)\) is not independent of the input \(\tilde{u}(t)\) (Pintelon & Schoukens, 2012). The noise source \(\tilde{n}_n(t)\), on the contrary, is typically assumed to be independent of the input \(u(t)\).

The obtained model \(G_{bla}(e^{j \omega T_x})\) does not only depend on the nonlinear system, but also on the class of input signals \(U\) that is used. The class \(U\) fixes both the probability density function (e.g. Gaussian inputs, uniform input distribution, etc.) and the power spectrum (or power spectral density) of the signals that are used to estimate it. Note that the BLA is equal to zero for zero-mean Gaussian signals applied to an even static nonlinear function (Bussgang, 1952).

The BLA can be estimated nonparametrically using the so-called robust and fast method (Schoukens et al., 2004, 2005a; Pintelon & Schoukens, 2012; Schoukens et al., 2012a). Afterwards, a parametric model can be estimated on top of the nonparametric one, where the nonparametric disturbance model is used as a frequency weighting (Pintelon & Schoukens, 2012). One can also choose to estimate a parametric model directly (Ljung, 1999; Enqvist & Ljung, 2005; Pintelon & Schoukens, 2012). This requires the user to make a careful selection of the disturbance model that captures both the nonlinear and noise distortion contributions.

### 4.2 \(\epsilon\)-Approximation

The Riemann equivalence class of asymptotically normally distributed excitation signals \(S_U\) is very useful to characterize a nonlinear system. There are some drawbacks, however, with the \(S_U\) class when it comes to structure detection and system analysis of some classes of systems. It is more convenient to consider infinitely small signals. For such excitation signals, the linearization of a cascade equals the cascade of the linearization. It is clear that such a signal cannot be used in practice, but it will allow us to formalize some system properties, especially for the Hammerstein-Wiener and feedback structures considered in this paper.

The class of \(\epsilon\)-excitations is defined in (Schoukens et al., 2015a) as follows:

**Definition 2. Class of \(\epsilon\)-Excitations \(S_\epsilon\).** The signal \(u_\epsilon(t), t = 1, \ldots, N\), belongs to the class of \(\epsilon\)-excitations \(S_\epsilon\) if it belongs to the class of Riemann equivalent excitation signals \(S_U\) and

\[
\sigma_u^2 = E\{\tilde{u}_\epsilon^2\} = \epsilon^2,
\]

\(\epsilon\)-excitations are (asymptotically) normally distributed signals with a limited variance. In this paper results are given for \(\epsilon\) converging towards zero.

**Definition 3. \(\epsilon\)-Approximation:** Consider the best linear approximation \(G_{bla}\) obtained for a random excitation \(u = u_{DC} + u_\epsilon\), and \(u_\epsilon \in S_\epsilon\), where \(u_{DC}\) is a constant offset. Define:

\[
\lim_{\epsilon \to 0} G_{bla}(e^{j \omega T_x})|_{u_\epsilon \in S_\epsilon} = G_\epsilon(e^{j \omega T_x})
\]

The definition of \(G_\epsilon\) can be extended to deterministic signals as discussed and formalized in (Mäkilä & Partington, 2003). The properties of the \(\epsilon\)-approximation are discussed in more detail in (Schoukens et al., 2015a).

### 5 Maximum Likelihood

A least squares cost function is minimized to obtain the model parameters:

\[
\hat{\theta} = \arg \min_\theta \sum_{t=1}^{N} (y(t) - \hat{y}(t, \theta))^2,
\]

where \(\hat{y}(t, \theta)\) is the modeled output, depending on the parameters \(\theta\).

Note that the cost function that is used in eq. (18) results in a maximum likelihood estimate if white Gaussian additive noise satisfying Assumption 1 is present at the output. A weighted version (using either a nonparametric noise model or a monic parametric noise model) of the cost function needs to be used to obtain a (sample) maximum likelihood estimate in the case of additive colored Gaussian noise. The estimated sample variance of the noise can be used as a weighting function in a frequency domain cost function. A maximum likelihood estimator is asymptotically efficient (Cramér, 1946; Ljung, 1999; Pintelon & Schoukens, 2012). This means that the maximum likelihood estimator achieves the lowest asymptotic mean squared error possible with a consistent estimator.

Unfortunately, this cost function is, in most of the cases considered in this paper, non-convex with respect to the parameters \(\theta\). A Levenberg-Marquardt algorithm (Levenberg, 1944; Marquardt, 1963; Fletcher, 1991; Pintelon & Schoukens, 2012) is used to minimize the cost function in a numerically stable way. This algorithm converges to a local minimum of the cost function. Hence, good initial values of the parameters are very important to
ensure the good quality of the final estimates. Such initial estimates can be obtained using the algorithms that are described in the following sections. Many of these initialization algorithms result in a consistent estimate, but they are not efficient.

## 6 Single Branch Models

The BLA of a single branch block-oriented system (Hammerstein, Wiener, Wiener-Hammerstein) provides valuable information about the system. It is shown in the following sections that the BLA of these systems is equal to a scaled version of the linear dynamics that are present in the system. This allows the user to perform an important part of the model selection problem, i.e. the selection of the number of poles and zeros in the system dynamics, in the LTI framework. Although the case of the Hammerstein-Wiener system structure is a bit more involved, the BLA approach proves to provide a good initial estimate for Hammerstein-Wiener identification.

### 6.1 Hammerstein

The Hammerstein structure is, together with the Wiener structure, one of the most simple block-oriented model structures. A Hammerstein structure [Hammerstein, 1930] consists of a static nonlinear block followed by an LTI block (see Figure 1). The Hammerstein structure is used to model nonlinear systems for which the static nonlinearity is present at the input of the system, such as a nonlinear actuator followed by a linear process, or some chemical processes and physiological systems (Giri & Bai, 2010).

Many different Hammerstein identification algorithms exist in the literature and are listed in (Giri & Bai, 2010). The first Hammerstein identification algorithm was introduced in (Narendra & Gallman, 1966), and further improved in (Bai & Li, 2004) and (Bai & Li, 2010). Several other Hammerstein identification algorithms were developed in the last decades. A non-exhaustive list is given below where the methods are classified depending on their properties: kernel-based and mixed parametric-nonparametric identification algorithms (Hasiewicz et al., 2012; Mzyk, 2014; Risuleo et al., 2015), parametric approaches (Chang & Luus, 1971; Crama et al., 2004; Schoukens et al., 2007; Falck et al., 2010; Risuleo et al., 2015), overparametrization approaches (Bai, 1998; Falck et al., 2010; Risuleo et al., 2015), blind identification algorithms (Bai, 2002; Vanbeylen et al., 2008), a set-membership approach (Szafraniec, 2009), and (continuous-time) instrumental variables approach (Laurain et al., 2012). The MIMO (multiple input multiple output) Hammerstein case is considered in (Verhaegen & Westwick, 1996; Goethals et al., 2005). The works of (Giri et al., 2011) and (Wang et al., 2012) extend the Hammerstein structure to contain dynamic backlash or hysteresis nonlinearities.

### Best Linear Approximation

The BLA of a Hammerstein system is a scaled version of the LTI dynamics that are present in \( S(q) \) if input signals belonging to the Riemann equivalence class of asymptotically normally distributed excitation signals are used.

**Theorem 2.** The BLA \( G_{bla}(e^{j\omega T_x}) \) of a Hammerstein system, excited by an input signal belonging to the Riemann equivalence class of asymptotically normally distributed excitation signals (Definition 1), is asymptotically given by:

\[
G_{bla}(e^{j\omega T_x}) = \alpha S(e^{j\omega T_x}), \quad \alpha \in \mathbb{R}.
\]  

**Proof.** The BLA is given by eq. (11):

\[
G_{bla}(e^{j\omega T_x}) = \frac{S_{\delta U}(e^{j\omega T_x})}{S_{\delta U}(e^{j\omega T_x})} = S(e^{j\omega T_x}) \frac{S_{\delta U}(e^{j\omega T_x})}{S_{\delta U}(e^{j\omega T_x})} = \alpha S(e^{j\omega T_x}),
\]

using the notation of Figure 1, where \( \frac{S_{\delta U}(e^{j\omega T_x})}{S_{\delta U}(e^{j\omega T_x})} = \alpha \) due to Bussgang’s theorem (Bussgang, 1952).

The gain \( \alpha \) depends not only on the system characteristics, but also on the amplitude, power spectrum, and the offset (DC value) of the input signal.

### Identification using the BLA

Many different approaches use the BLA, or similar correlation methods, to identify a Hammerstein system (Billings & Fakhouri, 1978; Hunter & Korenberg, 1986; Westwick & Kearney, 2003; Crama et al., 2004). The identification of a Hammerstein system using the BLA is quite straightforward. The BLA approach decouples the identification of the LTI block and the static nonlinear block.

Firstly, the LTI dynamics are estimated using the BLA, this results in a consistent estimate of the LTI block \( \hat{S}(q) \), up to an unknown gain exchange \( \alpha \) (see eq. (19)). Secondly, the static nonlinearity is estimated using a parametrization which is linear in the parameters. This results in a problem that is linear in the parameters:

\[
y(t) = S(q) \left[ f_i(u(t)) \right] = S(q) \sum_{i=1}^{n} \gamma_i f_i(u(t))
\]  

Finally, the estimated nonlinearity can be used to recover the actual input signal from the output signal. This is known as the inverse problem of Hammerstein identification.
where \( n_f \) denotes the number of basis functions \( f_i(.) \) that is used to describe the static nonlinearity. Note that eq. (25) is linear in the parameters \( \gamma_i \) when \( S(q) \) is replaced by its estimate \( \hat{S}(q) \). Hence, the parameters \( \gamma_i \) can easily be obtained by a linear least squares estimation. This results in a consistent, but not efficient, estimator of the model parameters (Schoukens, 2015). An efficient estimate can be obtained with the nonlinear optimization described in Section 5.

6.2 Wiener

A Wiener structure (Wiener, 1958) consists of an LTI block followed by a static nonlinear block (see Figure 1). A Wiener structure is often used to model systems where the nonlinear behavior is present at the output of the system. Some examples of such structures are systems with sensor nonlinearities, overflow valves and some physiological systems (Giri et al., 2010).

Several Wiener system identification algorithms are listed in (Giri & Bai, 2010). A wide variety of Wiener identification algorithms have been developed in the last decades. Some are grouped here based on their properties. A non-exhaustive list of methods contains: nonparametric (or semi-parametric) identification algorithms (Hasiewicz et al., 2012; Lindsten et al., 2013; Mzyk, 2014), parametric approaches (Billings & Fakhouri, 1977; Hunter & Korenberg, 1986; Wigren, 1993; Crama & Schoukens, 2001; Westwick & Kearney, 2003), a minimal Lipschitz approach (Pelckmans, 2011), (orthogonal) basis function expansion approaches (Lacy & Bernstein, 2003; Aljamaan et al., 2014), blind identification algorithms (Vanbeylen et al., 2009), a recursive approach (Greblicki, 2001), a set-membership approach (Sznaier, 2009), and for systems that contain backlash nonlinearities (Giri et al., 2014). The MIMO Wiener system case is considered in (Westwick & Verhaegen, 1996; Januczak, 2007). Most of the approaches consider that the noise source is present at the output of the system only, however some methods allow for process noise to be present in the middle of the Wiener system, in between the linear dynamics and the nonlinearity (Hagenblad et al., 2008; Lindsten et al., 2013; Wahlberg et al., 2014, 2015).

Best Linear Approximation

Similar to the Hammerstein case, the BLA of a Wiener system is a scaled version of the LTI dynamics that are present in \( G(q) \) if input signals belonging to the Riemann equivalence class of asymptotically normally distributed excitation signals are used.

**Theorem 3.** The BLA \( G_{bla}(e^{j\omega T}) \) of a Wiener system, excited by an input signal belonging to the Riemann equivalence class of asymptotically normally distributed excitation signals (Definition 1), is asymptotically given by:

\[
G_{bla}(e^{j\omega T}) = \alpha G(e^{j\omega T}), \quad \alpha \in \mathbb{R}. \tag{26}
\]

**Proof.** The proof is completely similar to the proof of Theorem 2, see also (Enqvist & Ljung, 2005; Pintelon & Schoukens, 2012; Schoukens, 2015).

Similar to the Hammerstein case, the gain \( \alpha \) depends not only on the system characteristics, but also on the amplitude, power spectrum, and the offset (DC value) of the input signal.

**Identification using the BLA**

The identification of Wiener systems using the BLA is again very similar to the Hammerstein case. BLA or correlation based identification methods for Wiener systems are presented in (Billings & Fakhouri, 1978; Hunter & Korenberg, 1986; Crama & Schoukens, 2001; Westwick & Kearney, 2003).

A two-step identification approach results in a consistent estimate of the model parameters. First, the LTI block is estimated using the BLA. Second, the static nonlinear block is identified starting from the estimated intermediate signal \( \hat{x}(t) = \hat{G}(q)[u(t)] \) and the system output. One can use his/her favorite model structure (polynomial, neural network, ...) to describe the static nonlinearity. This results in a consistent, but not efficient, estimator of the model parameters (Schoukens, 2015). An efficient estimate can be obtained with the nonlinear optimization described in Section 5.

6.3 Hammerstein-Wiener

Hammerstein and Wiener models were discussed in the previous sections. A more general block structure is the Hammerstein-Wiener model (static nonlinear block - LTI block - static nonlinear block, see Figure 1). These models are used in a wide range of applications such as chemical processes (Giri & Bai, 2010), ionospheric dynamics (Palanthandalam-Madapusi et al., 2005), submarine detection (Abrahamsson et al., 2007), and RF (radio frequency) power amplifier modeling (Taringou et al., 2010).

Different identification approaches are proposed to identify a Hammerstein-Wiener model or models that are very similar to Hammerstein-Wiener block structures. Amongst those are: iterative approaches (Zhu, 2002; Voros, 2004), overparametrization methods (Bai, 1998). 

\[
= \sum_{i=1}^{n_f} \gamma_i S(q)[f_i(u(t))], \tag{25}
\]
Section 5.

The BLA (or ϵ-approximation) is used as a starting point for the identification of a Hammerstein-Wiener system in (Crama & Schoukens, 2001). First, the BLA is estimated resulting in an estimate $\hat{G}(q)$ of the LTI block. Next, the front static nonlinearity and the inverse of the back static nonlinearity are identified by minimizing:

$$V_N = \sum_{t=1}^{N} \left( \sum_{i=1}^{n_f} \gamma_i \hat{G}(q) [f_i(u(t))] - \sum_{j=1}^{n_g} \delta_j g_j(y(t)) \right)^2$$

where it is assumed that the static nonlinearity $f(u)$ and the inverse of the static nonlinearity $g(r)$ can be represented by a linear combination of $n_f$ and $n_g$ nonlinear basis functions respectively. This problem is linear in the parameters, it can be solved using a total least squares approach (Van Huffel & Lemmerling, 2002). In a next step the estimate of the inverse of the back static nonlinear block can be replaced by an estimate of the forward static nonlinearity $g(r)$. These estimates serve as an initialization for the nonlinear optimization described in (Brouri et al., 2013).

Best Linear Approximation

Contrary to Hammerstein and Wiener systems, the BLA of a Hammerstein-Wiener system cannot be simplified to a simple function of the system dynamics, even if input signals belonging to the Riemann equivalence class of asymptotically normally distributed excitation signals are used. The intermediate signal $r(t)$ (see Figure 1) is not Gaussian due to the presence of the first static nonlinearity. Hence, Bussgang’s theorem cannot be applied. However, (Wong et al., 2012) shows that the BLA will not differ too much from a scaled version of the LTI block $G(q)$ in many practical cases (i.e. if the memory length of $G(q)$ is large w.r.t. the correlation length of the intermediate signal $x(t)$).

The theoretical framework of the ϵ-approximation (a small signal analysis of the system, see (Schoukens et al., 2015a)) can also be used to replace the BLA framework. The ϵ-approximation of a Hammerstein-Wiener system is a scaled version of the LTI dynamics $G(q)$ present in the system (Schoukens et al., 2015a).

Identification using the BLA

The BLA (or ϵ-approximation) is used as a starting point for the identification of a Hammerstein-Wiener system in (Crama & Schoukens, 2004). First, the BLA is estimated resulting in an estimate $\hat{G}(q)$ of the LTI block. Next, the front static nonlinearity and the inverse of the back static nonlinearity are identified by minimizing:

$$V_N = \sum_{t=1}^{N} \left( \sum_{i=1}^{n_f} \gamma_i \hat{G}(q) [f_i(u(t))] - \sum_{j=1}^{n_g} \delta_j g_j(y(t)) \right)^2$$

where it is assumed that the static nonlinearity $f(u)$ and the inverse of the static nonlinearity $g(r)$ can be represented by a linear combination of $n_f$ and $n_g$ nonlinear basis functions respectively. This problem is linear in the parameters, it can be solved using a total least squares approach (Van Huffel & Lemmerling, 2002). In a next step the estimate of the inverse of the back static nonlinear block can be replaced by an estimate of the forward static nonlinearity $g(r)$. These estimates serve as an initialization for the nonlinear optimization described in Section 5.

6.4 Wiener-Hammerstein

A Wiener-Hammerstein model has a structure that is a bit more involved than the Wiener or the Hammerstein model structure. The static nonlinearity is now sandwiched in between two LTI blocks (see Figure 1). The presence of the two LTI blocks results in a problem that is harder to identify: the main challenge in identifying a Wiener-Hammerstein system lies in the separation of the dynamics over the front and the back LTI block.

The approaches described in (Vandersteen et al., 1997; Tiels et al., 2015) estimate the two LTI blocks in a non-parametric way using carefully designed input signals. Volterra (and Tensor decomposition) based approaches are presented in (Weiss et al., 1998; Tan & Godfrey, 2002; Kibangou & Favrie, 2006; Ben Ahmed et al., 2013), an iterative approach is discussed by (Voros, 2007), a recursive EIV method is presented in (Mu & Chen, 2014) and evolutionary-based approaches are presented in (Dewhirst et al., 2010; Naitali & Girij, 2016). The MISO (multiple input single output) Wiener-Hammerstein case is discussed in (Boutayeb & Darouach, 1995), and the MIMO case in (Katayama & Ase, 2016). Many approaches use the BLA, or a similar correlation analysis, as a starting point for the algorithm e.g. (Billings & Fakhouri, 1978; Billings, 1986; Korenberg & Hunter, 1986; Westwick & Kenniey, 2003; Crama & Schoukens, 2005; Lauwers, 2011; Tan et al., 2012; Ströberg & Schoukens, 2012; Sjöberg et al., 2012; Westwick & Schoukens, 2012; Schoukens et al., 2014b; Vanbeylen, 2014; Tiels et al., 2014; Schoukens et al., 2014a; Tiels et al., 2015; Giordano & Sjöberg, 2015; Ase & Katayama, 2015; Katayama & Ase, 2016).

Best Linear Approximation

Similar to the Hammerstein and Wiener case, the BLA of a Wiener-Hammerstein system is a scaled version of the LTI dynamics that are present in $G(q) S(q)$ when input signals belonging to the Riemann equivalence class of asymptotically normally distributed excitation signals are used.

Theorem 4. The BLA $G_{bla}(e^{j\omega T_s})$ of a Wiener-Hammerstein system excited by an input signal belonging to the Riemann equivalence class of asymptotically normally distributed excitation signals (Definition 1) is asymptotically given by:

$$G_{bla}(e^{j\omega T_s}) = \alpha G(e^{j\omega T_s}) S(e^{j\omega T_s}), \quad \alpha \in \mathbb{R}. \quad (28)$$

Proof. The proof is completely similar to the proof of Theorem 2, see also (Schoukens, 2015).
Identification using the BLA

There are many BLA-based Wiener-Hammerstein identification methods. In this section we focus on the method that is presented in (Sjöberg & Schoukens, 2012; Sjöberg et al., 2012). The challenge of separating the front and the back dynamics is reformulated as a pole and zero allocation problem: which poles and zeros belong to the front dynamic block and which poles and zeros belong to the back dynamic block?

The BLA is estimated and parametrized in a first step, and the poles and zeros of the BLA are computed. These poles and zeros are, due to Theorem 4 and under the assumption that the product \( G(q)S(q) \) does not contain any pole-zero cancellations, the poles and zeros of the LTI blocks of the Wiener-Hammerstein system:

\[
G_{bla} = \alpha G(q)S(q),
\]

\[
\alpha \prod_{i=1}^{n_{p}} (p_i - q^{-1}) \prod_{i=1}^{n_{z}} (z_i - q^{-1}).
\]

The poles \( p_i \) and the zeros \( z_i \) of the parametrized BLA are the combined poles and zeros of the LTI blocks \( G(q) \) and \( S(q) \) of the Wiener-Hammerstein system.

These poles and zeros of the BLA need to be assigned to either \( G(q) \) or \( S(q) \):

\[
G_{bla} = \alpha \hat{G}(q)\hat{S}(q),
\]

\[
\hat{G}(q) = \prod_{i=1}^{n_{p}+n_{d}} (z_i - q^{-1})^{\beta_{z_i}}, \quad \prod_{i=1}^{n_{p}+n_{d}} (p_i - q^{-1})^{\beta_{p_i}},
\]

\[
\hat{S}(q) = \prod_{i=1}^{n_{p}+n_{d}} (z_i - q^{-1})^{-1-\beta_{z_i}}, \quad \prod_{i=1}^{n_{p}+n_{d}} (p_i - q^{-1})^{-1-\beta_{p_i}}.
\]

\( \beta_{p_i} \) and \( \beta_{z_i} \) are binary parameters that assign the pole or zero to either \( \hat{G}(q) \) (\( \beta_{p_i} = 1 \)), or to \( \hat{S}(q) \) (\( \beta_{p_i} = 0 \)). The parameters \( \beta_{p_i} \) and \( \beta_{z_i} \) are grouped in the binary parameter vector \( \beta \in \{0, 1\}^{(n_{p}+n_{d})} \times 1 \).

There is a finite number of different realizations of the parameter vectors \( \beta \). All possible realizations of the parameter vector are evaluated. To evaluate a realization of \( \beta \), a static nonlinearity is estimated for every realization \( k \). This problem is linear in the parameters if the static nonlinearity \( \hat{f}_k(x(t)) \) is described by a linear combination of basis functions. It is solved using linear least squares regression, resulting in the estimate \( \hat{f}_k \). Finally, all the estimated models are ranked based on their mean squares error \( mse_k \):

\[
mse_k = \frac{1}{N} \sum_{t=1}^{N} (y(t) - \hat{y}_k(t))^2,
\]

\[
\hat{y}_k(t) = \hat{S}_k(q) \left[ \hat{f}_k \left( \hat{G}_k(q) [u(t)] \right) \right],
\]

where \( N \) is the number of data points in the measured input-output record. The model with the lowest error is selected.

It is important to note that each complex conjugate pole or zero pair is assigned either to \( \hat{G}(q) \) or \( \hat{S}(q) \) in the combinations. The single elements of the pair are never assigned separately. This is required as the LTI blocks consist of transfer functions with real coefficients only. This also introduces a possible disadvantage of the approach: it can happen that two real poles (or zeros) are combined into a complex conjugate pole (or zero) pair during the parametrization of the BLA since the solution handles real poles and zeros as single elements. This pair cannot be correctly assigned if these poles (or zeros) originate from the two different LTI subsystems.

The total number of realizations of \( \beta \) depends on the number of poles and the number of zeros, \( n_{p} \) and \( n_{z} \), respectively. The total number is at least \( 2^{n_{p}+n_{z}} \) and at most \( 2^{2(n_{p}+n_{z})} \). The minimum number of realizations is obtained when all poles and zeros are part of a complex conjugate pair, the maximum number of realizations is obtained when all poles and zeros are real. This number increases very rapidly with the model order, which makes the brute-force scan approach in (Sjöberg & Schoukens, 2012) computationally expensive, since it equals the total number of least squares regressions that need to be performed to scan all possible pole-zero allocations.

A discrete optimization over the binary parameter vector \( \beta \) can be used to speed up the algorithm (Schoukens et al., 2014d). The evaluation of all possible pole and zero allocations can be avoided by using a fractional approach (Vanbeylen, 2014; Giordano & Sjöberg, 2015), or by using a higher order correlation, e.g. the so-called QBLA, to separate the dynamics over the front and the back LTI block (Billings & Fakhouri, 1978; Schoukens et al., 2008; Westwick & Schoukens, 2012; Schoukens et al., 2014b). An equivalent subspace version of the method presented in (Sjöberg & Schoukens, 2012) is developed in (Ase & Katayama, 2015), and extended to the MIMO case in (Katayama & Ase, 2016).

Although it is shown in (Sjöberg & Schoukens, 2012) that this approach results in a consistent estimate of the system parameters, a nonlinear optimization of all the parameters simultaneously as described in Section 5, following the linear initialization step, can be used to reduce the variance of the estimates.

7 Parallel Branch Model Structures

In the case of parallel block-oriented structures (parallel Hammerstein, parallel Wiener and parallel Wiener-Hammerstein, see Figure 1) the BLA will provide not only information about the LTI blocks that are present...
in the system, but also about the number of parallel branches in the system. By combining the BLAs obtained at different setpoints of the systems the selection of the number of parallel branches and the selection of the number of poles and zeros that are present in the system can be performed (Schoukens et al., 2011; Schoukens & Rolain, 2012c; Schoukens et al., 2015b). This allows the user to perform the selection of the number of poles and zeros in the system dynamics in the LTI framework, and to select the number of parallel branches in the system at an early stage of the identification procedure.

The BLA can also be used to construct a set of orthonormal basis functions to represent the system dynamics (Tiele & Schoukens, 2013). These basis functions are based on a number of user-selected pole locations. The poles of the BLA are a natural choice as a starting point to construct the orthonormal basis functions.

### 7.1 Parallel Hammerstein and Parallel Wiener

Parallel Hammerstein and parallel Wiener models are generalized forms of the Hammerstein and Wiener structure. They consist of a parallel connection of more than one Hammerstein (or Wiener) system (see Figure 1). Parallel Hammerstein models are sometimes called Uryson models, memory polynomial, or generalized Hammerstein models in the literature, see for example (Gallman, 1975; Billings, 1980; Doyle et al., 2002; Glarnouchi & Hammi, 2009).

**Parallel Hammerstein:** A popular approach that is used to identify a parallel Hammerstein system is to feed the input signal to a parallel connection of some nonlinear basis functions that are each followed by a finite impulse response filter (FIR) to realize a branch. The corresponding LTI multiple-input-single-output identification problem is then solved using one’s favorite FIR identification method in (Gadringer et al., 2007). In earlier work (Gallman, 1975), Hermite polynomials were used as orthogonal basis functions to model the static nonlinearity for Gaussian input signals. These approaches result in a linear least squares identification problem, which is easy to solve. However, they suffer from two drawbacks: i) the user gets no physical insight in the number of parallel paths in the device under test (DUT), as the number of parallel branches in the model is set by the degree of the nonlinearity, ii) for systems with long memory effects, a large number of parameters is needed due to the FIR-nature of the dynamic model. The model is therefore not parsimonious.

**Parallel Wiener:** Previously proposed parallel Wiener estimation methods suffer from some disadvantages. Some methods rely on an estimate of the Volterra kernel of the system under test (Kibangou & Favier, 2009). This requires a very large amount of data for the identification. Other methods are limited to the use of finite impulse response models for the linear subsystems (Lyzell & Enqvist, 2012a; Westwick & Kearney, 1997; Korenberg, 1991). These approaches typically result in an unwanted high number of parallel branches.

### Best Linear Approximation

Similar to the single branch case, the BLA of a parallel Hammerstein and parallel Wiener system is a simple function of the LTI dynamics that are present in the system.

**Theorem 5.** The BLA $G_{bla}(e^{j\omega T_s})$ of a parallel Hammerstein system excited by an input signal belonging to the Riemann equivalence class of asymptotically normally distributed excitation signals (Definition 1) is asymptotically given by:

$$G_{bla}(e^{j\omega T_s}) = \sum_{k=1}^{n_h} \alpha_k S[k](e^{j\omega T_s}), \quad \alpha_k \in \mathbb{R}. \quad (36)$$

**Proof.** The BLA of a Hammerstein system is given by eq. (19). This holds for each separate branch in the parallel Hammerstein system. The output $y$ is given by a summation of the outputs $y_k$ of each parallel Hammerstein branch. Hence, the BLA of a parallel Hammerstein system is given by the sum of the BLAs of each parallel branch. \qed

**Theorem 6.** The BLA $G_{bla}(e^{j\omega T_s})$ of a parallel Wiener system excited by an input signal belonging to the Riemann equivalence class of asymptotically normally distributed excitation signals and with a MISO static nonlinearity $g(.)$ of the parallel Wiener system that can be approximated arbitrarily well in least squares sense by a MISO polynomial is asymptotically given by:

$$G_{bla}(e^{j\omega T_s}) = \sum_{k=1}^{n_h} \alpha_k G[k](e^{j\omega T_s}), \quad \alpha_k \in \mathbb{R}. \quad (37)$$

**Proof.** See (Schoukens, 2015). \qed

An important observation with respect to eq. (36) and eq. (37) is that the input dependent gain $\alpha_k$ only appears in the numerator. For example, the BLA in eq. (36) is given by:

$$G_{bla}(q) = \sum_{k=1}^{n_h} \alpha_k S[k](q), \quad (38)$$

where:

$$S[k](q) = \frac{B_s[k](q)}{A_s[k](q)}. \quad (39)$$
Introducing a common denominator results in the following BLA expression:
\[
G_{bla}(q) = \frac{\sum_{k=1}^{n_{br}} \alpha_k B^{[k]}_i(q) \prod_{j=1, j \neq k}^{n_{br}} A^{[j]}_i(q)}{\prod_{k=1}^{n_{br}} A^{[k]}_i(q)}.
\] (40)

A similar result can be obtained for the parallel Wiener case.

This means that the poles of the identified BLA are also the poles of the LTI blocks that are present in the system. The zeros of the BLA of a parallel Hammerstein (or parallel Wiener) system may change when the amplitude, power spectrum, or the offset (DC value) of the input signal changes. Indeed, the denominator stays the same when the gains \(\alpha_k\) change, but the numerator coefficients depend on the gains \(\alpha_k\). This is discussed in further detail in [Schoukens et al., 2015a, Schoukens, 2015; Schoukens et al., 2015b].

### Parallel Wiener and parallel Hammerstein identification using the BLA - SVD Approach

The main difficulty in identifying a parallel Wiener or parallel Hammerstein system is separating the dynamics over the different parallel branches. This problem can be made easier by allowing more parallel branches to be present in the estimated model (see e.g. [Gallman, 1975; Billings, 1980; Doyle et al., 2002; Lyzell & Enqvist, 2012a; Westwick & Kearney, 1997; Korenberg, 1991]). A singular value decomposition (SVD) based approach to identify a parallel Wiener or parallel Hammerstein system with a low number of parallel branches is presented in [Schoukens et al., 2010, 2011; Schoukens & Rolain, 2012b; Schoukens et al., 2013a, 2015b]. This section explains the driving idea of this approach.

The approach proposed in [Schoukens et al., 2013a, Schoukens & Rolain, 2012b; Schoukens et al., 2015b] starts with an estimation of the BLA of the considered system for different operating conditions. The different operating conditions have been obtained using input signals with different power spectra. This includes the use of different magnitudes, different offsets, or different coloring of the power spectra.

Next, the measured BLAs are parametrized using a different LTI model for each operating condition. A common denominator model is used for all operating conditions simultaneously:
\[
\hat{G}_{bla}^{[i_r]}(q, \hat{\theta}_{bla}) = \frac{\hat{d}_0^{[i_r]} + \hat{d}_1^{[i_r]} q^{-1} + \ldots + \hat{d}_n^{[i_r]} q^{-n_d}}{c_0 + \hat{c}_1 q^{-1} + \ldots + \hat{c}_n q^{-n_c}},
\] (41)

where superscript \([i_r]\) denotes the setpoint of the system. This is indeed possible, as eq. (40) assures that the poles of the different measured BLAs are the same. A consistent estimate of the overall dynamics that are present in the nonlinear parallel Hammerstein or parallel Wiener system results. A matrix \(\hat{D}\) is constructed containing the stacked estimated numerator coefficients of the BLAs at the different operating conditions:
\[
\hat{D} = \begin{bmatrix}
\hat{d}_0^{[1]} & \hat{d}_1^{[1]} & \ldots & \hat{d}_n^{[1]} \\
\hat{d}_0^{[2]} & \hat{d}_1^{[2]} & \ldots & \hat{d}_n^{[2]} \\
\vdots & \vdots & \ddots & \vdots \\
\hat{d}_0^{[R]} & \hat{d}_1^{[R]} & \ldots & \hat{d}_n^{[R]}
\end{bmatrix}.
\] (42)

Starting from the parametrized BLAs, a decomposition of the overall dynamics at the different operating conditions is calculated. It uses the singular value decomposition (SVD) of the \(\hat{D}\) matrix:
\[
\hat{D} = \hat{U}_{bla} \hat{\Sigma}_{bla} \hat{V}_{bla}^T.
\] (43)

where the columns of \(\hat{V}_{bla}\) are the right singular vectors which act as an orthonormal basis for the row space of the \(\hat{D}\)-matrix, \(\hat{\Sigma}_{bla}\) is a diagonal matrix containing the singular values, and the columns of \(\hat{U}_{bla}\) constitute an orthonormal basis for the column space.

The column vectors in \(\hat{V}_{bla}\) provide an estimate of the numerator coefficients for each branch \(k\):
\[
\hat{S}^{[k]}(q) = \frac{\hat{\delta}_0^{[k]} + \hat{\delta}_1^{[k]} q^{-1} + \ldots + \hat{\delta}_n^{[k]} q^{-n_d}}{c_0 + \hat{c}_1 q^{-1} + \ldots + \hat{c}_n q^{-n_c}},
\] (44)

where \(\hat{\delta}_j^{[k]}\) is the element of the \(j\)-th row and \(k\)-th column of the matrix \(\hat{V}_{bla}\).

The number of parallel branches in the system is obtained based on the estimated rank of the decomposed matrix (based on the singular values obtained in \(\hat{\Sigma}_{bla}\)).

Once an estimate of the LTI blocks of the parallel Hammerstein or parallel Wiener system have been obtained, the estimation of the static nonlinearities in the system boils down to a simple linear least squares problem as in the single branch Hammerstein or Wiener case.

Although it is shown in [Schoukens et al., 2015b; Schoukens, 2015] that this approach results in a consistent estimate of the system parameters, a nonlinear optimization of all the parameters simultaneously as described in Section 5 can be used to reduce the variance of the estimates.

### 7.2 Wiener-Schetzen

A Wiener-Schetzen model (see Figure 4) is a subset of parallel Wiener models where the dynamics are
OBFs are stable, discrete-time, proper, rational transfer functions $F_i(e^{j\omega})$ that are orthogonal with respect to the inner product

$$\langle F_1, F_2 \rangle = \frac{1}{2\pi} \int_{-\pi}^{\pi} F_1(e^{j\omega}) F_2^*(e^{j\omega}) d\omega. \quad (45)$$

The outputs $x_1(t)$ and $x_2(t)$ of two OBFs $F_1$ and $F_2$ that share the same input are orthogonal if the input is white. This follows from the generalized expected value $E\{x_1(t)x_2(t)\}$, which can be written as

$$E\{x_1(t)x_2(t)\} = \frac{1}{2\pi} \int_{-\pi}^{\pi} F_1(e^{j\omega}) F_2^*(e^{j\omega}) |U(e^{j\omega})|^2 d\omega \quad (46)$$

using Parseval’s theorem \cite{Heuberger2005}, and which reduces to the inner product of two orthonormal transfer functions if the input is white ($|U(e^{j\omega})|^2$ is constant). Although in general no claims can be made about the orthogonality for non-white inputs, the considered OBFs are quite robust with respect to the coloring of the OBFs \cite{Heuberger2005, Chapter 6}.

Popular choices for the OBFs are Laguerre, Kautz, and Takenaka-Malmquist basis functions \cite{Heuberger2005}. The considered OBFs can be constructed from a set of poles $\{\xi_i\}$, and in its most general form are the Takenaka-Malmquist basis functions:

$$F_i(q) = \sqrt{1 - |\xi_i|^2} \prod_{k=1}^{i-1} \frac{q - \xi_k^* q}{q - \xi_k}, \quad (47)$$

for $i = 1, 2, \ldots; \xi_i \in \mathbb{C}$; and $|\xi_i| < 1$. Note that choosing all poles $\xi_i$ in the origin results in an FIR model. Choosing all poles identical and real results in a Laguerre basis. A two-parameter Kautz basis is obtained when choosing a pole structure $\{\xi_1, \xi_2, \xi_1, \xi_2, \ldots\}$. Note that this pole structure can include multiple repetitions of $\xi_1, \xi_2$, where $\xi_1$ and $\xi_2$ are either real or complex conjugated. Generalized orthonormal basis functions (GOBFs) are obtained when a finite set of $n_\xi$ poles is periodically repeated, i.e.

$$\xi_{i+(k-1)n_\xi} = \xi_i, \quad i = 1, 2, \ldots, n_\xi, \quad k = 1, 2, \ldots \quad (48)$$

A proper choice of the pole locations is important in order to accurately represent the linear dynamics with a limited number of OBFs. Choosing the poles equal to the true poles of the underlying linear dynamics of the (parallel) Wiener system is optimal in this sense. The poles of the BLA are thus excellent candidates to be used in constructing the OBFs. Possible mismatches between the true poles and the estimates from the BLA can be compensated for by periodically repeating the set of pole locations.

Best Linear Approximation

Since a Wiener-Schetzen model is a parallel Wiener model with a MISO polynomial nonlinearity, its BLA is given by Theorem 6. In particular, the BLA of a Wiener-Schetzen model is a linear combination of the dynamics in the different branches. The poles of the BLA are thus also the poles of the LTI dynamics that are present in the system.

Identification using the BLA

Wiener-Schetzen models appear in the literature under various names. In the original ideas of Wiener \cite{Wiener1958}, (continuous-time) Laguerre OBFs are used and the model is called a Laguerre system (see also \cite{Boyd1985}). Kibangou et al. \cite{Kibangou2005} uses discrete-time Laguerre OBFs and uses the name Laguerre-Volterra filters. When using generalized orthonormal basis functions, the names GOB-Volterra filters \cite{Kibangou2005} and Wiener/Volterra models \cite{daRosa2007} are used, amongst others.

Several methods have been proposed to identify Wiener-Schetzen models. \cite{Gomez2004} proposes a non-iterative method, where it is assumed that the nonlinearity is invertible and that this inverse can be described in terms of known basis functions. An optimal selection of one of the parameters in a truncated two-parameter Kautz OBF expansion is proposed in \cite{daRosa2007}. This optimal selection is done by minimizing an upper bound of the error made by approximating the system’s Volterra kernels by the truncated OBF expansion. The Volterra kernels of the system are assumed to be known in this approach. A model structure very similar to a Wiener-Schetzen model is identified in \cite{Totterman2009} and \cite{Gomez2012}, where the nonlinearity is expanded in terms of support vector machines (SVMs).

Figure 4. A Wiener-Schetzen model ($F_1(q), \ldots, F_n(q)$ are OBFs, $g(x_1, \ldots, x_n)$ is a multivariate polynomial).
estimates. Fast convergence rates can be obtained (Tiels & Schoukens 2013b, 2014).

Once the OBFs are constructed, the intermediate signals $x_1(t), \ldots, x_n(t)$ can be computed. Retrieving the polynomial coefficients then boils down to a linear regression. Note that by choosing OBFs and Hermite polynomials (Schetzen 2006), which are orthogonal for Gaussian inputs, this regression problem is optimally conditioned for white Gaussian inputs. Note that this is only true for infinitely long data records; for a finite data record, approximate orthogonality is obtained.

Compared to the BLA/SVD approach in the previous section, only one experiment suffices in this case. The model is also linear in the parameters, but this comes at the cost of a large number of parameters, since the number of polynomial coefficients increases combinatorially with the number of OBFs and the degree of nonlinearity. Dimension reduction techniques, such as presented in (Schoukens et al. 2013a, 2013b) can be used to reduce the number of basis functions, and hence, the number of parameters.

7.3 Parallel Wiener-Hammerstein

Parallel Hammerstein and parallel Wiener model structures are restricted to systems with dominant output dynamics (Hammerstein) or input dynamics (Wiener). A further increase in flexibility of the model structure is obtained here by considering the parallel Wiener-Hammerstein structure (see Figure 1). Previously published methods (Baumgartner & Rugh 1975, Wysocki & Rugh 1976, Billings & Fakhouri 1979, 1982) studied a subclass of the parallel Wiener-Hammerstein structure. Identification methods based on repeated sine measurements (Baumgartner & Rugh 1975, Wysocki & Rugh 1976), or white Gaussian inputs (Billings & Fakhouri 1979) are available for this model structure. It is shown in (Palm 1978, 1979) that a wide class of Volterra systems can be approximated arbitrary well using a parallel Wiener-Hammerstein model structure. However, no method is presented there to identify such models.

The parallel Wiener-Hammerstein system class that is used here is a more general system class than the $S_M$ system class that is used in (Baumgartner & Rugh 1975, Wysocki & Rugh 1976, Billings & Fakhouri 1979) in the following sense. The $S_M$ model has $M$ parallel branches, and the $m$-th branch contains a monomial nonlinearity that is fixed and equal to $(.)^m$. This restricts the model to have a polynomial nonlinearity only, and to contain only one branch for each degree of this polynomial nonlinearity. Thus a parallel Wiener-Hammerstein containing two parallel branches, each with different LTI subsystems, and with different polynomial nonlinearities can, in general, not be modeled by a $S_M$ model. The methods that are presented in (Schoukens et al. 2013b, 2014c, 2015b) also make some extra assumptions on the parallel Wiener-Hammerstein system. However, even when these assumptions are met, the considered system class still allows for a much more complex system behavior than the $S_M$ class does.

Best Linear Approximation

Similar to the parallel Hammeest and parallel Wiener case, the BLA of a parallel Wiener-Hammerstein system is a simple function of the LTI dynamics that are present in the parallel branches of the system.

Theorem 7. The BLA $G_{bla}(e^{j \omega T_s})$ of a parallel Wiener-Hammerstein system excited by an input signal belonging to the Riemann equivalence class of asymptotically normally distributed excitation signals (Definition 1) is asymptotically given by:

$$G_{bla}(e^{j \omega T_s}) = \sum_{k=1}^{n_{br}} \alpha_k G_k^{[k]}(e^{j \omega T_s}) S_k^{[k]}(e^{j \omega T_s}), \quad \alpha_k \in \mathbb{R}. \quad (49)$$

Proof. The BLA of a Wiener-Hammerstein system is given by eq. (28). This holds for each separate branch in the parallel Wiener-Hammerstein system. The output $y$ is given by a summation of the outputs $y_k$ of each parallel Wiener-Hammerstein branch. Hence, the BLA of a parallel Wiener-Hammerstein system is given by the sum of the BLA of each parallel branch.

Identification using the BLA - SVD Approach

The identification algorithm for the parallel Wiener-Hammerstein case is a combination of the identification algorithms presented in Sections 6.4 and 7.1. It is described in detail in (Schoukens et al. 2015b, 2015). Firstly, the BLA decomposition approach is applied as described in Section 7.1 for the parallel Hammerstein and parallel Wiener case to split the BLA dynamics over the different parallel branches.

Next, the poles and zeros of each branch are allocated to the front or the back simultaneously using the brute-force pole-zero allocation scan described in Section 6.4. The static nonlinearity of the model is estimated during this step.

Although it is shown in (Schoukens et al. 2015b, 2015) that this approach results in a consistent estimate of the system parameters, a nonlinear optimization of all the parameters simultaneously as described in Section 5 can be used to reduce the variance of the estimates.
Two disadvantages of this approach are 1) the high computational load due to the brute-force pole-zero allocation scan, and 2) the introduction of a MIMO static nonlinearity in the final model although the original system is decoupled of nature. The appearance of the MIMO nonlinearity is due to equivalence transforms and the SVD approach to split the dynamics (Schoukens et al., 2015b; Schoukens, 2015).

The brute-force pole-zero allocation scan can be made faster by allocating the poles and zeros of each branch separately instead of doing the allocation for all branches simultaneously (Schoukens, 2015).

The coupled, MIMO nonlinearity can be decoupled into a number of SISO (single input single output) static nonlinearities in a second step using tensor decomposition methods as is shown in Schoukens & Rolain, 2012a; Tiels & Schoukens, 2013a; Schoukens et al., 2014c; Dreesen et al., 2015).

8 Feedback Model Structures

The BLA or \(\epsilon\)-approximation can be used in many cases for the identification of feedback block-oriented model structures. Both frameworks allow the user to perform the selection of the number of poles and zeros in the system dynamics in the LTI framework, in the three cases considered in this section.

It is commonly assumed for nonlinear feedback systems to have at least one sample delay in either the forward or the backwards path of the feedback loop to avoid the presence of nonlinear algebraic loops during simulation. It is shown in Relan & Schoukens, 2016 under band-limited assumptions that the approximation error made by introducing such a delay in the model can be kept arbitrary low by selecting a sufficiently high sampling frequency.

8.1 Simple Feedback Structure

The simple feedback structure consists of the feedback connection of an LTI block and a static nonlinear block (see Figure 1). One can think of two possible simple feedback structures: LTI block in the feedforward and the nonlinearity in feedback or vice-versa. It can be shown that both combinations are equivalent from an input-output point of view (see Section 2 and Schoukens et al., 2005b). A BLA-based identification algorithm for the simple feedback structure is proposed in Paduart et al., 2004; Paduart, 2008.

Best Linear Approximation

Contrary to most of the previous cases, the BLA of a simple feedback system cannot be simplified to an easy function of the system dynamics, even if input signals belonging to the Riemann equivalence class of asymptotically normally distributed excitation signals are used. The input signal of the nonlinearity \(y(t)\) is not Gaussian due to the presence of the nonlinear feedback loop. Hence, Bussgang’s theorem cannot be applied. However, in many practical cases it is fair to assume that the nonlinearity can be linearized to a gain \(\alpha\) and thus that the BLA will be close to something of the form:

\[
G_{bla}(e^{j\omega T_s}) \approx \frac{G(e^{j\omega T_s})}{1 + \alpha G(e^{j\omega T_s})}, \quad \alpha \in \mathbb{R}. \quad (50)
\]

The \(\epsilon\)-approximation, contrary to the BLA, results in a simple and exact analytical expression of the linear approximation as a function of the system dynamics (Schoukens et al., 2015a):

\[
G_{\epsilon}(e^{j\omega T_s}) = \frac{G(e^{j\omega T_s})}{1 + \gamma G(e^{j\omega T_s})}, \quad \gamma \in \mathbb{R}. \quad (51)
\]

The gains \(\alpha\) and \(\gamma\) depend not only on the system characteristics, but they can also depend on the amplitude, power spectrum, and the offset (DC value) of the input signal.

Note that in the feedback case (also in the Wiener-Hammerstein feedback case in Section 8.2), only the poles of the \(\epsilon\)-approximation shift with a changing gain \(\gamma\):

\[
G_{\epsilon}(q) = \frac{B(q)}{A(q) + \gamma B(q)}. \quad (52)
\]

where

\[
G(q) = \frac{B(q)}{A(q)}. \quad (53)
\]

This is discussed in further detail in Schoukens et al., 2015a.

Identification using the BLA

The identification algorithm that is presented here follows the lines of the algorithm presented in Paduart et al., 2004; Paduart, 2008.

Firstly, it is important to note that an arbitrary gain \(\beta\) can be shifted from the static nonlinear block to the linear block in the model. Indeed we have that:

\[
y(t) = G(q)[u(t) - f(y(t))] = G(q)[u(t) - f(y(t)) + \beta y(t) - \beta y(t)]
\]
(1 − βG(q))[y(t)] = G(q)[u(t) − ̂f(y(t))]  
(54)

\[ y(t) = \frac{G(q)}{1-\beta G(q)} [u(t) - ̂f(y(t))] \]

\[ y(t) = ̂G(q)[u(t) - ̂f(y(t))] \]

where ̂f(y(t)) = f(y(t)) − βy(t) and ̂G(q) = \frac{G(q)}{1-\beta G(q)}.

This indicates that the BLA results in a good initial estimate of the linear subsystem up to the equivalence transforms present in the model structure:

\[ ̂G(q) = G_{bla}(q) ≈ \frac{G(q)}{1 + \alpha G(q)}. \]  
(55)

The static nonlinearity is estimated next. The static nonlinearity is assumed to be represented as a linear combination of nonlinear basis functions. It is estimated by cutting the feedback loop:

\[ ̂G(q)[u(t)] - y(t) = ̂G(q)\sum_{i=1}^{n_f} \gamma_i f_i(y(t)), \]

\[ = \sum_{i=1}^{n_f} \gamma_i ̂G(q)[f_i(y(t))], \]  
(56)

where \( n_f \) denotes the number of basis functions that is used to describe the static nonlinearity. Note that this equation is linear in the parameters \( \gamma_i \). Hence, the parameters \( \gamma_i \) can easily be obtained by a linear least squares estimation.

Note that the noisy output is used as a regressor during the initial identification of the static nonlinearity in eq. (56). This results in a very simple initialization scheme which works well for high signal-to-noise ratios.

Finally, a nonlinear optimization (see Section 5) of all the parameters of the simple feedback structure together is applied such that a consistent estimate is obtained. The construction of the Jacobian of the simple feedback structure is more involved and time-consuming (compared to the single and parallel branch case) due to its recursive dependency on the output at previous time instances. This holds as well for the Wiener-Hammerstein feedback case and the LFR case in Sections 8.2 and 8.3.

8.2 Wiener-Hammerstein Feedback

The Wiener-Hammerstein feedback structure is a generalized version of the simple feedback structure. It consists of a Wiener-Hammerstein system in the feedback path combined with an LTI system in the feedback path (see Figure 1). Both BLA-based approaches [Schoukens et al., 2005b, 2008] and a convex relaxation based approach [Sou et al., 2008] have been developed for this system structure. It can be shown that this structure is equivalent to an LTI system in the forward path and a Wiener-Hammerstein system in the feedback path [Schoukens et al., 2005b, 2008].

Best Linear Approximation

As for the simple feedback structure case, the BLA of a Wiener-Hammerstein feedback structure cannot be simplified to an easy function of the system dynamics if input signals belonging to the Riemann equivalence class of asymptotically normally distributed excitation signals are used. However, in many practical cases it is fair to assume that the BLA will be close to something of the form:

\[ G_{bla}(e^{jωT_s}) ≈ \frac{α G^{[1]}(e^{jωT_s}) G^{[2]}(e^{jωT_s}) G^{[3]}(e^{jωT_s})}{1 + α G^{[1]}(e^{jωT_s}) G^{[2]}(e^{jωT_s}) G^{[3]}(e^{jωT_s})}, \]  
(57)

The \( \epsilon \)-approximation, contrary to the BLA, results in a simple and exact analytical expression of the linear approximation in function of the system dynamics [Schoukens et al., 2015a]:

\[ G_ε(e^{jωT_s}) ≈ \frac{γ G^{[1]}(e^{jωT_s}) G^{[2]}(e^{jωT_s}) G^{[3]}(e^{jωT_s})}{1 + γ G^{[1]}(e^{jωT_s}) G^{[2]}(e^{jωT_s}) G^{[3]}(e^{jωT_s})}, \]  
(58)

The gains \( α \) and \( γ \) depend not only on the system characteristics, but can also depend on the amplitude, power spectrum, and the offset (DC value) of the input signal.

Identification using the BLA

The BLA-based identification algorithm is outlined in [Schoukens et al., 2005b, 2008]. The key is the observation that the inverse of the BLA is approximately given by:

\[ \frac{1}{G_{bla}(q)} ≈ G^{[3]}(q) + \frac{β}{G^{[1]}(q)G^{[2]}(q)}, \]  
(59)

where \( β = \frac{1}{α} \).

In a first step, the BLA of the system is identified at two or more different setpoints of the system:

\[ ̂G_{bla}^{[j]}(q) ≈ \frac{α^{[j]} G^{[1]}(q) G^{[2]}(q) G^{[3]}(q)}{1 + α^{[j]} G^{[1]}(q) G^{[2]}(q) G^{[3]}(q)}. \]  
(60)

A nonparametric initial estimate ̂G^{[3]}(e^{jωT_s}) of the linear feedback system is obtained as the mean of the inverses
theory to estimate the LFR model. The method presented by {Van Mulders et al., 2013} starts from a polynomial nonlinear state space model and transforms it into a LFR model. The block-oriented, BLA-based method that is presented in {Vanbeylen, 2013} is studied in some more detail below.

**Best Linear Approximation**

The BLA of a general LFR system cannot be simplified to an easy function of the system dynamics, even if input signals belonging to the Riemann equivalence class of asymptotically normally distributed excitation signals are used. The input signal of the nonlinearity \( y(t) \) is not Gaussian due to the presence of the nonlinear feedback loop. Hence, Bussgang’s theorem cannot be applied. However, in many practical cases it is fair to assume that the nonlinearity can be linearized to a gain \( \alpha \) and thus that the BLA will be close to something of the form:

\[
G_{bla}(e^{j\omega T_s}) \approx G[1](e^{j\omega T_s}) + \frac{\alpha G[2](e^{j\omega T_s})G[3](e^{j\omega T_s})}{1 + \alpha G[4](e^{j\omega T_s})}, \quad \alpha \in \mathbb{R}. 
\]

The \( \epsilon \)-approximation, contrary to the BLA, results in a simple and exact analytical expression of the linear approximation in function of the system dynamics \( \text{Schoukens et al., 2015a} \):

\[
G_\epsilon(e^{j\omega T_s}) = G[1](e^{j\omega T_s}) + \frac{\gamma G[2](e^{j\omega T_s})G[3](e^{j\omega T_s})}{1 + \gamma G[4](e^{j\omega T_s})}, \quad \gamma \in \mathbb{R}. 
\]

The gains \( \alpha \) and \( \gamma \) depend not only on the system characteristics, but they can also depend on the amplitude, power spectrum, and the offset (DC value) of the input signal.

Note that in the LFR case both the poles and zeros of the \( \epsilon \)-approximation shift with a changing gain \( \gamma \):

\[
G_\epsilon(q) = G[1](q) + \frac{\gamma G[2](q)G[3](q)}{1 + \gamma G[4](q)}, \quad \text{where} \quad G[1](q) = \frac{B[i](q)}{A[i](q)}. 
\]

This results in the following expression, containing both a dependency on \( \gamma \) in numerator and denominator:

\[
G_\epsilon(q) = \frac{B[1](q)A[2](q)A[3](q)A[4](q) + \gamma B[4](q)}{A[1](q)A[2](q)A[3](q)A[4](q) + \gamma B[4](q)}.
\]

8.3 **Linear Fractional Representation**

The linear fractional representation (LFR) model structure is studied in \{Vandersteen & Schoukens, 1999\} \{Van Mulders et al., 2013\} \{Vanbeylen, 2013\}. It is a combination of a Wiener-Hammerstein feedback system with a linear feedforward branch in parallel (see Figure 1). It is the most general block-oriented system that contains only one static nonlinear block. All other block-oriented systems with only one static nonlinearity can be reformulated to an LFR structured system.

The method in \{Vandersteen & Schoukens, 1999\} relies on a large set of two-tone measurements and Volterra
\[
\gamma A^{[1]}(q)B^{[2]}(q)B^{[3]}(q)A^{[4]}(q) + A^{[1]}(q)A^{[2]}(q)A^{[3]}(q)(A^{[4]}(q) + \gamma B^{[4]}(q)) \tag{69}
\]

This is discussed in further detail in (Schoukens et al., 2015a).

**Identification using the BLA - Ricatti Approach**

The identification approach for the LFR model that is presented in (Vanbeylen, 2013) shares some of its aspects with the approaches that are explained above: it relies on at least two measured BLAs, and it needs to perform a scan over several possible solutions to allocate the dynamics over the different LTI-blocks in the model. The main aspects of the identification algorithm are described below.

First, the BLA is estimated at two different setpoints of the system. This BLA is converted to linear state space models.

Next an equivalence transformation for each state space model is estimated by solving a nonsymmetric algebraic Riccati equation. The details of this equation can be found in (Vanbeylen, 2013). This results in different possible solutions.

Each solution is evaluated by constructing the corresponding LFR model. The model with the best fit is selected. This step is very similar to the pole-zero allocation scan during the Wiener-Hammerstein identification in Section 6.4.

A nonlinear optimization (see Section 5) of all the parameters of the closed loop simple feedback structure together is applied such that a consistent estimate is obtained.

### 9 Pros and cons of the model structures

The model structures and identification methods that are explained in the previous sections all have their pros and cons. Some model structures have the ability to accurately model the input/output behavior of a large class of systems, but use a large number of parameters in return. Some identification methods can easily generate starting values from a single experiment, others require computationally expensive steps on multiple data sets. This section provides an overview of the model structures in terms of their descriptive power, number of parameters, and general difficulty to estimate the model. These criteria are explained below.

In the context of this paper, a universal approximator is a model structure that can arbitrarily well approximate the input/output behavior of systems belonging to the class of Wiener systems in mean-square sense (Schoukens et al., 2015a). This means that there exists a uniformly bounded Volterra series whose output converges in mean-square sense to the system’s output (Schetzen, 2006). The class of Wiener systems should not be confused with the block-oriented Wiener system class of Section 6.2. Since systems belonging to the class of Wiener systems respond to a periodic input with a periodic steady-state output with the same period (Boyd et al., 1984), these systems are also called PISPO (period in same period out) systems. This class of systems includes systems with saturation and discontinuous nonlinearities, but it excludes chaotic behavior, sub-harmonics, and hysteresis. Parallel Wiener and parallel Wiener-Hammerstein models are universal approximators without any constraint on the excitation class (Schetzen, 2006). Other approximation properties of these model structures have been shown in the past (Palm, 1979; Boyd & Chua, 1985; Korenberg, 1991).

A large number of independent parameters leads to a large variance on the estimated parameters (Söderström & Stoica, 1989) when a model is identified from noisy measurements. Therefore, a model is preferably parameter-parsimonious, as decreasing the variability on the estimated parameters requires larger data sets, thereby increasing the computational time of the optimization step in Section 5.

The optimization step in Section 5 converges to a local optimum of the (non-convex) cost function. Good initial estimates are thus required. A consistent initial estimate is preferable, as in that case a good initial estimate is guaranteed as more estimation data are used. The optimization step itself is consistent for each of the considered model structures.

Some identification approaches require multiple data sets, e.g. BLA estimates at different setpoints, which is not always possible due to time constraints.

Table 1 provides an overview of these properties as well as some general difficulties of the identification algorithms that are presented in the previous sections.

Only the parallel Wiener, Wiener-Schetzen, and parallel Wiener-Hammerstein model structures are universal approximators. Nevertheless, these feedforward structures cannot capture chaotic behavior, sub-harmonics, or hysteresis, since this requires feedback. The Linear Fractional Representation (LFR) is the most general block-oriented structure that contains only one static nonlinear block. The Wiener, Hammerstein, Wiener-Hammerstein, simple feedback, and Wiener-Hammerstein feedback structures are thus all special cases of the LFR.

The single branch models of Section 6 all have a low number of parameters, since these structures consist of
a limited series connection of SISO blocks. The parallel Hammerstein model structure consists of a number of parallel connections of Hammerstein models and thus still has a fairly low number of parameters. The parallel Wiener and Wiener-Schetzen model structures both have a MISO static nonlinear block, which typically results in a high number of parameters. For the parallel Wiener structure, however, the proposed identification method results in a low number of parallel branches, i.e. a low number of inputs to the nonlinear block, and thus a medium number of parameters. The parallel Wiener-Hammerstein structure has a MIMO static nonlinear block, resulting in a high number of parameters. The feedback structures of Section 8 consist of a limited (at most five) SISO blocks, and thus have a low number of parameters. The LFR structure, however, is typically parameterized in state-space form \(\text{[Vanbeylen, 2013]}\), which introduces at least \(n^2\) redundant parameters, where \(n\) is the model order. This is due to the existence of an \(n\) by \(n\) transformation matrix that leaves the input/output behavior unchanged. This transformation matrix should be nonsingular, but otherwise has elements that can be freely chosen. The LFR structure can thus have a medium number of parameters, although the number of independent parameters is still low.

Most of the proposed identification methods result in consistent initial estimates. This is not the case for the Hammerstein-Wiener identification method and the methods for the feedback structures of Section 8, since these methods use the noise-corrupted output either in some of the regressors (Hammerstein-Wiener and simple feedback) or to reconstruct internal signals that constitute some of the regressors (Wiener-Hammerstein feedback and LFR). Moreover, an approximation to the BLA is made in the Hammerstein-Wiener structure and all feedback structures. This approximation can be avoided by replacing the BLA-framework by the theoretical framework of the \(\epsilon\)-approximation \(\text{[Schoukens et al., 2015a]}\).

The presented methods on model structures that include a Wiener-Hammerstein branch all involve a brute-force scan to separate the dynamics of the Wiener-Hammerstein branch over its two LTI blocks. A similar scan over the solutions of a nonsymmetric algebraic Riccati equation is performed in the LFR identification method. These scans are mildly computationally expensive if only one Wiener-Hammerstein branch is involved (Wiener-Hammerstein, Wiener-Hammerstein feedback, and LFR structures). The scan is computationally expensive for the parallel Wiener-Hammerstein structure, since there, a scan needs to be performed for each branch and each time a MIMO nonlinearity needs to be estimated.

10 Guidelines for the User

Two important steps in the system identification process are the input design and model structure selection step. This section provides some user guidelines for the input design and the model structure selection problem.

10.1 Model Structure Selection

The use of the BLA framework is not limited to obtaining a block-oriented model of a nonlinear system. The BLA framework also offers a powerful analysis tool that allows the user, when combined with a well-thought-out input design strategy, to acquire a lot of insight in the nonlinear system at very low cost. The questions "How nonlinear is the system?", "Do I need a nonlinear model?", and "What kind of nonlinear effects do I observe in the system?" can be (partly) answered within the BLA framework \(\text{[Lauwers et al., 2008]}\) \(\text{[Pintelon & Schoukens, 2012]}\) \(\text{[Schoukens et al., 2012a, 2015a]}\), an overview can be found in \(\text{[Schoukens et al., 2016]}\).

The BLA framework offers the user the possibility to obtain both a nonparametric frequency domain estimate of the noise disturbance present at the output and the system combined, and a nonparametric frequency domain estimate of the nonlinear disturbances generated by the system \(\text{[Pintelon & Schoukens, 2012]}\) \(\text{[Schoukens et al., 2012a]}\). These can be used to quantify the level of the nonlinearities in the system and whether or not the nonlinearities or the noise are the dominant disturbance source. Hence, it also allows the user to make a choice on whether or not a nonlinear model should be estimated, based on a quantitative measure. It could be that a linear model is accurate enough.

Exciting the nonlinear system under test at different setpoints of the system (e.g. different input amplitudes or mean values of the input signal) can be very useful to detect what kind of nonlinear effects are present in the system. It is presented in \(\text{[Schoukens et al., 2015a]}\) that a pole shift in the BLAs over the different setpoints indicates the presence of nonlinear feedback loops, while a zero shift indicates the presence of parallel nonlinear signal paths in the system. It are exactly these pole and zero shifts that are exploited by some of the presented identification algorithms to generate initial estimates of the linear dynamic blocks of the selected block-oriented structure. The BLA-based model structure selection approach can be combined with other model structure selection approaches, e.g. \(\text{[Haber & Unbehauen, 1990]}\) \(\text{[Pearson, 2003]}\), to guide the user to a suitable model structure.

It is often not possible to perform an elaborate measurement campaign on the nonlinear system that needs to be modeled. A thorough system analysis and model structure selection step are not always possible under these
Table 1
Overview of the considered model structures and their identification using the BLA, as explained in Sections 6–8. Hammerstein is abbreviated as H, Wiener as W, and Linear Fractional Representation as LFR.

| Model structure | Section | Universal approximator | Number of parameters | Consistent initial values | General difficulties |
|-----------------|---------|------------------------|----------------------|---------------------------|---------------------|
| Hammerstein (H) | 5.1     | no                     | low                  | yes                       | /                   |
| Wiener (W)      | 5.2     | no                     | low                  | yes                       | /                   |
| H-W             | 5.3     | no                     | low                  | no¹                      | assumes an invertible output nonlinearity, iterative procedure |
| W-H             | 5.4     | no                     | low                  | yes                       | mildly computationally expensive brute-force scan |
| Parallel H      | 6.1     | no                     | low/medium           | yes                       | BLA needed at different setpoints |
| Parallel W      | 6.1     | yes/no²                | medium               | yes                       | BLA needed at different setpoints |
| Wiener-Schetzen | 6.2     | yes                    | high³                | yes                       | /                   |
| Parallel W-H    | 6.3     | yes/no²                | high⁴                | yes                       | BLA needed at different setpoints, computationally expensive brute-force scan |
| Simple feedback | 7.1     | no                     | low                  | no                        | /                   |
| W-H feedback    | 7.2     | no                     | low                  | no                        | BLA needed at different setpoints, mildly computationally expensive brute-force scan |
| LFR             | 7.3     | no                     | low/medium           | no                        | BLA needed at different setpoints, mildly computationally expensive scan over solutions of a NARE⁵ |

¹: only consistent in absence of noise and model errors.
²: although the model structure is a universal approximator, the proposed identification methods cannot deal with the general class of Volterra systems; some limitations apply.
³: multiple input single output polynomial.
⁴: multiple input multiple output static nonlinearity.
⁵: nonsymmetric algebraic Riccati equation.

circumstances. Fortunately, most of the presented modeling approaches are relatively simple and can be applied at a fairly low computational cost. This allows the user to try out multiple model structures and compare the quality of the resulting model on a validation data set.

10.2 Input Design

A nonlinear model is in many practical cases only an approximation of the system under study. Therefore, it is of great importance to use input signals that represent a realistic set of excitations to the system. The system should be excited within a realistic range of operation, both in amplitude and frequency. Also the amplitude distribution (e.g. the crest factor of the input) can play an important role. These considerations often exclude relatively simple excitation signals such as sine inputs and impulses.

The BLA framework prefers, but is not limited to, different realizations of periodic, random signals to perform a nonparametric analysis of the nonlinear system under test. An increase in the number of signal realizations decreases the variance on the BLA estimate due to both the nonlinear and noise distortion, while an increase of the number of periods only decreases the variance due to the noise distortion. A typical choice would be to use 7 different realizations and 2 (steady state) periods of a periodic, random signal [Pintelon & Schoukens 2012, Schoukens et al. 2012a]. Some of the presented identification algorithms require the BLA to be measured at different setpoints of the system. An experimental algorithm is introduced in [Esfahani et al. 2016] to find a set of input signal rms-values and dc offsets such that the distortion on the BLA is minimized.

Multisine signals (and more specifically random phase multisines) are excellent candidate input signals since they are periodic and offer a full control of the amplitude spectrum of the input signal [Pintelon & Schoukens 2012]. Small multisine signals can be superposed on a more realistic input signal for the system under test to
obtain a random input signal containing a rich frequency content that can be used for a nonparametric analysis of the nonlinear system under test. An example of this can be found in (Widanage et al., 2011; Marconato et al., 2014).

11 Benchmark and Practical Results

This section provides an overview of the benchmark setups and real-life systems on which the BLA-based block-oriented identification methods have been applied. These results illustrate the good performance and the practical usefulness of the BLA-based block-oriented nonlinear modeling approaches.

The Wiener-Hammerstein identification techniques in Section 6.4 obtained excellent results on the Wiener-Hammerstein benchmark (Schoukens et al., 2009b; Hjalmarsson et al., 2012; Sjöberg et al., 2012). The simple feedback structure and the Wiener-Hammerstein feedback structure have been successfully applied on an electronic simulator of a hardening spring available as the Silverbox benchmark (Schoukens et al., 2006b; Paduart, 2008; Wigren & Schoukens, 2013). The coupled electric drives benchmark in (Wigren & Schoukens, 2013) are an example of a system that is difficult to identify using the BLA based approaches due to the presence of the purely even nonlinearity (see Section 4).

Other systems, not available as a benchmark, that have been successfully modeled by the block-oriented BLA-based approaches are: the insulin-glucose regulatory system (Wiener, Wiener-Schetzen and LFR models) (Marconato et al., 2014; Vanbeylen et al., 2014), a RF crystal detector (Wiener-Hammerstein feedback model) (Schoukens et al., 2008), a logarithmic amplifier (Wiener and parallel Wiener models) (Schoukens & Rolain, 2012c), a valve audio amplifier (Hammerstein and parallel Hammerstein models) (Schoukens et al., 2011) and a Doherty power amplifier (parallel Wiener-Hammerstein model) (Schoukens, 2015).

12 Conclusion

This paper gives an overview of the different block-oriented systems that can be identified based on the best linear approximation, and it gives an overview of the identification algorithms that have been developed in the past. A wide range of systems can be modeled using the best linear approximation approaches: single branch systems, parallel branch systems and nonlinear feedback systems.

The best linear approximation framework allows the user to extract important information about the system (e.g. model order of the dynamics, system structure information), it guides the user in taking good modeling decisions, and it proves to be a good starting point for nonlinear system identification algorithms for a wide range of block-oriented model structures. The best linear approximation provides in some cases a direct estimate of the linear dynamic block of the system. In other cases the best linear approximation can be combined with a pole-zero allocation scan, a singular value decomposition, orthogonal basis functions or a Ricatti-equation based approach to identify the nonlinear system under test.

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