EFFICIENT REDUCED-BASIS TREATMENT OF NONAFFINE AND NONLINEAR PARTIAL DIFFERENTIAL EQUATIONS

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Abstract. In this paper, we extend the reduced-basis approximations developed earlier for linear elliptic and parabolic partial differential equations with affine parameter dependence to problems involving (a) nonaffine dependence on the parameter, and (b) nonlinear dependence on the field variable. The method replaces the nonaffine and nonlinear terms with a coefficient function approximation which then permits an efficient offline-online computational decomposition. We first review the coefficient function approximation procedure: the essential ingredients are (i) a good collateral reduced-basis approximation space, and (ii) a stable and inexpensive interpolation procedure. We then apply this approach to linear nonaffine and nonlinear elliptic and parabolic equations; in each instance, we discuss the reduced-basis approximation and the associated offline-online computational procedures. Numerical results are presented to assess our approach.

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1. INTRODUCTION

The design, optimization, control, and characterization of engineering components or systems often requires repeated, reliable, and real-time prediction of selected performance metrics, or “outputs”, $s$; here superscript “e” shall refer to “exact” and we shall later introduce a “truth approximation” which will bear no superscript. Typical “outputs” include forces, critical stresses or strains, flowrates, or heat fluxes. These outputs are typically functionals of a field variable, $u^e(\mu)$ – such as temperatures or velocities – associated with a parametrized partial differential equation that describes the underlying physics; the parameters, or “inputs”, $\mu$, serve to identify a particular configuration of the component or system – geometry, material properties, boundary conditions, and loads. The relevant system behavior is thus described by an implicit input-output relationship, $s^e(\mu)$, evaluation of which demands solution of the underlying partial differential equation (PDE).

Keywords and phrases. Reduced-basis methods, parametrized PDEs, non-affine parameter dependence, offline-online procedures, elliptic PDEs, parabolic PDEs, nonlinear PDEs.

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The abstract formulation for an elliptic problem can be stated as follows: given any \( \mu \in \mathcal{D} \subset \mathbb{R}^P \), we evaluate \( s^e(\mu) = \ell(u^e(\mu)) \), where \( u^e(\mu) \in X^e \) is the solution of

\[
a(u^e(\mu), v; \mu) = f(v; \mu), \quad \forall v \in X^e.
\]

Here \( \mathcal{D} \) is the parameter domain in which our \( P \)-tuple (input) parameter \( \mu \) resides; \( X^e(\Omega) \) is an appropriate Hilbert space; \( \Omega \) is a bounded domain in \( \mathbb{R}^d \) with Lipschitz continuous boundary \( \partial \Omega \); \( f(\cdot; \mu), \ell(\cdot) \) are \( X^e \)-continuous linear functionals; and \( a(\cdot; \mu) \) is a \( X^e \)-continuous bilinear form.

In actual practice, of course, we do not have access to the exact solution; we thus replace \( u^e(\mu) \) with a “truth” approximation, \( u(\mu) \), which resides in (say) a suitably fine piecewise-linear finite element approximation space \( X \subset X^e \) of very large dimension \( \mathcal{N} \). Our “truth” approximation is thus: given any \( \mu \in \mathcal{D} \), we evaluate

\[
s(\mu) = \ell(u(\mu)),
\]

where \( u(\mu) \in X \) is the solution of

\[
a(u(\mu), v; \mu) = f(v; \mu), \quad \forall v \in X.
\]

We shall assume – hence the appellation “truth” – that the discretization is sufficiently rich such that \( u(\mu) \) and \( u^e(\mu) \) and hence \( s(\mu) \) and \( s^e(\mu) \) are indistinguishable at the accuracy level of interest. The reduced-basis approximation shall be built upon this reference (or “truth”) finite element approximation, and the reduced-basis error will thus be evaluated with respect to \( u(\mu) \in X \). Our formulation must be stable and efficient as \( \mathcal{N} \to \infty \).

We now turn to the abstract formulation for the controlled parabolic case. For simplicity, in this paper we will directly consider a time-discrete framework associated to the time interval \( [0, T] \). We divide \( I \equiv [0, T] \) into \( K \) subintervals of equal length \( \Delta t = \frac{T}{K} \) and define \( t^k = k\Delta t, \, 0 \leq k \leq K \), and \( I \equiv \{t^0, \ldots, t^K\} \); for notational convenience, we also introduce \( \mathcal{K} \equiv \{1, \ldots, K\} \). We shall consider Euler-Backward for the time integration; we can also readily treat higher-order schemes such as Crank-Nicolson [12]. The “truth” approximation is thus: given any \( \mu \in \mathcal{D} \), we evaluate the output

\[
s(\mu, t^k) = \ell(u(\mu, t^k)), \quad \forall k \in \mathcal{K},
\]

where \( u(\mu, t^k) \in X \) satisfies

\[
m(u(\mu, t^k), v) + \Delta t a(u(\mu, t^k), v; \mu) = m(u(\mu, t^{k-1}), v) + \Delta t f(v; \mu) b(t^k), \quad \forall v \in X, \, \forall k \in \mathcal{K},
\]

with initial condition (say) \( u(\mu, t^0) = u_0(\mu) = 0 \). Here, \( f(\cdot; \mu) \) and \( \ell(\cdot) \) are \( Y^e \)-continuous \( (X^e \subset Y^e) \) linear functionals, \( m(\cdot, \cdot) \) is a \( Y^e \)-continuous bilinear form, and \( b(t^k) \) is the control input. We note that the output, \( s(\mu, t^k) \), and the field variable, \( u(\mu, t^k) \), are now functions of the discrete time \( t^k, \forall k \in \mathcal{K} \).

Our goal is the development of numerical methods that permit the rapid yet accurate and reliable prediction of these PDE-induced input-output relationships in real-time or in the limit of many queries – relevant, for example, in the design, optimization, control, and characterization contexts. To achieve this goal we will pursue the reduced-basis method. The reduced-basis method was first introduced in the late 1970s for the nonlinear analysis of structures [1, 25] and subsequently abstracted and analyzed [5, 11, 28, 33] and extended [16, 18, 26] to a much larger class of parametrized partial differential equations. The foundation of the reduced basis method is the realization that, in many instances, the set of all solutions \( u(\mu) \) (say, in the elliptic case) as \( \mu \) varies can be approximated very well by its projection on a finite and low dimensional vector space: for sufficiently well chosen \( \mu_i \), there exist coefficients \( c_i = c_i^N(\mu) \) such that the finite sum \( \sum_{i=1}^N c_i u(\mu_i) \) is very close to \( u(\mu) \) for any \( \mu \).
More recently, the reduced-basis approach and also associated a posteriori error estimation procedures have been successfully developed for (i) linear elliptic and parabolic PDEs that are affine in the parameter [13, 14, 20, 21, 29, 39] – the bilinear form $a(w, v; \mu)$ can be expressed as

$$a(w, v; \mu) = \sum_{q=1}^{Q} \Theta^q(\mu) a^q(w, v),$$

where the $\Theta^q : \mathcal{D} \to \mathbb{R}$ and $a^q(w, v), 1 \leq q \leq Q,$ are parameter dependent functions and parameter-independent bilinear forms, respectively; and (ii) elliptic PDEs that are at most quadratically nonlinear in the first argument [24, 38, 40] – in particular, $a(w, v; \mu)$ satisfies (6) and is at most quadratic in $w$ (but of course linear in $v$).

In these cases a very efficient offline-online computational strategy relevant in the many-query and real-time contexts can be developed. The operation count for the online stage – in which, given a new parameter value, we calculate the reduced-basis output and associated error bound – depends on a low power of the dimension of the “truth” finite element approximation.

Unfortunately, if $a$ is not affine in the parameter this computational strategy breaks down; the online complexity will still depend on $\mathcal{N}$. For example, for general $g(x; \mu)$ (here $x \in \Omega$ and $\mu \in \mathcal{D}$), the bilinear form

$$a(w, v; \mu) \equiv \int_{\Omega} \nabla w \cdot \nabla v + \int_{\Omega} g(x; \mu) w v$$

will not admit an efficient (online $\mathcal{N}$-independent) computational decomposition. In a recent CRAS note [4], we introduce a technique that recovers the efficient offline-online decomposition even in the presence of nonaffine parameter dependence. In this approach, we develop a “collateral” reduced-basis expansion $g_M(x; \mu)$ for $g(x; \mu)$ and then replace $g(x; \mu)$ in (7) with some necessarily affine approximation $g_M(x; \mu) = \sum_{m=1}^{M} \varphi_M m(\mu)q_m(x)$. (Note since we shall also have another reduced-basis expansion for $u(\mu)$, the term “collateral” is used to distinguish the two reduced-basis expansions.) The essential ingredients are (i) a “good” collateral reduced-basis approximation space, $W^\perp_M = \text{span}\{q_m(x), 1 \leq m \leq M\}$, (ii) a stable and inexpensive ($\mathcal{N}$-independent) interpolation procedure by which to determine the $\varphi_M m(\mu), 1 \leq m \leq M$, and (iii) an effective a posteriori estimator with which to quantify the newly introduced error terms. In this paper we shall expand upon the brief presentation in [4] and furthermore address the treatment of nonaffine parabolic problems; we shall also extend the technique to elliptic and parabolic problems in which $g$ is a nonaffine nonlinear function of the field variable $u$ – we hence treat certain classes of nonlinear problems.

A large number of model order reduction (MOR) techniques [2, 7, 8, 22, 27, 32, 36, 41] have been developed to treat nonlinear time-dependent problems. One approach is linearization [41] and polynomial approximation [8, 27]. However, inefficient representation of the nonlinear terms and fast exponential growth (with the degree of the nonlinear approximation order) of the computational complexity render these methods quite expensive, in particular for strong nonlinearities; other approaches for highly nonlinear systems (such as piecewise-linearization) [32, 35] suffer from similar drawbacks. It is also important to note that most MOR techniques focus only on temporal variations; the development of reduced-order models for parametric applications – our focus here – is much less common [6, 9].

This paper is organized as follows: In Section 2 we present a short review of the “empirical interpolation method” – coefficient function approximation – introduced in [4]. The abstract problem formulation, reduced-basis approximation, and computational considerations for linear coercive elliptic and linear coercive parabolic problems with nonaffine parameter dependence are then discussed in Sections 3 and 4, respectively. We extend these results in Section 5 to monotonic nonlinear elliptic PDEs and in Section 6 to monotonic nonlinear parabolic PDEs. Numerical results are included in each section in order to confirm and assess our theoretical results. (Note that, due to space limitations, we do not present in this paper associated a posteriori error estimators; the reader is referred to [4, 12, 23, 37] for a detailed development of this topic.)
2. Empirical interpolation

In this section we describe the empirical interpolation method for constructing the coefficient-function approximation of parameter-dependent functions. We further present a priori and a posteriori error analyses of the method. Finally, we provide a numerical example to illustrate various features of the method.

2.1. Coefficient-function procedure

We begin by summarizing the results in [4]. We consider the problem of approximating a given $\mu$-dependent function $g(\cdot; \mu) \in L^\infty(\Omega) \cap C^0(\Omega), \forall \mu \in D$, of sufficient regularity by a reduced-basis expansion $g_M(\cdot; \mu)$; here, $L^\infty(\Omega) \equiv \{v | \text{ess sup}_{x \in \Omega} |v(x)| < \infty\}$. To this end, we introduce the nested sample sets $S^M_0 = \{\mu^0_1 \in D, \ldots, \mu^0_M \in D\}$, and associated nested reduced-basis spaces $W^M_0 = \text{span} \{\xi_m \equiv g(x; \mu^0_m), 1 \leq m \leq M\}$, in which our approximation $g_M$ shall reside. We then define the best approximation

$$g^*_M(\cdot; \mu) \equiv \arg \min_{z \in W^M_0} \|g(\cdot; \mu) - z\|_{L^\infty(\Omega)}$$

and the associated error

$$\epsilon^*_M(\mu) \equiv \|g(\cdot; \mu) - g^*_M(\cdot; \mu)\|_{L^\infty(\Omega)}.$$  

We note that if $g(\cdot; \mu) \notin C^0(\Omega)$, the best approximation $g^*_M(\cdot; \mu)$ may not be unique. But all the examples considered in this paper have $g(\cdot; \mu) \in C^0(\Omega)$; otherwise, we could use domain decomposition for $g(\cdot; \mu)$. More generally, we can work in a Banach space $B$ that in our context will be $L^\infty(\Omega) \cap C^0(\Omega)$ or $L^2(\Omega)$. Then $g \in C^0(D; B)$ and the forthcoming construction of $S^M_0$ is affected with respect to the $B$ norm.

The construction of $S^M_0$ and $W^M_0$ is based on a greedy selection process. To begin, we choose our first sample point to be $\mu^0_1 = \arg \max_{\mu \in \Xi^0} \|g(\cdot; \mu)\|_{L^\infty(\Omega)}$, and define $S^1_0 = \{\mu^1_1\}, \xi_1 \equiv g(x; \mu^1_1)$, and $W^1_0 = \text{span} \{\xi_1\}$; here $\Xi^0$ is a suitably large but finite-dimensional parameter set in $D$. Then, for $M \geq 2$, we set $\mu^M_M = \arg \max_{\mu \in \Xi^0} \epsilon^*_M(\cdot; \mu)$, and define $S^M_0 = S^{M-1}_0 \cup \mu^M_M, \xi_M = g(x; \mu^M_M), \text{ and } W^M_0 = \text{span} \{\xi_m, 1 \leq m \leq M\}$. In essence, $W^M_0$ comprises basis functions from the parametrically induced manifold $M^y = \{g(\cdot; \mu) | \mu \in D\}$. Thanks to our truth approximation, the optimization for $g^*_M(\cdot; \mu)$ and hence $\epsilon^*_M(\mu)$ is a standard linear program.

We note that the determination of $\mu^M_M$ requires the solution of a linear program for each parameter point in $\Xi^0$; the computational cost thus depends strongly on the size of $\Xi^0$. In the parabolic case this cost may be prohibitively large – at least in our current implementation – if the function $g$ is time-varying either through an explicit dependence on time or (for nonlinear problems) an implicit dependence via the field variable $u(\mu, t^k)$. As we shall see, in these cases the parameter sample $\Xi^0$ is in effect replaced by the parameter-time sample $\Xi^0 \equiv \Xi^0 \times I$; even for modest $K$ the computational cost can be very high. Thus we propose an alternative construction of $S^M_0$: we replace the $L^\infty(\Omega)$-norm in our best approximation by the $L^2(\Omega)$-norm; our next sample point is now given by $\mu^M_M = \arg \max_{\mu \in \Xi^0} \inf_{z \in W^M_{M-1}} \|g(\cdot; \mu) - z\|_{L^2(\Omega)}$, which is relatively inexpensive to evaluate – the computational cost to evaluate $\inf_{z \in W^M_{M-1}} \|g(\cdot; \mu) - z\|_{L^2(\Omega)}$ is $O(MN^2) + O(M^3)$. The following analysis is still rigorous for this alternative (or “surrogate”) construction of $S^M_0$, since we are working in a finite-dimensional space and hence all norms are equivalent; in fact, the $L^\infty(\Omega)$ and $L^2(\Omega)$ procedures yield very similarity convergence results in practice (see Sect. 2.3).

We begin the analysis of our greedy procedure with the following lemma.

**Lemma 2.1.** Suppose that $M_{\max}$ is chosen such that the dimension of span $M^y$ exceeds $M_{\max}$; then, for any $M \leq M_{\max}$, the space $W^M_0$ is of dimension $M$.

**Proof.** It directly follows from our hypothesis on $M_{\max}$ that $\epsilon_0 \equiv \epsilon^*_M(\mu^0_{M+1}) > 0$; our “arg max” construction then implies $\epsilon^*_M(\mu^0_M) \geq \epsilon_0, 2 \leq M \leq M_{\max}$, since $\epsilon^*_M(\mu^0_M) \geq \epsilon^*_M(\mu^0_{M+1}) \geq \epsilon_0$. Now we prove Lemma 2.1 by induction. Clearly, $\dim(W^2_0) = 1$; assume $\dim(W^M_0) = M - 1$; then if $\dim(W^M_0) \neq M$, we have $g(\cdot; \mu^M_M) \in W^M_{M-1}$ and thus $\epsilon^*_M(\mu^M_M) = 0$; however, the latter contradicts $\epsilon^*_M(\mu^M_M) \geq \epsilon_0 > 0$. □
We now construct nested sets of interpolation points \( T_M = \{x_1, \ldots, x_M\}, 1 \leq M \leq M_{\text{max}} \). We first set

\[
x_1 = \arg \sup_{x \in \Omega} |\xi_1(x)|, \quad q_1 = \xi_1(x)/\xi_1(x_1), \quad B_{11}^1 = q_1(x_1) = 1.
\]

Hence, our first interpolation point \( x_1 \) is the maximum point of the first basis function. Then for \( M = 2, \ldots, M_{\text{max}} \), we solve the linear system

\[
\sum_{j=1}^{M-1} \sigma_{ij}^{M-1} q_j(x_i) = \xi_M(x_i), \quad 1 \leq i \leq M - 1;
\]

we calculate

\[
r_M(x) = \xi_M(x) - \sum_{j=1}^{M-1} \sigma_{ij}^{M-1} q_j(x),
\]

and set

\[
x_M = \arg \sup_{x \in \Omega} |r_M(x)|, \quad q_M(x) = r_M(x)/r_M(x_M), \quad B_{ij}^M = q_j(x_i), \quad 1 \leq i, j \leq M.
\]

In essence, the interpolation point \( x_M \) and the basis function \( q_M(x) \) are the maximum point and the normalization of the residual function \( r_M(x) \) which results from the interpolation of \( \xi_M(x) \) by (11).

It remains to demonstrate

**Lemma 2.2.** The construction of the interpolation points is well-defined, and the functions \( \{q_1, \ldots, q_M\} \) form a basis for \( W_M^p \). In addition, the matrix \( B^M \) is lower triangular with unity diagonal.

**Proof.** We shall proceed by induction. Clearly, we have \( W_1^p = \text{span} \{q_1\} \). Next we assume \( W_{M-1}^p = \text{span} \{q_1, \ldots, q_{M-1}\} \); if (i) \( B_{M-1}^{M-1} \) is invertible and (ii) \( |r_M(x_M)| > 0 \), then our construction may proceed and we may form \( W_M^p = \text{span} \{q_1, \ldots, q_M\} \). To prove (i), we just note from the construction procedure that \( B_{ij}^{M-1} = r_j(x_i)/r_j(x_j) = 0 \) for \( i < j \); that \( B_{ij}^{M-1} = r_j(x_i)/r_j(x_j) = 1 \) for \( i = j \); and that \( |B_{ij}^{M-1}| = |r_j(x_i)/r_j(x_j)| \leq 1 \) for \( i > j \) since \( x_j = \arg \sup_{x \in \Omega} |r_j(x)|, 1 \leq j \leq M \). Hence, \( B^M \) is lower triangular with unity diagonal. To prove (ii) (and hence also that the \( x_i, 1 \leq i \leq M, \) are distinct), we observe that \( |r_M(x_M)| \geq \varepsilon_{M-1}(\mu_M^p) \geq \varepsilon_0 > 0 \) since \( \varepsilon_{M-1}(\mu_M^p) \) is the error associated with the best approximation. \( \square \)

Furthermore, from the invertibility of \( B^M \), we immediately derive

**Lemma 2.3.** For any \( M \)-tuple \( (\alpha_i)_{i=1, \ldots, M} \) of real numbers, there exists a unique element \( w \in W_M^p \) such that \( w(x_i) = \alpha_i, 1 \leq i \leq M \).

It remains to develop an efficient procedure for obtaining a good collateral reduced-basis expansion \( g_M(\cdot; \mu) \). Based on the approximation space \( W_M^p \) and set of interpolation points \( T_M \), we can readily construct an approximation to \( g(x; \mu) \). Indeed, our coefficient function approximation is the interpolant of \( g \) over \( T_M \) as provided for from Lemma 2.3:

\[
g_M(x; \mu) = \sum_{m=1}^M \varphi_M m(\mu) q_m(x),
\]

where \( \varphi_M(\mu) \in \mathbb{R}^M \) is given by

\[
\sum_{j=1}^M B_{ij}^M \varphi_M j(\mu) = g(x_i; \mu), \quad 1 \leq i \leq M;
\]

note that \( g_M(x_i; \mu) = g(x_i; \mu), 1 \leq i \leq M \). We define the associated error as

\[
\varepsilon_M(\mu) \equiv \|g(\cdot; \mu) - g_M(\cdot; \mu)\|_{L^\infty(\Omega)}.
\]
It remains to understand how well \( g_{\mu} (x; \mu) \) approximates \( g(x; \mu) \).

### 2.2. Error analysis

#### 2.2.1. A priori stability: Lebesgue constant

To begin, we define a “Lebesgue constant” \([10,31,34]\) \( \Lambda_{\mu} = \sup_{x \in \Omega} \sum_{m=1}^{M} |V_{m}^{M} (x)| \). Here, the \( V_{m}^{M} (x) \in W^{g}_{M} \) are characteristic functions satisfying \( V_{m}^{M} (x_{m}) = \delta_{mn}, 1 \leq m, n \leq M \), the existence and uniqueness of which is guaranteed by Lemma 2.3; here \( \delta_{mn} \) is the Kronecker delta symbol. It can be shown that

**Lemma 2.4.** The set of all characteristic functions \( \{V_{m}^{M}\}_{m=1}^{M} \) is a basis for \( W^{g}_{M} \). Furthermore, the two bases \( q_{m}, 1 \leq m \leq M \), and \( V_{m}^{M}, 1 \leq m \leq M \), are related by

\[
q_{i} (x) = \sum_{j=1}^{M} B_{j}^{M} V_{j}^{M} (x), \quad 1 \leq i \leq M. \tag{17}
\]

**Proof.** It is immediate from the definition of the \( V_{m}^{M} \) that the set of all characteristic functions \( \{V_{m}^{M}\}_{m=1}^{M} \) is linearly independent. This set thus constitutes a basis for \( W^{g}_{M} \), in fact a nodal basis associated with the set \( \{x_{m}\}_{m=1}^{M} \). Then, we consider \( x = x_{m}, 1 \leq n \leq M \), and note that

\[
\sum_{j=1}^{M} B_{j}^{M} V_{j}^{M} (x_{n}) = \sum_{j=1}^{M} B_{j}^{M} \delta_{jn} = B_{n}^{M} = q_{i} (x_{n}), \quad 1 \leq i \leq M; \tag{17}
\]

it thus follows from Lemma 2.3 that (17) holds.

We observe that \( \Lambda_{M} \) depends on \( W^{g}_{M} \) and \( T_{M} \), but not on \( \mu \). We can further prove

**Lemma 2.5.** The interpolation error \( \varepsilon_{M} (\mu) \) satisfies \( \varepsilon_{M} (\mu) \leq \varepsilon_{M}^{*} (\mu) (1 + \Lambda_{M}), \ \forall \mu \in D \).

**Proof.** We first introduce \( \varepsilon_{M}^{*} (x; \mu) = g(x; \mu) - g_{M}^{*} (x; \mu) \). It then follows that

\[
g_{M} (x; \mu) - g_{M}^{*} (x; \mu) = \sum_{m=1}^{M} \left( g_{M} (x_{m}; \mu) - g_{M}^{*} (x_{m}; \mu) \right) V_{m}^{M} (x)
\]

\[
= \sum_{m=1}^{M} ((g_{M} (x_{m}; \mu) - g(x_{m}; \mu)) + (g(x_{m}; \mu) - g_{M}^{*} (x_{m}; \mu))) V_{m}^{M} (x)
\]

\[
= \sum_{m=1}^{M} \varepsilon_{M}^{*} (x_{m}; \mu) V_{m}^{M} (x). \tag{18}
\]

Furthermore, from the definition of \( \varepsilon_{M} (\mu) \) and \( \varepsilon_{M}^{*} (\mu) \), and the triangle inequality, we obtain

\[
\varepsilon_{M} (\mu) = \| g(\cdot; \mu) - g_{M}(\cdot; \mu) \|_{L^{\infty}(\Omega)} \leq \varepsilon_{M}^{*} (\mu) + \| g_{M}(\cdot; \mu) - g_{M}^{*}(\cdot; \mu) \|_{L^{\infty}(\Omega)}.
\]

This yields, from (18),

\[
\varepsilon_{M} (\mu) - \varepsilon_{M}^{*} (\mu) \leq \| g_{M}(\cdot; \mu) - g_{M}^{*}(\cdot; \mu) \|_{L^{\infty}(\Omega)}
\]

\[
= \max_{i \in [1, \ldots, M]} \| e_{i}^{*} (x_{i}; \mu) \|_{L^{\infty}(\Omega)}
\]

the desired result then immediately follows from \( |e_{i}^{*} (x_{i}; \mu)| \leq \varepsilon_{M}^{*} (\mu), \ 1 \leq i \leq M. \) \( \square \)

We can further show
Proposition 2.6. The Lebesgue constant \( \Lambda_M \) satisfies \( \Lambda_M \leq 2^M - 1 \).

Proof. We first recall two crucial properties of the matrix \( B^M \): (i) \( B^M \) is lower triangular with unity diagonal and \( q_m(x_m) = 1, 1 \leq m \leq M \), and (ii) all entries of \( B^M \) are of modulus no greater than unity \( -\|q_m\|_{L^\infty(\Omega)} \leq 1, 1 \leq m \leq M \). Hence, from (17) we can write

\[
|V^M_m(x)| = |q_m(x) - \sum_{i=m+1}^{M} B^M_{i,m} V^M_i(x)| \\
\leq 1 + \sum_{i=m+1}^{M} |V^M_i(x)|, \quad 1 \leq m \leq M - 1.
\]

It follows that, starting from \( |V^M_M(x)| = |q_M(x)| \leq 1 \), we can deduce \( |V^M_{M+1-m}(x)| \leq 1 + |V^M_M(x)| + \ldots + |V^M_{M+2-m}(x)| \leq 2^{m-1}, 2 \leq m \leq M \), and thus obtain \( \sum_{m=1}^{M} |V^M_m(x)| \leq 2^M - 1 \).

Proposition 2.6 is very pessimistic and of little practical value (though \( \varepsilon_M^*(\mu) \) does often converge sufficiently rapidly that \( \varepsilon_M^*(\mu) 2^M \to 0 \) as \( M \to \infty \)); this is not surprising given analogous results in the theory of polynomial interpolation [10, 31, 34]. In applications, the actual asymptotic behavior of \( \Lambda_M \) is much lower than the upper bound of Proposition 2.6; however, Proposition 2.6 does provide a theoretical basis for some stability.

2.2.2. A posteriori estimators

Given an approximation \( g_M(x; \mu) \) for \( M \leq M_{\text{max}} - 1 \), we define \( \mathcal{E}_M(x; \mu) \equiv \hat{\varepsilon}_M(x; \mu) q_{M+1}(x) \), where \( \hat{\varepsilon}_M(x; \mu) \equiv |g(x_{M+1}; \mu) - g_M(x_{M+1}; \mu)| \). In general, \( \varepsilon_M(\mu) \geq \hat{\varepsilon}_M(\mu) \), since \( \varepsilon_M(\mu) = ||g(\cdot; \mu) - g(\cdot; \mu)||_{L^\infty(\Omega)} \geq |g(x; \mu) - g_M(x; \mu)| \) for all \( x \in \Omega \), and thus also for \( x = x_{M+1} \). However, we can prove

Proposition 2.7. If \( g(\cdot; \mu) \in W_{M+1}^q \), then (i) \( g(x; \mu) - g_M(x; \mu) = \pm \mathcal{E}_M(x; \mu) \), and (ii) \( ||g(\cdot; \mu) - g_M(\cdot; \mu)||_{L^\infty(\Omega)} = \hat{\varepsilon}_M(\mu) \).

Proof. By our assumption \( g(\cdot; \mu) \in W_{M+1}^q \), there exists \( \kappa(\mu) \in \mathbb{R}^{M+1} \) such that \( g(x; \mu) - g_M(x; \mu) = \sum_{m=1}^{M+1} \kappa_m(\mu) q_m(x) \). We now consider \( x = x_i, 1 \leq i \leq M + 1 \), and arrive at

\[
\sum_{m=1}^{M+1} \kappa_m(\mu) q_m(x_i) = g(x_i; \mu) - g_M(x_i; \mu), \quad 1 \leq i \leq M + 1.
\]

It thus follows that \( \kappa_m(\mu) = 0, 1 \leq m \leq M \), since \( g(x_i; \mu) - g_M(x_i; \mu) = 0, 1 \leq i \leq M \), and the matrix \( q_m(x_i) = B_m^{(i)} \) is lower triangular, and that \( \kappa_{M+1}(\mu) = g(x_{M+1}; \mu) - g_M(x_{M+1}; \mu) \) since \( q_{M+1}(x_{M+1}) = 1 \); this concludes the proof of (i). The proof of (ii) then directly follows from \( ||q_{M+1}||_{L^\infty(\Omega)} = 1 \).

Of course, in general \( g(\cdot; \mu) \not\in W_{M+1}^q \), and hence our estimator \( \hat{\varepsilon}_M(\mu) \) is unfortunately a lower bound. However, if \( \varepsilon_M(\mu) \to 0 \) very fast, we expect that the effectivity,

\[
\eta_M(\mu) \equiv \frac{\hat{\varepsilon}_M(\mu)}{\varepsilon_M(\mu)}
\]

shall be close to unity; furthermore, the estimator is very inexpensive – one additional evaluation of \( g(\cdot; \mu) \) at a single point in \( \Omega \). (Note we can readily improve the rigor of our bound at only modest additional cost: if we assume that \( g(\cdot; \mu) \in W_{M+k}^q \), then \( \hat{\varepsilon}_M = 2^{k-1} \max_{1 \leq i \leq k} |g(x_{M+k}; \mu) - g_M(x_{M+k}; \mu)| \) is an upper bound for \( \varepsilon_M(\mu) \) (see Props. 2.6 and 2.7.).

We refer to [4, 12, 23] for the incorporation of these error estimators into output bounds for reduced basis approximations of nonaffine partial differential equations.
Figure 1. (a) Parameter sample set $S_M^g$, $M_{\text{max}} = 51$, and (b) interpolation points $x_m$, $1 \leq m \leq M_{\text{max}}$, for the function $G(x; \mu)$ of (20).

2.3. Numerical results

We consider the function $g(\cdot; \mu) = G(\cdot; \mu)$, where

$$G(x; \mu) \equiv \frac{1}{\sqrt{(x(1) - \mu(1))^2 + (x(2) - \mu(2))^2}}$$

for $x = (x(1), x(2)) \in \Omega \equiv [0, 1]^2$ and $\mu \in \mathcal{D} \equiv [-1, -0.01]^2$. We choose for $\Xi^g$ a deterministic grid of $40 \times 40$ parameter points over $\mathcal{D}$. We take $\mu^0 = (-0.01, -0.01)$ and then pursue the empirical interpolation procedure described in Section 2.1 to construct $S_M^g$, $W_M^g$, $T_M$, and $B^M$, $1 \leq M \leq M_{\text{max}}$, for $M_{\text{max}} = 51$. We see that the parameter points in $S_M^g$, shown in Figure 1a, are mainly distributed around the corner $(-0.01, -0.01)$ of the parameter domain; and that the interpolation points in $T_M$, plotted in Figure 1b, are largely allocated around the corner $(0, 0)$ of the physical domain $\Omega$. This is because $G(x; \mu)$ varies most significantly at $\mu = (-0.01, -0.01)$ (upper right corner of Fig. 1a) and $x = (0, 0)$ (lower left corner of Fig. 1b).

We now introduce a parameter test sample $\Xi_{\text{prop}}^g$ of size $Q_{\text{Test}} = 225$, and define $\varepsilon_{\text{max}} = \max_{\mu \in \Xi_{\text{prop}}^g} \varepsilon_M(\mu)$, $\varepsilon_{\text{M, max}} = \max_{\mu \in \Xi_{\text{prop}}^g} \varepsilon_M(\mu)$, $\eta_M(\mu) = Q_{\text{Test}}^{-1} \sum_{\mu \in \Xi_{\text{prop}}^g} (\varepsilon_M(\mu)/\varepsilon_M^*(\mu))$, $\tau_M = Q_{\text{Test}}^{-1} \sum_{\mu \in \Xi_{\text{prop}}^g} \eta_M(\mu)$, and $\kappa_M$; here $\eta_M(\mu)$ is the effectivity defined in (19), and $\kappa_M$ is the condition number of $B^M$. We present in Table 1 $\varepsilon_{\text{M, max}}$, $\varepsilon_M^*(\mu)$, $\tau_M$, $\Lambda_M$, $\kappa_M$, and $\kappa_M$ as a function of $M$. We observe that $\varepsilon_{\text{M, max}}$ and $\varepsilon^*_M(\mu)$ converge rapidly with $M$; that the Lebesgue constant provides a reasonably sharp measure of the interpolation-induced error; that the Lebesgue constant grows very slowly – and hence $\varepsilon_M(\mu)$ will be only slightly larger than the min max result $\varepsilon^*_M(\mu)$; that the error estimator effectivity is reasonably close to unity (note the last column of the Tab. 1 in [4] (analogous to Tab. 1 here) contains an error – we purported in [4] to report the average of $\eta_M(\mu)$ $\equiv \varepsilon_M(\mu)/\varepsilon_M^*(\mu)$ over $\Xi_{\text{Test}}^g$, but in fact we reported the average over $\Xi_{\text{Test}}^g$ of $\varepsilon_M(\mu)/\varepsilon_M^*(\mu)$); and that $B^M$ is quite well-conditioned for our choice of basis (For the non-orthogonalized basis $\xi_m$, $1 \leq m \leq M$, the condition number of $B^M$ will grow exponentially with $M$.) These results are expected: although $G(x; \mu)$ varies rapidly as $\mu$ approaches 0 and $x$ approaches 0, $G(x; \mu)$ is nevertheless quite smooth in the prescribed parameter domain $\mathcal{D}$.

If we exploit the $L^2(\Omega)$-norm surrogate in our best approximation we can construct $S_M^g$ much less expensively. We present in Table 2 numerical results obtained from this alternative construction of $S_M^g$. The results are very similar to those in Table 1, which implies – as expected – that the approximation quality of our empirical interpolation approach is relatively insensitive to the choice of norm in the sample construction process.
Table 1. Numerical results for empirical interpolation of $G(x; \mu)$: $\varepsilon_{M, \text{max}}$, $\varepsilon_{M, \text{max}}^*$, $\bar{\eta}_M$, $\Lambda_M$, $\bar{\eta}_M$, and $\varkappa_M$ as a function of $M$.

<table>
<thead>
<tr>
<th>$M$</th>
<th>$\varepsilon_{M, \text{max}}$</th>
<th>$\varepsilon_{M, \text{max}}^*$</th>
<th>$\bar{\eta}_M$</th>
<th>$\Lambda_M$</th>
<th>$\bar{\eta}_M$</th>
<th>$\varkappa_M$</th>
</tr>
</thead>
<tbody>
<tr>
<td>8</td>
<td>1.72E–01</td>
<td>8.30E–02</td>
<td>0.68</td>
<td>1.76</td>
<td>0.17</td>
<td>3.65</td>
</tr>
<tr>
<td>16</td>
<td>1.42E–02</td>
<td>4.22E–03</td>
<td>0.67</td>
<td>2.63</td>
<td>0.10</td>
<td>6.08</td>
</tr>
<tr>
<td>24</td>
<td>1.01E–03</td>
<td>2.68E–04</td>
<td>0.49</td>
<td>4.42</td>
<td>0.28</td>
<td>9.19</td>
</tr>
<tr>
<td>32</td>
<td>2.31E–04</td>
<td>5.64E–05</td>
<td>0.48</td>
<td>5.15</td>
<td>0.20</td>
<td>12.86</td>
</tr>
<tr>
<td>40</td>
<td>1.63E–05</td>
<td>3.66E–06</td>
<td>0.54</td>
<td>4.98</td>
<td>0.60</td>
<td>18.37</td>
</tr>
<tr>
<td>48</td>
<td>2.44E–06</td>
<td>6.08E–07</td>
<td>0.37</td>
<td>7.43</td>
<td>0.29</td>
<td>20.41</td>
</tr>
</tbody>
</table>

Table 2. Numerical results for empirical interpolation of $G(x; \mu)$: $\varepsilon_{M, \text{max}}$, $\varepsilon_{M, \text{max}}^*$, $\bar{\eta}_M$, $\Lambda_M$, $\bar{\eta}_M$, and $\varkappa_M$ as a function of $M$; here $S_M^g$ is constructed with the $L^2(\Omega)$-norm as a surrogate for the $L^\infty(\Omega)$-norm.

<table>
<thead>
<tr>
<th>$M$</th>
<th>$\varepsilon_{M, \text{max}}$</th>
<th>$\varepsilon_{M, \text{max}}^*$</th>
<th>$\bar{\eta}_M$</th>
<th>$\Lambda_M$</th>
<th>$\bar{\eta}_M$</th>
<th>$\varkappa_M$</th>
</tr>
</thead>
<tbody>
<tr>
<td>8</td>
<td>2.69E–01</td>
<td>1.18E–01</td>
<td>0.66</td>
<td>2.26</td>
<td>0.23</td>
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</tr>
<tr>
<td>16</td>
<td>1.77E–02</td>
<td>3.96E–03</td>
<td>0.45</td>
<td>4.86</td>
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<td>24</td>
<td>8.07E–04</td>
<td>3.83E–04</td>
<td>0.43</td>
<td>3.89</td>
<td>0.28</td>
<td>13.53</td>
</tr>
<tr>
<td>32</td>
<td>1.69E–04</td>
<td>3.92E–05</td>
<td>0.45</td>
<td>7.07</td>
<td>0.47</td>
<td>16.60</td>
</tr>
<tr>
<td>40</td>
<td>2.51E–05</td>
<td>4.10E–06</td>
<td>0.43</td>
<td>6.40</td>
<td>0.25</td>
<td>18.84</td>
</tr>
<tr>
<td>48</td>
<td>2.01E–06</td>
<td>6.59E–07</td>
<td>0.30</td>
<td>8.86</td>
<td>0.18</td>
<td>21.88</td>
</tr>
</tbody>
</table>

3. NONAFFINE LINEAR COERCIVE ELLIPTIC EQUATIONS

We now incorporate the coefficient-function procedure described in Section 2.1 into our reduced-basis method to develop an efficient reduced-basis approximation for the nonaffine linear coercive elliptic equations. The presentation of this section is as follows: first, we introduce an abstract problem statement of the nonaffine elliptic equation as well as a model problem; we then develop our reduced-basis approximation for the abstract problem and discuss computational complexity; finally, we present numerical results obtained for the model problem.

3.1. Problem formulation

3.1.1. Abstract statement

We first define the Hilbert spaces $X^e \equiv H^1_0(\Omega)$ – or, more generally, $H_0^1(\Omega) \subset X^e \subset H^1(\Omega)$ – where $H^1(\Omega) = \{v \mid v \in L^2(\Omega), \nabla v \in (L^2(\Omega))^d\}$, $H^1_0(\Omega) = \{v \mid v \in H^1(\Omega), v|_{\partial \Omega} = 0\}$. The inner product and norm associated with $X^e$ are given by $(\cdot, \cdot)_{X^e}$ and $\|\cdot\|_{X^e} = (\cdot, \cdot)^{1/2}_{X^e}$, respectively; for example, $(w,v)_{X^e} \equiv \int_{\Omega} \nabla w \cdot \nabla v + \int_{\Omega} w v$, $\forall w, v \in X^e$. The truth approximation subspace $X$ shall inherit this inner product and norm: $(\cdot, \cdot)_X \equiv (\cdot, \cdot)^{1/2}_X$ and $\|\cdot\|_X \equiv \|\|\|_X$.

In this section, we are interested in a particular form for problem (1), in which

$$a(w, v; \mu) = a_0(w, v) + a_1(w, v, g(\cdot; \mu)), \quad (21)$$

and

$$f(v; \mu) = \int_{\Omega} vh(x; \mu), \quad (22)$$

where $a_0(\cdot, \cdot)$ is a (for simplicity, parameter-independent) bilinear form, $a_1 : X^e \times X^e \times L^\infty(\Omega)$ is a trilinear form, and $g(\cdot; \mu) \in L^\infty(\Omega)$, $h(\cdot; \mu) \in L^\infty(\Omega)$ are prescribed functions. For simplicity of exposition, we presume that $h(x; \mu) = g(x; \mu)$. 

We shall assume that \( a \) satisfies coercivity and continuity conditions
\[
0 < \alpha_0 \leq \alpha(\mu) \equiv \inf_{w \in X \setminus \{0\}} \frac{a(w, w; \mu)}{\|w\|_X^2}, \quad \forall \mu \in \mathcal{D},
\]
(23)
and \( a_1(\zeta_i, \zeta_j, g(\cdot; \mu)) \) depend on \( g(x; \mu) \) and must thus be evaluated online for every new parameter value \( \mu \); the samples, indeed, we shall continue the motivation discussed in Section 1. Specifically, we introduce the nested Lagrange-Hermitian spaces \([17]\). We then orthonormalize the \( \zeta \)
here \( \alpha(\mu) \) and \( \gamma(\mu) \) are the coercivity constant and the continuity constant, respectively. (We (plausibly) suppose that \( \alpha_0, \gamma_0 \) may be chosen independent of \( \mathcal{N} \)) We shall further assume that the trilinear form \( a_1 \) satisfies
\[
a_1(w, v, z) \leq \gamma_1 \|w\|_X \|v\|_X \|z\|_{L^\infty(\Omega)}, \quad \forall w, v, z \in L^\infty(\Omega),
\]
(25)
It is then standard, given that \( g(\cdot; \mu) \in L^\infty(\Omega) \), to prove existence and uniqueness of the exact solution and the truth approximation.

3.1.2. A model problem
We consider the following model problem defined on the unit square \( \Omega = [0, 1]^2 \in \mathbb{R}^2 \): Given the parameter input \( \mu = (\mu(1), \mu(2)) \in \mathcal{D} \equiv [-1, -0.01]^2 \), the field variable \( u(\mu) \in X \) satisfies (3), where \( X \subset X^e \equiv H_0^1(\Omega) \) is a piecewise-linear finite element approximation space of dimension \( N = 2601 \). Here \( a \) is given by (21) for
\[
a_0(w, v) = \int_{\Omega} \nabla w \cdot \nabla v, \quad a_1(w, v, g(\cdot; \mu)) = \int_{\Omega} g(x; \mu) \, w, \quad \forall \mu \in \mathcal{D},
\]
(26)
for \( g(x; \mu) = G(x; \mu) \) as defined in (20); and \( f \) is given by (22) for \( h(x; \mu) = g(x; \mu) = G(x; \mu) \). The output \( s(\mu) \) is evaluated as \( s(\mu) = \ell(u(\mu)) \) for \( \ell(v) = \int_{\Omega} v \). The solution \( u(\mu) \) develops a boundary layer in the vicinity of \( x = (0, 0) \) for \( \mu \) near the “corner” \((-0.01, -0.01)\).

3.2. Reduced-basis approximation

3.2.1. Discrete equations
We begin with motivating the need for the empirical interpolation approach in dealing with nonaffine problems; indeed, we shall continue the motivation discussed in Section 1. Specifically, we introduce the nested samples, \( S_N = \{\mu_1^N, \ldots, \mu_N^N \in \mathcal{D} \} \), \( 1 \leq N \leq N_{\max} \), and associated nested Lagrangian [28] reduced-basis spaces \( W_N^{\mathcal{N}} = \text{span} \{ \zeta_j, j = 1, \ldots, N \} \), \( 1 \leq N \leq N_{\max} \), where \( u(\mu) \) is the solution of (3) for \( \mu = \mu_0 \). (Note we may also consider Hermitian spaces built upon sensitivity derivatives of \( u \) with respect to \( \mu \) [15] or, more generally, Lagrange-Hermitian spaces [17].) We then orthonormalize the \( \zeta_j \), \( 1 \leq j \leq N \), with respect to \( (\cdot, \cdot)_X \) so that \( (\zeta_i, \zeta_j)_X = \delta_{ij} \), \( 1 \leq i, j \leq N \); the resulting algebraic system will be well-conditioned.

Were to follow the classical recipe, the reduced-basis approximation would be obtained by a standard Galerkin projection: given \( \mu \in \mathcal{D} \), we evaluate \( s_N(\mu) = \ell(u_N(\mu)) \), where \( u_N(\mu) \in W_N^0 \) is the solution of
\[
a_0(u_N(\mu), v) + a_1(u_N(\mu), v, g(\cdot; \mu)) = \int_{\Omega} g(x; \mu) v, \quad \forall v \in W_N^0.
\]
(27)
If we now express \( u_N(\mu) = \sum_{j=1}^N u_N j(\mu) \zeta_j \) and choose test functions \( v = \zeta_n \), \( 1 \leq n \leq N \), in (27), we obtain the \( N \times N \) linear algebraic system
\[
\sum_{j=1}^N (a_0(\zeta_i, \zeta_j) + a_1(\zeta_i, \zeta_j, g(\cdot; \mu))) u_N j(\mu) = \int_{\Omega} g(x; \mu) \zeta_i, \quad 1 \leq i \leq N.
\]
(28)
We observe that while \( a_0(\zeta_i, \zeta_j) \) is parameter-independent and can thus be pre-computed offline, \( \int_{\Omega} g(x; \mu) \zeta_i \) and \( a_1(\zeta_i, \zeta_j, g(\cdot; \mu)) \) depend on \( g(x; \mu) \) and must thus be evaluated online for every new parameter value \( \mu \); the
operation count for the online stage will thus scale as \( O(N^2N) \), where \( N \) is the dimension of the underlying truth finite element approximation space. The decrease in marginal cost in replacing the truth finite element approximation space with the reduced-basis approximation will be quite modest regardless of the dimension reduction \( N \ll N \).

To recover online \( N \)-independence, we appeal to the empirical interpolation method discussed in Section 2. We simply replace \( g(x; \mu) \) in (28) with the (necessarily) affine approximation \( g_M(x; \mu) = \sum_{m=1}^M \varphi_M(m(\mu))q_m(x) \) from (14) based upon the empirical interpolation approach described in Section 2. Our reduced-basis approximation is then: Given \( \mu \in \mathcal{D} \), find \( u_{N,M}(\mu) \in W^N_N \) such that

\[
\begin{align*}
a_0(u_{N,M}(\mu), v) + a_1(u_{N,M}(\mu), v, g_M(\cdot; \mu)) &= \int_{\Omega} g_M(x; \mu)v, \quad \forall v \in W^N_N; \\
\end{align*}
\]

we then evaluate the output estimate from

\[
s_{N,M}(\mu) = \ell(u_{N,M}(\mu)).
\]

We now express \( u_{N,M}(\mu) = \sum_{j=1}^N u_{N,M,j}(\mu) \zeta_j \), choose as test functions \( v = \zeta_n, \quad 1 \leq n \leq N \), and invoke (14) to obtain

\[
\sum_{j=1}^N \left( a_0(\zeta_i, \zeta_j) + \sum_{m=1}^M \varphi_M(m(\mu))a_1(\zeta_i, \zeta_j, q_m) \right) u_{N,M,j}(\mu) = \sum_{m=1}^M \varphi_M(m(\mu)) \int_{\Omega} \zeta_i q_m, \quad 1 \leq i \leq N,
\]

where \( \varphi_M(m(\mu)), \quad 1 \leq m \leq M \), is determined from (15). We indeed recover the online \( N \)-independence: the quantities \( a_0(\zeta_i, \zeta_j), a_1(\zeta_i, \zeta_j, q_m) \), and \( \int_{\Omega} \zeta_i q_m \) are all parameter independent and can thus be pre-computed offline, as discussed further in Section 3.2.3.

3.2.2. A priori theory

We consider here the convergence rate of \( u_{N,M}(\mu) \rightarrow u(\mu) \). In fact, it is a simple matter to demonstrate the optimality of \( u_{N,M}(\mu) \) in

**Proposition 3.1.** For \( \varepsilon_M(\mu) \) of (16) satisfying \( \varepsilon_M(\mu) \leq \frac{1}{2} \frac{\alpha(\mu)}{\phi_2(\mu)} \), we have

\[
\|u(\mu) - u_{N,M}(\mu)\|_X \leq \left( 1 + \frac{\gamma(\mu)}{\alpha(\mu)} \right) \inf_{w_N \in W^N_N} \|u(\mu) - w_N\|_X + \varepsilon_M(\mu) \left( \frac{\phi_1(\mu)\alpha(\mu) + 2\phi_2(\mu)\phi_3(\mu)}{\alpha^2(\mu)} \right);
\]

here \( \phi_1(\mu), \phi_2(\mu), \) and \( \phi_3(\mu) \) are given by

\[
\phi_1(\mu) = \frac{1}{\varepsilon_M(\mu)} \sup_{v \in X} \int_{\Omega} v(g(x; \mu) - g_M(x; \mu)) / \|v\|_X,
\]

\[
\phi_2(\mu) = \frac{1}{\varepsilon_M(\mu)} \sup_{w \in X} \sup_{v \in X} \frac{a_1(w, v; g(\cdot; \mu) - g_M(\cdot; \mu))}{\|w\|_X \|v\|_X},
\]

\[
\phi_3(\mu) = \sup_{v \in X} \int_{\Omega} v g_M(x; \mu) / \|v\|_X.
\]

**Proof.** For any \( w_N = u_{N,M}(\mu) + v_N \in W^N_N \), we have

\[
\begin{align*}
\alpha(\mu)\|w_N - u_{N,M}(\mu)\|^2_X &\leq \alpha(w_N - u_{N,M}(\mu), w_N - u_{N,M}(\mu); \mu) \\
&= \alpha(w_N - u(\mu), v_N; \mu) + \alpha(u(\mu) - u_{N,M}(\mu), v_N; \mu) \\
&\leq \gamma(\mu)\|w_N - u(\mu)\|_X \|v_N\|_X + \alpha(u(\mu) - u_{N,M}(\mu), v_N; \mu).
\end{align*}
\]
Note further from (3), (29), and (33)–(35) that the second term can be bounded by
\[
\alpha(u(\mu) - u_{N,M}(\mu), v_N; \mu) = \int_\Omega v_N g(x; \mu) - a(u_{N,M}(\mu), v_N; \mu)
\]
\[
= \int_\Omega v_N (g(x; \mu) - g_M(x; \mu)) - a_1(u_{N,M}(\mu), v_N; g(x; \mu) - g_M(x; \mu))
\]
\[
\leq \varepsilon_M(\mu) \phi_1(\mu) \|v_N\|_X + \varepsilon_M(\mu) \phi_2(\mu) \|v_N\|_X \|u_{N,M}(\mu)\|_X
\]
\[
\leq \varepsilon_M(\mu) \left( \frac{\phi_1(\mu) \alpha(\mu) + 2\phi_2(\mu) \phi_3(\mu)}{\alpha(\mu)} \right) \|v_N\|_X,
\]
where the last inequality derives from
\[
\alpha(\mu) \|u_{N,M}(\mu)\|_X^2 \leq a(u_{N,M}(\mu), u_{N,M}(\mu); \mu)
\]
\[
= \int_\Omega u_{N,M}(\mu) g_M(x; \mu) + a_1(u_{N,M}(\mu), u_{N,M}(\mu); g(x; \mu) - g_M(x; \mu))
\]
\[
\leq \phi_3(\mu) \|u_{N,M}(\mu)\|_X + \varepsilon_M(\mu) \phi_2(\mu) \|u_{N,M}(\mu)\|_X^2,
\]
and our hypothesis on \( \varepsilon_M(\mu) \). It then follows from (36) and (37) that
\[
\|w_N - u_{N,M}(\mu)\|_X \leq \frac{\gamma(\mu)}{\alpha(\mu)} \|w_N - u(\mu)\|_X + \varepsilon_M(\mu) \left( \frac{\phi_1(\mu) \alpha(\mu) + 2\phi_2(\mu) \phi_3(\mu)}{\alpha^2(\mu)} \right), \quad \forall w_N \in W_N^\mu.
\]
The desired result finally follows from (39) and the triangle inequality. (Note that \( \phi_1, \phi_2, \) and \( \phi_3 \) are bounded by virtue of our continuity requirements.)

We note from Proposition 3.1 that \( M \) should be chosen such that \( \varepsilon_M(\mu) \) is of the same order as the error in the best approximation, \( \inf_{w_N \in W_N^\mu} \|u(\mu) - w_N\|_X \), as otherwise the second term on the right-hand side of (32) may limit the convergence of the reduced-basis approximation. As regards the error in the best approximation, we note that \( W_N^\mu \) comprises “snapshots” on the parametrically induced manifold \( \mathcal{M}^\mu \equiv \{ u(\mu) \mid \forall \mu \in \mathcal{D} \} \subset X \). The critical observations are that \( \mathcal{M}^\mu \) is very low-dimensional and that \( \mathcal{M}^\mu \) is smooth under general hypotheses on stability and continuity. We thus expect that the best approximation will converge to \( u(\mu) \) very rapidly, and hence that \( N \) may be chosen small. (This is proven for a particularly simple case in [21].)

3.2.3. Offline-online procedure

We summarize here the procedure [3,18,20,29]. In the offline stage – performed only once – we first construct nested approximation spaces \( W_2^\mu \) and nested sets of interpolation points \( T_M \), \( 1 \leq M \leq M_{\text{max}} \); we then choose \( S_N^\mu \) and solve for (and orthonormalize) the \( \zeta_n, 1 \leq n \leq N \); we finally form and store \( a_0(\zeta_j, \zeta_i), a_1(\zeta_j, \zeta_i, q_m) \), \( f_2(\zeta_i, q_m) \), and \( \ell(\zeta_i), 1 \leq i \leq N, 1 \leq m \leq M_{\text{max}} \). (Note that \( S_N^\mu \) and \( W_N^\mu \) are constructed by a greedy selection process [24,29,40] that ensures “maximally independent” basis functions and hence a rapidly convergent reduced-basis approximation.) All quantities computed in the offline stage are independent of the parameter \( \mu \); note these quantities must be computed in a stable fashion which is consistent with the finite element quadrature points (see [23] p. 173, and [12] p. 132). In the online stage – performed many times for each new \( \mu \) – we first compute \( \varphi_M(\mu) \) from (15) at cost \( O(M^2) \) by appealing to the triangular property of \( B^M \); we then assemble and invert the (full) \( N \times N \) reduced-basis stiffness matrix \( a_0(\zeta_j, \zeta_i) + \sum \varphi_M(\mu) a_1(\zeta_j, \zeta_i, q_m) \) to obtain \( u_{N,M,j}, 1 \leq j \leq N, \) at cost \( O(N^2M) \) for assembly plus \( O(N^3) \) for inversion; we finally evaluate the reduced-basis output \( s_{N,M}(\mu) \) as \( s_{N,M}(\mu) = \sum_{j=1}^N u_{N,M,j} \ell(\zeta_j) \) at cost \( O(N) \). The operation count for the online stage is thus only \( O(M^2 + N^2M + N^3) \).

Hence, as required in the many-query or real-time contexts, the online complexity is independent of \( N \), the dimension of the underlying “truth” finite element approximation space. Since \( N, M \ll N \) we expect significant
3.2.4. Numerical results

We present here numerical results for the model problem of Section 3.1.2. We first define $(w, v) = \int_\Omega \nabla w \cdot \nabla v$; thanks to the Dirichlet conditions on the boundary, $(w, v)_X$ is appropriately coercive. We note that for our particular function, $g(x; \mu) = G(x; \mu)$ of (20), $S_g^M$, $W^M_g$, and hence $T^M$ and $B^M$ are already constructed in Section 2.3. The sample set $S_N^\mu$ and associated reduced-basis space $W_N^\mu$ are developed based on the adaptive sampling procedure described in [23,40].

We now introduce a parameter sample $\Xi_{\text{Test}} \subset D$ of size 225 (in fact, a regular $15 \times 15$ grid over $D$), and define $\epsilon_{u,N,M,\text{max,rel}} = \max_{\mu \in \Xi_{\text{Test}}} \| u(\mu) - u_{N,M}(\mu) \|_X / \| u_{\text{max}} \|_X$ and $\epsilon_{s,N,M,\text{max,rel}} = \max_{\mu \in \Xi_{\text{Test}}} | s(\mu) - s_{N,M}(\mu) | / | s_{\text{max}} |$; here $\| u_{\text{max}} \|_X = \max_{\mu \in \Xi_{\text{Test}}} \| u(\mu) \|_X$ and $| s_{\text{max}} | = \max_{\mu \in \Xi_{\text{Test}}} | s(\mu) |$. We present in Figure 2 $\epsilon_{u,N,M,\text{max,rel}}$ as a function of $N$ and $M$. We observe that the reduced-basis approximation converges very rapidly. We also note, consistent with Proposition 3.1, the “plateau” in the curves for $M$ fixed and the “drops” in the $N \to \infty$ asymptotes as $M$ is increased: for fixed $M$ the error in our coefficient function approximation $g_M(x; \mu)$ to $g(x; \mu)$ will ultimately dominate for large $N$; increasing $M$ renders the coefficient function approximation more accurate, which in turn leads to the drops in the asymptotic error. Figure 2 clearly suggests (for this particular problem) the optimal “$N - M$” strategy. We tabulate in Table 3 $\epsilon_{u,N,M,\text{max,rel}}$ and $\epsilon_{s,N,M,\text{max,rel}}$ for $M$ chosen roughly optimally — but conservatively, to ensure that we are not on a “plateau” for each $N$. We observe very rapid convergence of the reduced-basis approximation with $N, M$. (Note that the convergence of the output can be further improved by the introduction of adjoint techniques [23,24,29].)

Finally, we present in Table 4 the online computational times to calculate $s_{N,M}(\mu)$ as a function of $(N, M)$; the values are normalized with respect to the computational time for the direct calculation of the truth approximation output $s(\mu) = \ell(u(\mu))$. We achieve significant computational savings: for a relative accuracy of close to 0.024 percent (corresponding to $N = 8$, $M = 20$ in Tab. 3) in the output, the online saving is more than a factor of 2000.

![Figure 2](image-url)  

**Table 3.** Maximum relative error in the energy norm and output for the nonaffine elliptic example.

<table>
<thead>
<tr>
<th>$N$</th>
<th>$M$</th>
<th>$\epsilon_{u,N,M,\text{max,rel}}$</th>
<th>$\epsilon_{s,N,M,\text{max,rel}}$</th>
</tr>
</thead>
<tbody>
<tr>
<td>4</td>
<td>15</td>
<td>1.20 E –02</td>
<td>5.96 E –03</td>
</tr>
<tr>
<td>8</td>
<td>20</td>
<td>1.14 E –03</td>
<td>2.42 E –04</td>
</tr>
<tr>
<td>12</td>
<td>25</td>
<td>2.54 E –04</td>
<td>1.76 E –04</td>
</tr>
<tr>
<td>16</td>
<td>30</td>
<td>3.82 E –05</td>
<td>7.92 E –06</td>
</tr>
</tbody>
</table>
Table 4. Online computational times (normalized with respect to the time to solve for $s(\mu)$) for the nonaffine elliptic example.

<table>
<thead>
<tr>
<th>$N$</th>
<th>$M$</th>
<th>Online time for $s_{N,M}(\mu)$</th>
<th>(Online) time for $s(\mu)$</th>
</tr>
</thead>
<tbody>
<tr>
<td>4</td>
<td>15</td>
<td>2.39 E-04</td>
<td>1</td>
</tr>
<tr>
<td>8</td>
<td>20</td>
<td>4.33 E-04</td>
<td>1</td>
</tr>
<tr>
<td>12</td>
<td>25</td>
<td>5.41 E-03</td>
<td>1</td>
</tr>
<tr>
<td>16</td>
<td>30</td>
<td>6.93 E-03</td>
<td>1</td>
</tr>
</tbody>
</table>

4. Nonaffine linear parabolic equations

We now extend the results of the previous section to parabolic problems with nonaffine parameter dependence. The essential new ingredient is the presence of time; we shall “simply” treat time as an additional, albeit special, parameter. In what follows, we introduce an abstract statement of the nonaffine parabolic equation and develop the associated reduced-basis approximation; finally, we present numerical results obtained for a model problem.

4.1. Problem formulation

4.1.1. Abstract statement

The “truth” finite element approximation is based on (5) for $Y^e \equiv L^2(\Omega)$; as in Section 3, $a$ and $f$ are of the form (21) and (22), respectively. We shall make the following assumptions. First, we assume that the bilinear form $a(: ; \mu)$ is symmetric and satisfies the coercivity and continuity conditions (23) and (24), respectively. Second, we assume that the bilinear form $m(\cdot, \cdot)$ is symmetric

$$m(v, w) = m(w, v), \quad \forall w, v \in Y^e, \forall \mu \in D;$$

and $Y^e$-coercive,

$$0 < \sigma \equiv \inf_{v \in Y^e} \frac{m(v, v)}{\|v\|_{Y^e}^2}, \quad \forall \mu \in D;$$

and $Y^e$-continuous,

$$\sup_{w \in Y^e} \sup_{v \in Y^e} \frac{m(w, v)}{\|w\|_{Y^e} \|v\|_{Y^e}} \leq \rho < \infty, \quad \forall \mu \in D.$$  \hspace{1cm} (40)

(We plausibly suppose that $\rho$ and $\sigma$ may be chosen independent of $N$.) We also require that the linear forms $f(:; \mu) : X \to \mathbb{R}$ and $\ell(:) : X \to \mathbb{R}$ are bounded with respect to $\|\cdot\|_{Y^e}$; the former is perforce satisfied for the choice (22). Third, and finally, we assume that all linear and bilinear forms are independent of time – the system is thus linear time-invariant (LTI). It follows from our hypotheses that the finite element truth solution exists and is unique (see, e.g., [30]).

We note that the output and field variable are now functions of both the parameter $\mu$ and (discrete) time $t^k$. For simplicity of exposition, we assume here that $m$ does not depend on the parameter; however, dependence on the parameter is readily admitted [13]. We also note that the method presented here easily extends to nonzero initial conditions, to multiple control inputs and outputs, and to nonsymmetric problems such as the convection-diffusion equation [12].

4.1.2. Model problem

Our particular numerical example is the unsteady analogue of the model problem introduced in Section 3.1.2: we recall that $\mu \in D \equiv [-1, -0.01]^2$, that $\Omega = [0, 1]^2$, and that our “truth” approximation subspace $X \equiv H^1_0(\Omega)$ is of dimension $N = 2601$. The governing equation for $u(\mu, t^k) \in X$ is thus (5) with $a(w, v; \mu) = a_0(w, v) + a_1(w, v, g(:; \mu))$, $f(v; \mu) = \int_\Omega v g(:; \mu)$,

$$m(w, v) \equiv \int_\Omega w v.$$  \hspace{1cm} (42)
where \( g = G \) as given in (20). The output is given by \( s(\mu, t^k) = \ell(u(\mu, t^k)), \forall k \in \mathbb{K} \), where \( \ell(v) = \int_\Omega v \). We shall consider the time interval \( t = [0, 2] \) and a timestep \( \Delta t = 0.01 \); we thus have \( K = 200 \). Finally, we assume that we are given the periodic control input \( b(t) = \sin(2\pi t) \), \( t \in \mathbb{I} \).

4.2. Reduced-basis approximation

4.2.1. Fully discrete equations

We first introduce the nested sample sets \( S_N^n = \{ \tilde{\mu}_n^1 \in \bar{D}, \ldots, \tilde{\mu}_n^N \in \bar{D} \} \), \( 1 \leq N \leq N_{\text{max}} \), where \( \tilde{\mu} \equiv (\mu, t^k) \) and \( \bar{D} \equiv D \times \mathbb{I} \); note that the samples must now reside in the parameter-time space, \( \bar{D} \). We then define the associated nested Lagrangian [28] reduced-basis space

\[
W_N^u = \text{span}\{ u(\tilde{\mu}_n^u), 1 \leq n \leq N \}, \quad 1 \leq N \leq N_{\text{max}},
\]

(43)

where \( u(\tilde{\mu}_n^u) \) is the solution of (5) at time \( t = t^k_n \) for \( \mu = \mu_n^u \). (As in the elliptic case, the \( \zeta_n \) are actually orthonormalized relative to the \( \langle \cdot, \cdot \rangle_X \) inner product.)

Our reduced-basis approximation \( u_{N,M}(\mu, t^k) \) to \( u(\mu, t^k) \) is then obtained by a standard Galerkin projection: given \( \mu \in \mathbb{D} \), \( u_{N,M}(\mu, t^k) \in W_N^u \) satisfies

\[
m(u_{N,M}(\mu, t^k), v) + \Delta t \left[ a_0(u_{N,M}(\mu, t^k), v) + a_1(u_{N,M}(\mu, t^k), v; g_M(\cdot; \mu)) \right]
= m(u_{N,M}(\mu, t^{k-1}), v) + \Delta t \int_\Omega v g_M(x; \mu) b(t^k), \quad \forall v \in W_N^u, \forall k \in \mathbb{K},
\]

(44)

with initial condition \( u_{N,M}(\mu, t^0) = 0 \); here, \( g_M(x; \mu) \) is the coefficient function approximation defined in (14). We then evaluate the output estimate, \( s_{N,M}(\mu, t^k) \), from

\[
s_{N,M}(\mu, t^k) \equiv \ell(u_{N,M}(\mu, t^k)), \quad \forall k \in \mathbb{K}.
\]

(45)

The parameter-time sample set \( S_N^n \) and associated reduced-basis space \( W_N^u \) are constructed using a “greedy” adaptive sampling procedure in footnote 5; we refer the interested reader to [13] for a detailed discussion of this procedure.

The reduced-basis subspace defined in (43) is the span of solutions of our “truth approximation” \( u(\mu, t^k) \) at the sample points \( S_N^n \). In many cases, however, the control input \( b(t^k) \) is not known in advance and thus we cannot solve for \( u(\mu, t^k) \) – as often arises in optimal control problems. Fortunately, we may appeal to the LTI hypotheses in such cases and construct the space based on the impulse response [13].

As regards the convergence rate \( u_{N,M}(\mu, t^k) \rightarrow u(\mu, t^k) \), we can develop \( \text{apriori} \) estimates very similar in form to the elliptic case – the sum of a best approximation result and a perturbation due to the variational crime associated with the interpolation of \( g \). The result is given in Proposition A.1 in the Appendix. It is also clear from Proposition A.1 that \( M \) should be chosen such that \( \epsilon_M(\mu) \) is of the same order as the error in the best approximation, otherwise the perturbation term may limit the convergence of the reduced-basis approximation. As regards the best approximation, \( W_N^u \) comprises “snapshots” on the parametrically induced manifold \( M^n \equiv \{ u(\mu, t^k) \mid \forall (\mu, t^k) \in \bar{D} \} \) which is very low-dimensional and smooth under general hypotheses on stability and continuity; the best approximation \( u_{N,M}(\mu) \) should thus converge to \( u(\mu, t^k) \) very rapidly.

The offline-online procedure for nonaffine linear parabolic equations is a straightforward combination of the procedures developed for affine parabolic equations [13] and nonaffine elliptic equations (see Sect. 3). For example, the online effort is \( O(MN^2) \) to assemble the reduced-basis discrete system, \( O(N^3 + KN^2) \) to obtain the reduced-basis coefficients at \( t^k, 0 \leq k \leq K \), and \( O(KN) \) to compute the output at \( t^k, 0 \leq k \leq K \). (Recall that our system is LTI and hence the reduced-basis matrices are time-independent.)

4.2.2. Numerical results

We now present numerical results for our model problem of Section 4.1.2. The sample set \( S^n_M \) and associated basis \( W^n_M \) – and hence \( T_M \) and \( B_M \) – for the nonaffine function approximation are constructed as in Section 2.3.
Figure 3. Convergence of the reduced-basis approximation for the non-affine parabolic example.

Table 5. Maximum relative error in the energy norm and output for different values of $N$ and $M$ for the nonaffine parabolic problem.

<table>
<thead>
<tr>
<th>N</th>
<th>M</th>
<th>$\varepsilon_{u,N,M,\max,rel}$</th>
<th>$\varepsilon_{s,N,M,\max,rel}$</th>
</tr>
</thead>
<tbody>
<tr>
<td>5</td>
<td>8</td>
<td>4.12 E–02</td>
<td>4.23 E–02</td>
</tr>
<tr>
<td>10</td>
<td>16</td>
<td>3.12 E–03</td>
<td>3.03 E–03</td>
</tr>
<tr>
<td>20</td>
<td>24</td>
<td>1.97 E–04</td>
<td>1.79 E–04</td>
</tr>
<tr>
<td>30</td>
<td>32</td>
<td>2.46 E–05</td>
<td>7.65 E–06</td>
</tr>
<tr>
<td>40</td>
<td>40</td>
<td>4.27 E–06</td>
<td>2.21 E–06</td>
</tr>
<tr>
<td>50</td>
<td>48</td>
<td>7.48 E–07</td>
<td>1.29 E–07</td>
</tr>
</tbody>
</table>

We then generate the $S_N^u$ and associated reduced-basis space $W_N^u$ following the procedure in [24]; note for parabolic problems [12], we extract our snapshots from a parameter-time sample.

In the time-dependent case we define the maximum relative error in the energy norm as $\varepsilon_{u,N,M,\max,rel} = \max_{\mu \in \Omega_{\text{Test}}} |||e(\mu, t_K)|||/|||u(\mu, t_K)|||$, and the maximum relative output error as $\varepsilon_{s,N,M,\max,rel} = \max_{\mu \in \Omega_{\text{Test}}} |s(\mu, t_{s}(\mu)) - s_{N,M}(\mu, t_{s}(\mu))|/|s(\mu, t_{s}(\mu))|$, where $\Xi_{\text{Test}} \subset \Omega$ is the parameter test sample of size 225 introduced in Section 3.2.4, $\mu_u \equiv \arg\max_{\mu \in \Omega_{\text{Test}}} |||u(\mu, t_K)|||$, $t_{s}(\mu) = \arg\max_{t_k \in \mathbb{I}} |s(\mu, t_k)|$, and the energy norm is defined as $|||v(\mu, t_k)||| = \left(\int_{\mathbb{D}} (v(\mu, t_k), v(\mu, t_k)) + \sum_{k'=1}^{K} a(v(\mu, t_{k'}), v(\mu, t_{k'}); g(\cdot; \mu)) \Delta t) \right)^{1/2}$, $\forall v \in L^\infty(\mathbb{D} \times \mathbb{I}; X)$. We plot in Figure 3 $\varepsilon_{u,N,M,\max,rel}$ as a function of $N$ and $M$. The graph shows the same behavior already observed in the elliptic case: the error levels off at smaller and smaller values as we increase $M$. In Table 5, we present $\varepsilon_{u,N,M,\max,rel}$ and $\varepsilon_{s,N,M,\max,rel}$ as a function of $N$ and $M$; note that the tabulated $(N, M)$ values correspond roughly to the optimal “knees” of the $N - M$-convergence curves. It is interesting to compare Table 3 (elliptic) and Table 5 (parabolic): as expected, for the same accuracy, the requisite $M$ is roughly the same, since $G$ is time-independent; however, $N$ is larger for the parabolic case as $u$ is a function of $\mu$ and time.

In Table 6 we present, as a function of $N$ and $M$, the online computational times to calculate $s_{N,M}(\mu, t_k)$ and $\Delta_{N,M}(\mu, t_k)$, $\forall k \in \mathbb{K}$. The values are normalized with respect to the computational time for the direct calculation of the truth approximation output $s(\mu, t_k) = \ell(u(\mu, t_k))$, $\forall k \in \mathbb{K}$. The computational saving is quite significant: for a relative accuracy of roughly 0.02 percent ($N = 20, M = 24$) in the output, the online time to compute $s_{N,M}(\mu, t_k)$ is about 1/1000 the time to directly calculate $s(\mu, t_k)$. 


Table 6. Online computational times (normalized with respect to the time to solve for $s(\mu, t^k), \forall k \in K$) for the nonaffine parabolic problem.

<table>
<thead>
<tr>
<th>N</th>
<th>M</th>
<th>Online time for $s_{N,M}(\mu, t^k), \forall k \in K$</th>
<th>(Online) time for $s(\mu, t^k), \forall k \in K$</th>
</tr>
</thead>
<tbody>
<tr>
<td>5</td>
<td>8</td>
<td>6.96 E-04</td>
<td>1</td>
</tr>
<tr>
<td>10</td>
<td>16</td>
<td>7.61 E-04</td>
<td>1</td>
</tr>
<tr>
<td>20</td>
<td>24</td>
<td>1.05 E-03</td>
<td>1</td>
</tr>
<tr>
<td>30</td>
<td>32</td>
<td>1.25 E-03</td>
<td>1</td>
</tr>
<tr>
<td>40</td>
<td>40</td>
<td>1.68 E-03</td>
<td>1</td>
</tr>
<tr>
<td>50</td>
<td>48</td>
<td>2.06 E-03</td>
<td>1</td>
</tr>
</tbody>
</table>

5. NONLINEAR MONOTONIC ELLIPTIC EQUATIONS

In this section we extend the results of Section 3 to treat nonlinear monotonic elliptic equations. We first introduce an abstract statement of the nonlinear elliptic equation and as well as a model problem. We then develop the associated reduced-basis approximation based on a coefficient-function expansion for the nonlinear term; we then discuss numerical results obtained with the model problem.

5.1. Problem formulation

5.1.1. Abstract statement

We consider the following “exact” (superscript e) problem: for any $\mu \in D \subset \mathbb{R}^P$, find $s^e(\mu) = \ell(u^e(\mu))$, where $u^e(\mu) \in X^e$ satisfies the weak form of the $\mu$-parametrized nonlinear partial differential equation

$$a^L(u^e(\mu), v) + \int_{\Omega} g(u^e(\mu); x; \mu)v = f(v), \quad \forall v \in X^e. \quad (46)$$

Here $g(u^e; x; \mu)$ is a rather general nonaffine nonlinear function of the parameter $\mu$, spatial coordinate $x$, and field variable $u^e(x; \mu)$ (we present our assumptions later); and $a^L(\cdot, \cdot)$ and $f(\cdot)$ are $X^e$-continuous bounded bilinear and linear functionals, respectively – these forms are assumed to be parameter-independent for the sake of simplicity.

Next, we recall our reference (or “truth”) finite element approximation space $X(\subset X^e)$ of dimension $N$. Our truth approximation is then: given $\mu \in D$, we find

$$s(\mu) = \ell(u(\mu)), \quad (47)$$

where $u(\mu) \in X$ is the solution of the discretized weak formulation

$$a^L(u(\mu), v) + \int_{\Omega} g(u(\mu); x; \mu)v = f(v), \quad \forall v \in X. \quad (48)$$

We assume that $\|u^e(\mu) - u(\mu)\|_X$ is suitably small and hence that $N$ will typically be very large.

We shall make the following assumptions. First, we assume that the bilinear form $a^L(\cdot, \cdot) : X \times X \to \mathbb{R}$ is symmetric, $a^L(w, v) = a^L(v, w), \forall w, v \in X$. We shall also make two crucial hypotheses related to well-posedness.
Our first hypothesis is that the bilinear form $a^L$ satisfies a stability and continuity condition

$$0 < \alpha \equiv \inf_{v \in X} \frac{a^L(v, v)}{\|v\|_X^2},$$

(49)

and that $f \in L^2(\Omega)$. For the second hypothesis we require that $g : \mathbb{R} \times \Omega \times \mathcal{D} \rightarrow \mathbb{R}$ is continuous in its arguments, increasing in its first argument, and satisfies, $\forall g \in \mathbb{R}, \ y g(y; x; \mu) \geq 0$ for any $x \in \Omega$ and $\mu \in \mathcal{D}$. With these assumptions, the problems (46) and (48) are indeed well-posed.

We can prove that there exists a solution $u^c \in X^c$ to the problem (46) first by considering the problem (46) with $g$ replaced by

$$g_n(z; x; \mu) = \begin{cases} g(z; x; \mu) & \text{if } |z| \leq n \\ -n & \text{if } z < -n \\ n & \text{if } z > n \end{cases}$$

(51)

and then taking the limit using Fatou’s lemma (see [19]). In addition, the solution is unique: suppose indeed that (46) has two solutions, $u^n_1$ and $u^n_2$; this implies

$$a^L(u^n_1 - u^n_2, v) + \int_{\Omega} (g(u_1; x; \mu) - g(u_2; x; \mu)) v = 0, \quad \forall v \in H^1_0(\Omega);$$

by choosing $v = u^n_1 - u^n_2$, we arrive at

$$a^L(u^n_1 - u^n_2, u^n_1 - u^n_2) + \int_{\Omega} (g(u_1; x; \mu) - g(u_2; x; \mu)) (u^n_1 - u^n_2) = 0, \quad \forall v \in H^1_0(\Omega);$$

it follows from the coercivity of $a^L$ and monotonicity of $g$ in its first argument that $u^n_1 = u^n_2$, and hence the solution is unique.

5.1.2. A model problem

We consider the model problem $-\nabla^2 u + \mu(1) \frac{e^{\mu(1)u}}{\mu(1)} - \mu(2) \frac{e^{\mu(2)u}}{\mu(2)} = 100 \sin(2 \pi x^{(1)}_1) \cos(2 \pi x^{(2)}_2)$, where $(x^{(1)}, x^{(2)}) \in \Omega = [0, 1]^2$ and $\mu = (\mu(1), \mu(2)) \in \mathcal{D}^\mu \equiv [0.01, 10]^2$; we impose a homogeneous Dirichlet condition on the boundary $\partial \Omega$. The output of interest is the average of the field variable over the physical domain. The weak formulation is then stated as: given $\mu \in \mathcal{D}$, find $s(\mu) = \int_{\Omega} u(\mu)$, where $u(\mu) \in X = H^1_0(\Omega) \equiv \{ v \in H^1_0(\Omega) \mid v_{|\partial \Omega} = 0 \}$ is the solution of

$$\int_{\Omega} \nabla u \cdot \nabla v + \int_{\Omega} \mu(1) \frac{e^{\mu(1)u}}{\mu(1)} v = 100 \int_{\Omega} \sin(2 \pi x^{(1)}_1) \cos(2 \pi x^{(2)}_2) v, \quad \forall v \in X.$$  

(52)

Our abstract statement (47) and (48) then obtains for

$$a^L(w, v) = \int_{\Omega} \nabla w \cdot \nabla v, \quad f(v) = 100 \int_{\Omega} \sin(2 \pi x^{(1)}_1) \cos(2 \pi x^{(2)}_2) v, \quad \ell(v) = \int_{\Omega} v,$$

and

$$g(y; x; \mu) = \mu(1) \frac{e^{\mu(1)y}}{\mu(1)} - \mu(2) \frac{e^{\mu(2)y}}{\mu(2)}.$$  

(54)

Note that $\mu(1)$ controls the strength of the sink term and $\mu(2)$ the strength of the nonlinearity. Clearly, $g$ satisfies our hypotheses.

We present in Figure 4 two typical solutions obtained with the finite element “truth” approximation space $X$ of dimension $N = 2601$. We see that when $\mu = (0.01, 0.01)$, the solution has two negative peaks and two positive peaks with similar height (this solution is very similar to that of the problem in which $g(u; \mu)$ is zero). However,
as $\mu$ increases, the negative peaks remain largely unchanged while the positive peaks are strongly rectified as shown in Figure 4b for $\mu = (10, 10)$: as $\mu$ increases the exponential function $\mu(1)e^{\mu(2)u}$ damps the positive part of $u(\mu)$, but has no effect on the negative part of $u(\mu)$.

5.2. Reduced-basis approximation

5.2.1. Discrete equations

We first motivate the need for incorporating the empirical interpolation procedure into the reduced-basis method to treat nonlinear equations. If we were to directly apply the Galerkin procedure of the linear affine case, our reduced-basis approximation would satisfy

$$a^L(u_N(\mu),v) + \int_\Omega g(u_N(\mu);x;\mu)v = f(v), \quad \forall v \in W_N^u. \quad (55)$$

Observe that if $g$ is a low order [18,38] polynomial nonlinearity of $u$, we can then develop an efficient offline-online procedure. Unfortunately, this strategy can not be applied to high-order polynomial and non-polynomial nonlinearities: the operation count for the online stage will scale as some power of $N$.

We seek an online evaluation cost that depends only on the dimension of reduced-basis approximation spaces and the parametric complexity of the problems – and not on $N$. Towards that end, we first construct nested samples $S_M = \{\mu_0, \ldots, \mu_M \in \cal{D}\}$, associated nested approximation spaces $W_M^u = \text{span}\{\xi_m = g(u(\mu_m);x;\mu_m), 1 \leq m \leq M\} = \text{span}\{q_1, \ldots, q_M\}$, and nested sets of interpolation points $T_M = \{x_1, \ldots, x_M\}$ for $1 \leq M \leq M_{\text{max}}$ following the procedure of Section 2.1. Then for any given $w \in X$ and $M$, we approximate $g(w;x;\mu)$ by $g_M^w(x;\mu) = \sum_{m=1}^M \varphi_M^m(\mu)q_m(x)$, where $\sum_{j=1}^M B_{ij}^M \varphi_M^j(\mu) = g(w(x_i);x_i;\mu)$, $1 \leq i \leq M$.

We may now approximate $g(u_{N,M};x;\mu)$ as required in our reduced-basis projection for $u_{N,M}(\mu)$ by $g_M^{u_{N,M}}(x;\mu)$. Our reduced-basis approximation is thus: Given $\mu \in \cal{D}$, we evaluate

$$s_{N,M}(\mu) = \ell(u_{N,M}(\mu)), \quad (56)$$

where $u_{N,M}(\mu) \in W_N^u$ satisfies

$$a^L(u_{N,M}(\mu),v;\mu) + \int_\Omega g_M^{u_{N,M}}(x;\mu)v = f(v), \quad \forall v \in W_N^u. \quad (57)$$

We now turn to the computational complexity.
5.2.2. Offline-online procedure

The most significant new issue is efficient calculation of the nonlinear term \( g_N^M(x; \mu) \), which we now elaborate in some detail. We first expand our reduced-basis approximation and coefficient-function approximation as

\[
u_N(M) = \sum_{j=1}^{N} u_{N,M,j}(\mu) \zeta_j, \quad g_N^M(x; \mu) = \sum_{m=1}^{M} \varphi_M(m(\mu)) q_m.
\]

(58)

Inserting these representations into (57) yields

\[
\sum_{j=1}^{N} A_{ij}^N u_{N,M,j}(\mu) + \sum_{m=1}^{M} C_{im}^{N,M} \varphi_M(m(\mu)) = F_{Ni}, \quad 1 \leq i \leq N;
\]

(59)

where \( A^N \in \mathbb{R}^{N \times N}, C^{N,M} \in \mathbb{R}^{N \times M}, F_N \in \mathbb{R}^N \) are given by \( A_{ij}^N = a_i^L(\zeta_j, \zeta_i), 1 \leq i, j \leq N, \) \( C_{im}^{N,M} = \int_{\Omega} q_m(\zeta_i), 1 \leq i \leq N, 1 \leq m \leq M, \) and \( F_{Ni} = f(\zeta_i), 1 \leq i \leq N, \) respectively. Furthermore, \( \varphi_M(\mu) \in \mathbb{R}^M \) is given by

\[
\sum_{k=1}^{M} D_{mk}^N \varphi_M(k(\mu)) = g(u_{N,M}(x_m; \mu); x_m; \mu), \quad 1 \leq m \leq M
\]

\[
= g\left( \sum_{n=1}^{N} u_{N,M,n}(\mu) \zeta_n(x_m), x_m; \mu \right), \quad 1 \leq m \leq M.
\]

(60)

We then substitute \( \varphi_M(\mu) \) from (60) into (59) to obtain the following nonlinear algebraic system

\[
\sum_{j=1}^{N} A_{ij}^N u_{N,M,j}(\mu) + \sum_{m=1}^{M} D_{im}^N g\left( \sum_{n=1}^{N} \zeta_n(x_m) u_{N,M,n}(\mu); x_m; \mu \right) = F_{Ni}, \quad 1 \leq i \leq N,
\]

(61)

where \( D^{N,M} = C^{N,M}(B^M)^{-1} \in \mathbb{R}^{N \times M}. \)

To solve (61) for \( u_{N,M,j}(\mu), 1 \leq j \leq N, \) we may apply a Newton iterative scheme: given a current iterate \( \bar{u}_{N,M,j}(\mu), 1 \leq j \leq N, \) we find an increment \( \delta u_{N,M,j}(\mu), 1 \leq j \leq N, \) such that

\[
\sum_{j=1}^{N} (A_{ij}^N + \bar{E}_{ij}^N) \delta u_{N,M,j}(\mu) = F_{Ni} - \sum_{j=1}^{N} A_{ij}^N \bar{u}_{N,M,j}(\mu)
\]

\[
- \sum_{m=1}^{M} D_{im}^N g\left( \sum_{n=1}^{N} \zeta_n(x_m) \bar{u}_{N,M,n}(\mu); x_m; \mu \right), \quad 1 \leq i \leq N;
\]

(62)

here \( \bar{E}^N \in \mathbb{R}^{N \times N} \) must be calculated at every Newton iteration as

\[
\bar{E}_{ij}^N = \sum_{m=1}^{M} D_{im}^N g_1\left( \sum_{n=1}^{N} \zeta_n(x_m) \bar{u}_{N,M,n}(\mu); x_m; \mu \right) \zeta_j(x_m), \quad 1 \leq i, j \leq N;
\]

(63)

where \( g_1 \) is the partial derivative of \( g \) with respect to its first argument. Finally, the output can be evaluated as

\[
s_{N,M}(\mu) = \sum_{j=1}^{N} u_{N,M,j}(\mu)L_{N,j},
\]

(64)
where $L_N \in \mathbb{R}^N$ is the output vector with entries $L_{Nj} = \ell(j), 1 \leq j \leq N$. Based on this strategy, we can develop an efficient offline-online procedure for the rapid evaluation of $s_{N,M}(\mu)$ for each $\mu$ in $\mathcal{D}$.

The operation count of the online stage is essentially the predominant Newton update component (62): at each Newton iteration, we first assemble the right-hand side and compute $\bar{E}_N$ of (63) at cost $O(MN^2)$ – note we perform the sum in the parenthesis of (63) before performing the outer sum; we then form and invert the left-hand side (Jacobian) of (62) at cost $O(N^3)$. The online complexity depends only on $N$, $M$, and the number of Newton iterations; we thus recover $N$ independence of the online stage.

5.2.3. Numerical results

We first define $(w,v)_X = \int_\Omega \nabla w \cdot \nabla v$ and thus obtain $\alpha = 1$. We next construct $S^g_M$ and $W^g_M$ with the $L^2(\Omega)$-norm surrogate approach on $\Xi^g$, where $\Xi^g$ is a regular $12 \times 12$ grid over $\mathcal{D}$. We then generate the sample set $S^u_N$ and associated reduced-basis space $W^u_N$ using the adaptive sampling construction [23] over the grid $\Xi^g$ – but note for this nonlinear problem, our selection process is based directly on the energy norm of the true error (not an error estimate), $e(\mu) = u(\mu) - u_{N,M}(\mu)$, since the “truth” solutions $u(\mu)$ must be computed and stored for $\mu \in \Xi^g$ as part of the empirical interpolation procedure.

We now introduce a parameter test sample $\Xi_{\text{Test}}$ of size 225 (a regular $15 \times 15$ grid) and define $\epsilon_{N,M,\text{max,rel}}^u = \max_{\mu \in \Xi_{\text{Test}}} \|e_{N,M}(\mu)\|_X / \|u_{\text{max}}\|_X$ and $\epsilon_{N,M,\text{max,rel}}^s = \max_{\mu \in \Xi_{\text{Test}}} |s(\mu) - s_{N,M}(\mu)| / |s_{\text{max}}|$, where $\|u_{\text{max}}\|_X = \max_{\mu \in \Xi_{\text{Test}}} \|u(\mu)\|_X$ and $|s_{\text{max}}| = \max_{\mu \in \Xi_{\text{Test}}} |s(\mu)|$; note that $\Xi_{\text{Test}}$ is larger than (and mostly non-coincident with) $\Xi^g$. We present in Figure 5 $\epsilon_{N,M,\text{max,rel}}^u$ as a function of $N$ and $M$. We observe very rapid convergence of the reduced-basis approximation. Furthermore, the errors behave very similarly as in the linear example: the errors initially decrease, but then “plateau” in $N$ for a particular value of $M$; increasing $M$ effectively brings the error curves down. We also tabulate in Table 7 $\epsilon_{N,M,\text{max,rel}}^u$ and $\epsilon_{N,M,\text{max,rel}}^s$ for values of $(N,M)$ close to the “knees” of the convergence curves of Figure 5. We see that $s_{N,M}(\mu)$ converges very rapidly.

We present in Table 8 the online computational times to calculate $s_{N,M}(\mu)$ as a function of $(N,M)$. The values are normalized with respect to the computational time for the direct calculation of the truth approximation output $s(\mu) = \ell(u(\mu))$. The computational savings are much larger in the nonlinear case: for an relative accuracy of 0.0126 percent ($N = 12$, $M = 15$) in the output, the reduction in online cost is more than a factor of 3000; this is mainly because the matrix assembly of the nonlinear terms for the truth approximation is computationally very expensive. However we must also recall that, in the nonlinear case, the reduced-basis offline computations are much more extensive since we must solve the truth approximation over the large sample $\Xi^g$ when constructing $S^g_M$.
6. NONLINEAR PARABOLIC EQUATIONS

We now extend the results of the previous section to the time-dependent case and consider nonlinear parabolic problems. We briefly describe the abstract formulation of nonlinear parabolic problems and then develop the associated reduced-basis approximation. Finally, we discuss numerical results obtained for a model problem.

6.1. Problem formulation

6.1.1. Abstract statement

Our abstract statement is based on the nonlinear elliptic problem (48) discussed in Section 5: Given a parameter \( \mu \in \mathcal{D} \), we evaluate the output of interest

\[
s(\mu, t^k) = \ell(u(\mu, t^k)), \quad \forall k \in \mathcal{K}
\]

where the field variable \( u(\mu, t^k) \in X, \forall k \in \mathcal{K} \), satisfies the weak form of the nonlinear parabolic partial differential equation

\[
m(u(\mu, t^k), v) + \Delta t \ a^k(u(\mu, t^k), v) + \Delta t \int_{\Omega} g(u(\mu, t^k); x; \mu) v
\]

\[
= m(u(\mu, t^{k-1}), v) + \Delta t \ f(v) b(t^k), \quad \forall v \in X, \forall k \in \mathcal{K},
\]

with initial condition (say) \( u(\mu, t^0) = 0 \). We note that if an explicit scheme such as Euler-Forward is used, we then arrive at a linear system for \( u(\mu, t^k) \) but now burdened with a conditional stability restriction on \( \Delta t \). In that case, the discrete reduced-basis system is inherently linear.

As in the linear parabolic case and nonlinear elliptic case, the same assumptions are applied to \( m, a^k, g, f, \) and \( \ell \), it is a classical result of nonlinear analyses (truncation and monotonicity) to prove well-posedness of this problem (see [19]).
6.1.2. Model problem

Our particular numerical example is the unsteady analogue of the elliptic model problem introduced in Section 5.1.2: we have $\mu=(\mu(1),\mu(2)) \in D^\mu \equiv [0.01,10]^2$, the spatial domain is the unit square, $\Omega=[0,1]^2$, and our “truth” approximation finite element space $X=H^1_0(\Omega)$ has dimension $N=2601$. The field variable $u(\mu,t^k) \in X$ thus satisfies

$$
m(v,w) \equiv \int_\Omega v w, \quad a^L(v,w) \equiv \int_\Omega \nabla v \cdot \nabla w, \quad f(v) \equiv 100 \int_\Omega v \sin(2\pi x(1)) \cos(2\pi x(2)),
$$

and

$$
g(u(\mu,t^k);\mu) = \mu(1) \frac{e^{\mu(2)} u(\mu,t^k) - 1}{\mu(2)}. \quad (68)
$$

The output $s(\mu,t^k)$ is evaluated from (65) with $\ell(v) = \int_\Omega v$. We shall consider the time interval $I = [0,2]$ and a timestep $\Delta t = 0.01$; we thus have $K=200$. The control input is given by $b(t^k) = \sin(2\pi t^k)$, $t \in \mathbb{I}$.

6.2. Reduced-basis approximation

6.2.1. Fully discrete equations

We first introduce the nested sample sets $S^\mu_M = \{\tilde{\mu}^1_M \in \tilde{D}, \ldots, \tilde{\mu}^N_M \in \tilde{D}\}$, $1 \leq M \leq M_{\text{max}}$ and $S^\mu_N = \{\tilde{\mu}^1_N \in \tilde{D}, \ldots, \tilde{\mu}^N_N \in \tilde{D}\}$, $1 \leq N \leq N_{\text{max}}$, where $\tilde{\mu} \equiv (\mu,t^k)$ and $\tilde{D} \equiv D \times \mathbb{I}$. Note that, since $g(\cdot;\mu)$ is a function of the field variable $u(\mu,t^k)$, the sample set $S^\mu_M$ must now also reside in parameter-time space $\tilde{D}$; in general, $S^\mu_N \neq S^\mu_M$ and in fact $N \neq M$. We define the nested collateral reduced-basis space

$$
W^\mu_M = \{u(\mu,t^k) \in X: u(\mu,t^k) = \sum_{n=1}^{M} q_n(x) \varphi_n(\mu,t^k), \quad \varphi_n \in W^\mu_N, \quad \mu \in \tilde{D}\}
$$

and nested set of interpolation points $T_M = \{x_1,\ldots,x_M\}$, $1 \leq M \leq M_{\text{max}}$; here $u(\mu,t^k)$ is the solution of (66) at time $t = t^k$ for $\mu = \mu^0$. Next, we define the associated nested Lagrangian [28] reduced-basis space

$$
W^\mu_N = \{u(\mu,t^k) \in X: u(\mu,t^k) = \sum_{n=1}^{N} \varphi_n(\mu,t^k), \quad \varphi_n \in W^\mu_M, \quad \mu \in \tilde{D}\}
$$

and nested set of interpolation points $T_M = \{x_1,\ldots,x_M\}$, $1 \leq M \leq M_{\text{max}}$ (70)

where $u(\mu,t^k)$ is the solution of (66) at time $t = t^k$ for $\mu = \mu^0$.

Our reduced-basis approximation $u_{N,M}(\mu,t^k)$ to $u(\mu,t^k)$ is then given by: Given $\mu \in D$, $u_{N,M}(\mu,t^k) \in W^\mu_N$ satisfies

$$
m(u_{N,M}(\mu,t^k),v) + \Delta t \ a^L(u_{N,M}(\mu,t^k),v) + \Delta t \ \int_\Omega g_{M}^{u_{N,M}}(x;\mu,t^k) v = m(u_{N,M}(\mu,t^{k-1}),v) + \Delta t \ f(v) b(t^k), \quad \forall v \in W^\mu_N, \forall k \in \mathbb{K},
$$

with initial condition $u_{N,M}(\mu,t^0) = 0$; here, $g_{M}^{u_{N,M}}(x;\mu,t^k)$ is the approximation to $g(u_{N,M}(\mu,t^k);x;\mu)$ given by

$$
g_{M}^{u_{N,M}}(x;\mu,t^k) = \sum_{m=1}^{M} \varphi_{Mm}(\mu,t^k) q_m(x),
$$

where the coefficients $\varphi_{Mm}(\mu,t^k)$ are determined from

$$
\sum_{j=1}^{M} B_{ij}^{M} \varphi_{Mj}(\mu,t^k) = g(u_{N,M}(x;\mu,t^k);x;\mu), \quad 1 \leq i \leq M,
$$

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and $B_{ij}^M = q_j(x_i), 1 \leq i, j \leq M$. Finally, we evaluate the output from

$$s_{N,M}(\mu, t^k) = \ell(u_{N,M}(\mu, t^k)), \quad \forall k \in K.$$ (74)

(Note that, contrary to the previous sections, $\varphi_M(\mu, t^k)$ now also depends on time.)

At this point we should remark that our current approach of constructing the sample set $S_N^g$ and associated reduced-basis space $W_M^g$ in the nonlinear parabolic case is computationally very expensive. The reason, related to our greedy adaptive sampling procedure proposed in Section 2.1, is twofold. First, we need to calculate and store the “truth” solution $u(\mu, t^k)$ at all times $t^k \in I$ on the grid $\Xi^g$ in parameter space. In our numerical example $\Xi^g$ is of size 144 – we thus need to solve (66) 144 times and store $144 \times 200$ “truth” solutions $u(\mu, t^k)$. Second, as pointed out in Section 2.1, to determine the next sample point $\hat{\mu}^k$ in $\Xi^g \equiv \Xi^g \times I$, requires the solution of a linear program for all $\mu \in \Xi^g$ if $g$ is time-varying – as is inherently the case in the nonlinear context. (It should be noted that in the linear nonaffine parabolic case the function $g$ depends only on $x$ and $\mu$ and not on time.) Since this computation is too expensive in our current implementation, we revert to the least squares surrogate in this section – in choosing this approach we in fact rely on our numerical comparison in Section 2.1 showing that we can expect similar results.

6.2.2. Offline-online procedure

The offline-online decomposition follows directly from the corresponding procedures for linear nonaffine parabolic problems (see Sect. 4) and nonlinear elliptic problems (see Sect. 5). In summary, the operation count (per Newton iteration per timestep) in the online stage is $O(N^2 + N^3)$; the system is of course no longer “LTI”.

We remark that, in actual practice, $M$ can be quite large – and in fact much larger than $N$. We can reduce $M$ without sacrificing accuracy by splitting the time interval $I$ into several smaller subintervals $I_1, \ldots, I_I$ such that $I = \bigcup_{i=1}^I I_i$. We then construct, in the offline stage, $I$ separate samples sets $S_M^g$, $1 \leq i \leq I$, and associated reduced-basis spaces $W_{M,i}^g$, $1 \leq i \leq I$, on each interval $I_i$, $1 \leq i \leq I$. In the online stage we simply “switch” to the corresponding sample – and hence $T_M, B_M^g$, and $D^{N,M}$ as time progresses. This approach renders the offline computation more expensive (and online storage more extensive), but can increase the online efficiency considerably while retaining the desired accuracy.

6.2.3. Numerical results

We now present numerical results for our model problem of Section 6.1.2. We construct $S_M^g$ and hence $W_M^g$ with the surrogate least squares approach on $\Xi^g = \Xi \times I$, where $\Xi^g$ is a regular $12 \times 12$ grid over $D$. We generate the sample set $S_N^g$ and associated reduced-basis space $W^g$ using an adaptive sampling procedure – but note for this nonlinear parabolic problem, our selection process is based directly on the energy norm of the true error (not an error estimate), $e(\mu, t^k) = u(\mu) - u_{N,M}(\mu, t^k)$, since the “truth” solutions $u(\mu, t^k)$ are stored for $\mu \in \Xi^g$.

We now define the maximum relative error in the energy norm $\epsilon_{N,M,\max,rel}^g = \max_{\mu \in \Xi_{test}} ||e(\mu, t^K)||/ ||u(\mu, t^K)||$ and the maximum relative output error $\epsilon_{N,M,\max,rel}^g = \max_{\mu \in \Xi_{test}} |s(\mu, t_s(\mu)) - s_{N,M}(\mu, t_s(\mu))/ ||s(\mu, t_s(\mu))|||$. Here $\Xi_{test} \subset D$ is the parameter test sample of size 225 introduced in Section 5.2.3, $\mu_u \equiv \arg\max_{\mu \in \Xi_{test}} ||u(\mu, t^K)||$, $t_s(\mu) = \arg\max_{t \in I} |s(\mu, t^K)|$, and the energy norm is defined as $||v(\mu, t^K)|| = (m(v(\mu, t^K), v(\mu, t^K) + \sum_{k=1}^{K} a^{L}(v(\mu, t^K), v(\mu, t^K)) \Delta t)^{1/2}$, $\forall v \in L^\infty(D \times I; X)$.

We plot in Figure 6 $\epsilon_{N,M,\max,rel}^g$ as a function of $N$ for different values of $M$. We observe the same behavior as in the nonlinear elliptic case. We note, however, that $M$ is now much larger compared to the nonlinear elliptic model problem due to the time dependence; we recall that in the linear nonaffine elliptic and parabolic cases the required $M$ was the same since the nonaffine coefficient function did not depend on time. In Table 9, we present $\epsilon_{N,M,\max,rel}^g$ and $\epsilon_{N,M,\max,rel}^g$ as a function of $N$ and $M$. We observe very rapid convergence of the reduced-basis approximation; for $N = 20$ and $M = 80$ the error in the output is less than one percent.
Table 9. Relative error in the energy norm and output for the nonlinear parabolic problem.

<table>
<thead>
<tr>
<th>$N$</th>
<th>$M$</th>
<th>$\epsilon^u_{N,M,\max,\text{rel}}$</th>
<th>$\epsilon^s_{N,M,\max,\text{rel}}$</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>10</td>
<td>3.82E-01</td>
<td>1.00E-00</td>
</tr>
<tr>
<td>5</td>
<td>30</td>
<td>1.36E-02</td>
<td>1.91E-02</td>
</tr>
<tr>
<td>10</td>
<td>50</td>
<td>1.62E-03</td>
<td>1.46E-04</td>
</tr>
<tr>
<td>20</td>
<td>80</td>
<td>1.46E-04</td>
<td>1.67E-05</td>
</tr>
<tr>
<td>30</td>
<td>110</td>
<td>1.88E-05</td>
<td>5.16E-06</td>
</tr>
<tr>
<td>40</td>
<td>140</td>
<td>4.94E-06</td>
<td>1.56E-06</td>
</tr>
</tbody>
</table>

Table 10. Online computational times (normalized with respect to the time to solve for $s(\mu, t^k)$, $\forall k \in K$) for the nonlinear parabolic problem.

<table>
<thead>
<tr>
<th>$N$</th>
<th>$M$</th>
<th>Online time for $s_{N,M}(\mu, t^k)$, $\forall k \in K$</th>
<th>(Online) time for $s(\mu, t^k)$, $\forall k \in K$</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>10</td>
<td>6.62E-05</td>
<td>1</td>
</tr>
<tr>
<td>5</td>
<td>30</td>
<td>1.19E-04</td>
<td>1</td>
</tr>
<tr>
<td>10</td>
<td>50</td>
<td>1.74E-04</td>
<td>1</td>
</tr>
<tr>
<td>20</td>
<td>80</td>
<td>3.88E-04</td>
<td>1</td>
</tr>
<tr>
<td>30</td>
<td>110</td>
<td>7.20E-04</td>
<td>1</td>
</tr>
<tr>
<td>40</td>
<td>140</td>
<td>1.22E-03</td>
<td>1</td>
</tr>
</tbody>
</table>

In Table 10 we present, as a function of $N$ and $M$, the online computational times to calculate $s_{N,M}(\mu, t^k)$ and $\Delta^s_{N,M}(\mu, t^k)$, $\forall k \in K$. The values are normalized with respect to the computational time for the direct calculation of the truth approximation output $s(\mu, t^k) = \ell(u(\mu, t^k))$, $\forall k \in K$. The reduction in online response time is considerable. We again caution that the offline computations necessary in the nonlinear case are very extensive – primarily due to the sampling procedure for $S^o_{N,M}$. However, if a many-query context, or a clear demand for real-time response, can justify the offline cost, the reduced-basis methods can be very gainfully employed.
We consider here the rate at which $u_{N,M}(\mu, t^k)$ converges to $u(\mu, t^k)$ for the nonaffine linear parabolic case. As for the elliptic case, the interpolation-induced error will be measured through the functions $\phi_1(\mu), \phi_2(\mu)$ and $\phi_3(\mu)$ of (33)–(35) together with a comparison with respect to some best fit of $u(\mu, \cdot)$ by elements of $W_N^\mu$. The natural measure for the best fit is the “$m + \Delta t a$” norm. We thus introduce the projector $\pi_N$ defined by

$$m(v - \pi_N(v), w_N) + \Delta t a(v - \pi_N(v), w_N) = 0, \quad \pi_N(v) \in W_N^\mu, \forall w_N \in W_N^\mu, \forall v \in X. \quad (A.1)$$

We can then prove

**Proposition 6.1.** For $\varepsilon_M(\mu)$ of (16) satisfying $\varepsilon_M(\mu) < \alpha(\mu)/(4 \phi_2(\mu))$ (say), the error $e(\mu, t_k) \equiv u(\mu, t_k) - u_{N,M}(\mu, t_k)$ satisfies

$$\begin{align*}
\sigma \|e(\mu, t^k)\|_V^2 + \frac{\alpha(\mu)}{2} \Delta t \sum_{k'=1}^k \|e(\mu, t^{k'})\|_X^2 &\leq \Upsilon(\mu) \Delta t \sum_{k'=1}^k b(t^{k'})^2 \\
&+ 8\rho \|u(\mu, t^k) - \pi_N[u(\mu, t^k)]\|_V^2 + (8\gamma(\mu) + 4\alpha(\mu)) \Delta t \sum_{k'=1}^k \|u(\mu, t^{k'}) - \pi_N[u(\mu, t^{k'})]\|_X^2,
\end{align*} \quad (A.2)$$

where

$$\Upsilon(\mu) = \frac{18}{\alpha(\mu)} \varepsilon_M(\mu)^2 \left( \phi_1(\mu)^2 + \phi_2(\mu)^2 \frac{2\phi_3(\mu)^2}{\alpha(\mu)^2} \right),$$

and $\sigma$ and $\rho$ are given by (40) and (41), respectively.

**Proof.** To begin, we note from (5) and (44) that

$$m(e(\mu, t^k) - e(\mu, t^{k-1}), v) + \Delta t a(e(\mu, t^k), v; \mu)$$

$$= \Delta t \left( \int_\Omega \left[ v(g(x; \mu) - g_M(x; \mu)) \frac{b(t^k)}{t^k} + a_1(\mu, t^k) v; g(\cdot; \mu) - g_M(\cdot; \mu) \right] \right), \forall v \in W_N^\mu \quad (A.3)$$

with initial condition $e(\mu, t^0) = 0$, since $u(\mu, t^0) = u_{N,M}(\mu, t^0) = 0$ by assumption. It then follows that

$$m(e(\mu, t^k), v) + \Delta t a(e(\mu, t^k), v; \mu) = m(e(\mu, t^{k-1}), v) + \Delta t a(e(\mu, t^{k-1}), v; \mu) - \Delta t a(e(\mu, t^{k-1}), v; \mu)$$

$$+ \Delta t \left( \int_\Omega \left[ v(g(x; \mu) - g_M(x; \mu)) b(t^k) + a_1(\mu, t^k) v; g(\cdot; \mu) - g_M(\cdot; \mu) \right] \right), \forall v \in W_N^\mu. \quad (A.4)$$

Let us set now $w_N(\mu, t^k) = \pi_N[u(\mu, t^k)]$ in (A.1) and choose $v = e_N(\mu, t^k) \equiv w_N(\mu, t^k) - u_{N,M}(\mu, t^k)$ in (A.4). We then combine (A.1) and (A.4) to obtain

$$m(e_N(\mu, t^k), e_N(\mu, t^k)) + \Delta t a(e_N(\mu, t^k), e_N(\mu, t^k); \mu)$$

$$= m(e_N(\mu, t^{k-1}), e_N(\mu, t^k)) + \Delta t a(e_N(\mu, t^{k-1}), e_N(\mu, t^k); \mu) - \Delta t a(e_N(\mu, t^{k-1}), e_N(\mu, t^k); \mu)$$

$$+ \Delta t \left( \int_\Omega \left[ v(g(x; \mu) - g_M(x; \mu)) b(t^k) + a_1(\mu, t^k) v; g(\cdot; \mu) - g_M(\cdot; \mu) \right] \right)$$

$$= m(e_N(\mu, t^{k-1}), e_N(\mu, t^k)) - \Delta t a(u(\mu, t^{k-1}) - w_N(\mu, t^{k-1}), e_N(\mu, t^k); \mu)$$

$$+ \Delta t \left( \int_\Omega \left[ v(g(x; \mu) - g_M(x; \mu)) b(t^k) + a_1(u_N,M(\mu, t^k), v; g(\cdot; \mu) - g_M(\cdot; \mu)) \right] \right). \quad (A.5)$$
Multiplying both sides by 2 and then adding and subtracting appropriate terms, we arrive at

\[
m(e_N(\mu, t^k), e_N(\mu, t^k)) - m(e_N(\mu, t^{k-1}), e_N(\mu, t^{k-1})) + \Delta t a(e_N(\mu, t^k), e_N(\mu, t^k); \mu)
\]

\[
= -m(e_N(\mu, t^k) - e_N(\mu, t^{k-1}), e_N(\mu, t^k) - e_N(\mu, t^{k-1}))
\]

\[
- \Delta t a(u(\mu, t^{k-1}) - w_N(\mu, t^{k-1}) + e_N(\mu, t^k), u(\mu, t^{k-1}) - w_N(\mu, t^{k-1}) + e_N(\mu, t^k); \mu)
\]

\[
+ \Delta t a(u(\mu, t^{k-1}) - w_N(\mu, t^{k-1}), u(\mu, t^{k-1}) - w_N(\mu, t^{k-1}); \mu)
\]

\[
+ 2\Delta t \left( \int_{\Omega} v(g(x; \mu) - g_M(x; \mu)) b(t^k) + a_1(u_{N,M}(\mu, t^k), v; g(\cdot; \mu) - g_M(\cdot; \mu)) \right). \quad (A.6)
\]

It thus follows that

\[
m(e_N(\mu, t^k), e_N(\mu, t^k)) - m(e_N(\mu, t^{k-1}), e_N(\mu, t^{k-1})) + \Delta t a(e_N(\mu, t^k), e_N(\mu, t^k); \mu)
\]

\[
\leq \Delta t a(u(\mu, t^{k-1}) - w_N(\mu, t^{k-1}), u(\mu, t^{k-1}) - w_N(\mu, t^{k-1}); \mu)
\]

\[
+ 2\Delta t \left( \int_{\Omega} v(g(x; \mu) - g_M(x; \mu)) b(t^k) + a_1(u_{N,M}(\mu, t^k), v; g(\cdot; \mu) - g_M(\cdot; \mu)) \right). \quad (A.7)
\]

which after summing from \(k' = 1\) to \(k\) leads to

\[
m(e_N(\mu, t^k), e_N(\mu, t^k)) + \Delta t \sum_{k'=1}^{k} a(e_N(\mu, t^{k'}), e_N(\mu, t^{k'}); \mu)
\]

\[
\leq \Delta t \sum_{k'=1}^{k-1} a(u(\mu, t^{k'}) - w_N(\mu, t^{k'}), u(\mu, t^{k'}) - w_N(\mu, t^{k'}); \mu)
\]

\[
+ 2\Delta t \sum_{k'=1}^{k} \left( \phi_1(\mu) |b(t^{k'})| + \phi_2(\mu) \|u_{N,M}(\mu, t^{k'})\|_X \right) \varepsilon_M(\mu) \|e_N(\mu, t^{k'})\|_X. \quad (A.8)
\]

where the last inequality follows from (33) and (34). We take the square root of what we have obtained

\[
\left\{ m(e_N(\mu, t^k), e_N(\mu, t^k)) + \Delta t \sum_{k'=1}^{k} a(e_N(\mu, t^{k'}), e_N(\mu, t^{k'}); \mu) \right\}^{1/2}
\]

\[
\leq \left\{ \Delta t \sum_{k'=1}^{k-1} a(u(\mu, t^{k'}) - w_N(\mu, t^{k'}), u(\mu, t^{k'}) - w_N(\mu, t^{k'}); \mu) \right\}^{1/2}
\]

\[
+ \left\{ 2\Delta t \sum_{k'=1}^{k} \left( \phi_1(\mu) |b(t^{k'})| + \phi_2(\mu) \|u_{N,M}(\mu, t^{k'})\|_X \right) \varepsilon_M(\mu) \|e_N(\mu, t^{k'})\|_X \right\}^{1/2}, \quad (A.9)
\]
so a triangular inequality and (A.1) together give

\[
\left\{ m(u(\mu, t^k) - u_{N,M}(\mu, t^k), u(\mu, t^k) - u_{N,M}(\mu, t^k)) \right\}^{1/2}
\]

\[
+ \Delta t \sum_{k'=1}^{k} a(u(\mu, t^{k'}), u_{N,M}(\mu, t^{k'}))
\]

\[
\leq 2 \left\{ m(u(\mu, t^k) - w_N(\mu, t^k), u(\mu, t^k) - w_N(\mu, t^k)) \right\}^{1/2}
\]

\[
+ \Delta t \sum_{k'=1}^{k} a(u(\mu, t^{k'}), w_N(\mu, t^{k'}))
\]

\[
+ \left\{ 2 \Delta t \sum_{k'=1}^{k} \left( \phi_1(\mu) |b(t^{k'})| + \phi_2(\mu) \|u_{N,M}(\mu, t^{k'})\|_X \right) \varepsilon_M(\mu) \|e_{N,M}(\mu, t^k)\|_X \right\}^{1/2}
\]

We now note that \(\|e_{N,M}(\mu, t^k)\|_X \leq \|u(\mu, t^k) - w_N(\mu, t^k)\|_X + \|e(\mu, t^k)\|_X\) and recall the Young inequality (for \(c \in \mathbb{R}, d \in \mathbb{R}, g \in \mathbb{R}^+\))

\[
2 |c| |d| \leq \frac{1}{g^2} c^2 + g^2 d^2,
\]

which we apply four times: first, with \(c = \varepsilon_M(\mu) \phi_1(\mu) |b(t^k)|, d = \|u(\mu, t^k) - w_N(\mu, t^k)\|_X\), and \(g^2 = \alpha(\mu)\); second, with \(c = \varepsilon_M(\mu) \phi_1(\mu) |b(t^k)|, d = \|e(\mu, t^k)\|_X\), and \(g^2 = \alpha(\mu)/8\); third, with \(c = \varepsilon_M(\mu) \phi_2(\mu) \|u_{N,M}(\mu, t^k)\|_X\), \(d = \|u(\mu, t^k) - w_N(\mu, t^k)\|_X\), and \(g^2 = \alpha(\mu)\); and fourth, with \(c = \varepsilon_M(\mu) \phi_2(\mu) \|u_{N,M}(\mu, t^k)\|_X, d = \|e(\mu, t^k)\|_X\), and \(g^2 = \alpha(\mu)/8\). We can then bound the last term of (A.8) by

\[
2 \Delta t \sum_{k'=1}^{k} \left( \phi_1(\mu) |b(t^{k'})| + \phi_2(\mu) \|u_{N,M}(\mu, t^{k'})\|_X \right) \varepsilon_M(\mu) \|e_{N,M}(\mu, t^k)\|_X
\]

\[
\leq \varepsilon_M(\mu)^2 \frac{9}{\alpha(\mu)} \left( \phi_1(\mu)^2 \Delta t \sum_{k'=1}^{k} b(t^{k'})^2 + \phi_2(\mu)^2 \Delta t \sum_{k'=1}^{k} \|u_{N,M}(\mu, t^{k'})\|_X^2 \right)
\]

\[
+ 2 \Delta t \alpha(\mu) \sum_{k'=1}^{k} \|u(\mu, t^{k'}) - w_N(\mu, t^{k'})\|_X^2 + \Delta t \frac{\alpha(\mu)}{4} \sum_{k'=1}^{k} \|e(\mu, t^{k'})\|_X^2.
\]

We next use \(v = u_{N,M}(\mu, t^k)\) in (44), invoke the Cauchy-Schwarz inequality for \(m(u_{N,M}(\mu, t^k), u_{N,M}(\mu, t^k))\) and apply (A.11) with \(c = m^{1/2}(u_{N,M}(\mu, t^k), u_{N,M}(\mu, t^k)), d = m^{1/2}(u_{N,M}(\mu, t^{k-1}), u_{N,M}(\mu, t^{k-1}))\), and \(g = 1\),
to get
\[
m(u_{N,M}(\mu, t^k), u_{N,M}(\mu, t^k)) - m(u_{N,M}(\mu, t^{k-1}), u_{N,M}(\mu, t^{k-1})) \\
+ 2 \Delta t a(u_{N,M}(\mu, t^k), u_{N,M}(\mu, t^k); \mu) \\
\leq 2 \Delta t \int_\Omega (u_{N,M}(\mu, t^k)g_M(x; \mu)) b(t^k) \\
+ 2 \Delta t a_1(u_{N,M}(\mu, t^k), u_{N,M}(\mu, t^k); g(x; \mu) - g_M(x; \mu)) \\
\leq 2 \Delta t \phi_3(\mu) \|u_{N,M}(\mu, t^k)\|_X |b(t^k)| + 2 \Delta t \varepsilon_M(\mu) \phi_2(\mu) \|u_{N,M}(\mu, t^k)\|_X^2 \\
\leq \frac{\Delta t}{\alpha(\mu) - 2 \phi_2(\mu) \varepsilon_M(\mu)} \phi_3(\mu)^2 b(t^k)^2 + \Delta t \alpha(\mu) \|u_{N,M}(\mu, t^k)\|_X^2, \\ 
\text{(A.13)}
\]

where the second inequality follows from (34) and (35), and the last inequality from (A.11) with \( c = \phi_3(\mu) b(t^k) \), \( d = \|u_{N,M}(\mu, t^k)\|_X \), and \( \varphi = \alpha(\mu) - 2 \phi_2(\mu) \varepsilon_M(\mu) \); note that \( \varphi > 0 \) from our assumption on \( \varepsilon_M(\mu) \). Invoking (23) and summing (A.13) from \( k' = 1 \) to \( k \) we obtain
\[
m(u_{N,M}(\mu, t^k), u_{N,M}(\mu, t^k)) + \Delta t \sum_{k'=1}^k a(u_{N,M}(\mu, t^{k'}), u_{N,M}(\mu, t^{k'}); \mu) \\
\leq \frac{\phi_3(\mu)^2}{\alpha(\mu) - 2 \phi_2(\mu) \varepsilon_M(\mu)} \Delta t \sum_{k'=1}^k b(t^{k'})^2. \\ 
\text{(A.14)}
\]

From the coercivity of \( m \) and \( a \), and our assumption on \( \varepsilon_M(\mu) \) it then directly follows that
\[
\Delta t \sum_{k'=1}^k \|u_{N,M}(\mu, t^{k'})\|_X^2 \\
\leq \frac{\phi_3(\mu)^2}{\alpha(\mu) - 2 \phi_2(\mu) \varepsilon_M(\mu)} \Delta t \sum_{k'=1}^k b(t^{k'})^2 \\
\leq \frac{2 \phi_3(\mu)^2}{\alpha(\mu)^2} \Delta t \sum_{k'=1}^k b(t^{k'})^2. \\ 
\text{(A.15)}
\]

From (A.10) and invoking (A.12) and (A.15) we obtain
\[
m(c(\mu, t^k), e(\mu, t^k) + \Delta t \sum_{k'=1}^k a(e(\mu, t^{k'}), e(\mu, t^{k'}); \mu) \\
\leq 8m(u(\mu, t^k) - w_N(\mu, t^k), u(\mu, t^k) - w_N(\mu, t^k)) \\
+ 8\Delta t \sum_{k'=1}^k a(u(\mu, t^{k'}), u(\mu, t^{k'}); u(\mu, t^{k')}; \mu) \\
+ 4 \Delta t \alpha(\mu) \sum_{k'=1}^k \|u(\mu, t^{k'}) - w_N(\mu, t^{k'})\|_X^2 + \Delta t \alpha(\mu)^2 \sum_{k'=1}^k \|e(\mu, t^{k'})\|_X^2 \\
+ \Upsilon(\mu) \Delta t \sum_{k'=1}^k b(t^{k'})^2, \\ 
\text{(A.16)}
\]

where
\[
\Upsilon(\mu) = \frac{18}{\alpha(\mu)} \varepsilon_M(\mu)^2 \left( \phi_1(\mu)^2 + \phi_2(\mu)^2 \frac{2 \phi_3(\mu)^2}{\alpha(\mu)^2} \right).
\]
The desired result then directly follows from the fact $w_N(\mu, t^k)$ is the projection of $u(\mu, t^k)$ with respect to the $m + \Delta t a$ norm.

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**References**


