



A New Approach for Solving Volterra Integral Equations Using the Reproducing Kernel Method

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Abstract

This paper is concerned with a technique for solving Volterra integral equations in the reproducing kernel Hilbert space. In contrast with the conventional reproducing kernel method, the Gram-Schmidt process is omitted here and satisfactory results are obtained. The analytical solution is represented in the form of series. An iterative method is given to obtain the approximate solution. The convergence analysis is established theoretically. The applicability of the iterative method is demonstrated by testing some various examples.

Keywords : Reproducing kernel method; Volterra integral equations; Gram-Schmidt orthogonalization process.

1 Introduction

IN the last decades, the reproducing kernel method (RKM) has been a promising method which applied more and more for solving various problems such as ordinary differential equations, partial differential equations, differential-difference equations, integral equations, and so on (see e.g. [1]-[16] and references therein). Among many literatures related to RKM for solving various problems and even among a bunch of extensive works related to RKM for solving integral equations, we just mention some more interesting problems. An approximate solution of the Fredholm integral equation of the first kind in

the reproducing kernel space was presented by Du and Cui [8, 9], solution of a system of the linear Volterra integral equations was discussed by Yang et al. [10], solvability of a class of Volterra integral equations with weakly singular kernel using RKM was investigated in [11, 12, 13], Geng [14] explained how to solve a Fredholm integral equation of the third kind in the reproducing kernel space, and Ketabchi et al. [7] obtained some error estimates for solving Volterra integral equations using RKM. In [7] and some other places, a general technique for solving Volterra integral equations was discussed in the reproducing kernel space. This general technique is based on the Gram-Schmidt (GS) orthogonalization process. In this study, we aim to explain how to construct a reproducing kernel method without using this process. For this purpose, we consider the following nonlinear Volterra integral equation

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$$u(x) = F(x, u(x)), \quad (1.1)$$

where

$$F(x, u(x)) = f(x) + \int_0^x k(x, \xi) \mathcal{N}(u(\xi)) d\xi,$$

$x \in [0, 1]$, in which functions f and k and the nonlinear operator \mathcal{N} are considered such that Eq. (1.1) has a unique solution. Furthermore, we need to assume that F is continuous. The rest of the paper is organized as follows. In the next Section, some preliminaries are represented. The method implementation is discussed in Section 3. Section 4 is devoted to convergence analysis of the method. For confirming the theoretical results, some numerical examples are provided in Section 5. The paper will be closed by a brief conclusion in the last Section.

2 Preliminaries

In this section, we follow the recent work by Cui et al. [16] and represent some useful materials.

Definition 2.1 Let H be a Hilbert space of functions $f : \Omega \rightarrow \mathbb{R}$. Denote by $\langle \cdot, \cdot \rangle$ the inner product and let $\|\cdot\| = \sqrt{\langle \cdot, \cdot \rangle}$ be the induced norm in H . The function $R : \Omega \times \Omega \rightarrow \mathbb{R}$ is called a reproducing kernel of H if the followings are satisfied

- (i) $R_y(x) = R(x, y) \in H, \forall y \in \Omega,$
- (ii) $f(y) = \langle f(x), R_y(x) \rangle, \forall f \in H,$ for all $y \in \Omega.$

Definition 2.2 A Hilbert space H of functions on a set Ω is called a reproducing kernel Hilbert space if there exists a reproducing kernel R of H .

Remark 2.1 The existence of the reproducing kernel of a Hilbert space is due to the Riesz Representation Theorem. It is known that the reproducing kernel of a Hilbert space is unique.

Theorem 2.1 [5] The reproducing kernel R of reproducing kernel Hilbert space H is positive definite.

Definition 2.3 The function space $W_m[0, 1]$ is defined as follows

$$W_m[0, 1] = \{u | u^{(m-1)} \in AC[0, 1], u^{(m)} \in L^2[0, 1]\},$$

$AC[0, 1]$ is an absolutely continuous function in $[0, 1]$.

The inner product and norm in $W_m[0, 1]$ are defined respectively by

$$\begin{aligned} \langle u, v \rangle_{W_m} &= \\ &= \sum_{i=0}^{m-1} u^{(i)}(0)v^{(i)}(0) + \int_0^1 u^{(m)}(x)v^{(m)}(x)dx, \\ \forall u, v \in W_m[0, 1], \end{aligned}$$

$$\|u\|_{W_m} = \sqrt{\langle u, u \rangle_{W_m}}, \forall u \in W_m[0, 1].$$

In general, the function space $W_m[0, 1]$ is a reproducing kernel space and its reproducing kernel R_m has the following reproducing property $u(\cdot) = \langle u(x), R_m(x, \cdot) \rangle_{W_m}, \forall u \in W_m[0, 1]$. For $m = 1$, the function space $W_1[0, 1]$ is a reproducing kernel space and its reproducing kernel is

$$R(x, y) = \begin{cases} 1 + x, & x \leq y, \\ 1 + y, & x > y. \end{cases}$$

3 The method implementation

We rewrite Eq. (1.1) as follows

$$Lu(x) = f(x) + \int_0^x k(x, \xi) \mathcal{N}(u(\xi)) d\xi,$$

where $L : W_1[0, 1] \rightarrow W_1[0, 1]$ is an invertible bounded linear operator, \mathcal{N} is a nonlinear operator, and f is an arbitrary function in $W_1[0, 1]$. $W_1[0, 1]$ is a reproducing kernel space defined according to the highest derivatives involved in (1.1). We choose a countable set of points $\{x_i\}_{i=1}^\infty$ in the interval $[0, 1]$, and define

$$\phi_i(x) = R(x, x_i), \quad \psi_i(x) = L^* \phi_i(x),$$

where L^* is the adjoint operator of L . Obviously, in this paper L is the identity operator and therefore $\psi_i(x) = R(x, x_i)$.

Theorem 3.1 Let $\{x_i\}_{i=1}^\infty$ be dense in the interval $[0, 1]$. If Eq. (1.1) has a unique solution, then it can be represented as

$$u(x) = \sum_{j=1}^\infty a_j \psi_j(x), \tag{3.2}$$

where the coefficients a_j are determined by solving the following semi-infinite system of linear equations

$$Ba = F, \tag{3.3}$$

in which

$$B = [\psi_j(x_i)], \quad i, j = 1, 2, \dots,$$

$$a = [a_1, a_2, \dots]^T,$$

and

$$F = [F(x_1, u(x_1)), F(x_2, u(x_2)), \dots]^T.$$

Proof. Since $\{x_i\}_{i=1}^\infty$ is dense in the interval $[0, 1]$, then $\psi_j(x)$ is a complete system in $W_1[0, 1]$, see e.g. [16]. So the analytical solution can be represented as Eq. (3.2). Since

$$\begin{aligned} &\langle \psi_i(x), \psi_j(x) \rangle_{w_1} = \\ &\langle L^* \phi_i(x), \psi_j(x) \rangle_{w_1} = \\ &\langle \phi_i(x), L\psi_j(x) \rangle_{w_1} = L\psi_j(x)|_{x=x_i} = \psi_j(x_i), \end{aligned}$$

and

$$\begin{aligned} &\langle u(x), \psi_j(x) \rangle_{w_1} = \\ &\langle u(x), L^* \phi_j(x) \rangle_{w_1} = \\ &\langle Lu(x), \phi_j(x) \rangle_{w_1} = F(x_j, u(x_j)), \end{aligned}$$

according to the best approximation in Hilbert spaces [5], the coefficients a_j are determined by (3.3).

The approximate solution of the problem is the m -term intercept of the analytical solution which can be determined by solving a $m \times m$ system of linear equations. We need to construct an iterative method for solving (3.3). For this purpose, we choose the number of points m , the number of iterations n and put the initial function $u_{0,m}(x) = 0$. Then, the approximate solution of Eq. (1.1) is defined by

$$u_{n,m}(x) = \sum_{j=1}^m a_j^n \psi_j(x) = F(x_j, u_{n-1,m}(x_j)). \tag{3.4}$$

Remark 3.1 *There exists a unique solution for equations (3.4) due to the strictly positive definiteness property of the reproducing kernel.*

The results of this section can be summarized in the following algorithm.

Algorithm1.

1. Choose m collocation points in the interval $[0, 1]$.
2. Set $B = [\psi_j(x_i)]$, $i, j = 1, 2, \dots, m$.
3. Choose the number of iterations n .
4. Set $i = 0$.
5. Set the initial function $u_{0,m}(x) = 0$.
6. Set $i = i + 1$.
7. Set $F = [F(x_j, u_{i-1,m}(x_j))]^T$, $j = 1, \dots, m$.

8. Solve $Ba = F$.
9. Set $u_{i,m}(x) = \sum_{j=1}^m a_j^i \psi_j(x)$.
10. If $i < n$, then go to step 6, else stop the algorithm.

The conventional reproducing kernel method which used the GS orthogonalization process is represented in the following algorithm [7].

Algorithm2.

1. Choose m collocation points in the domain set $[0, 1]$.
3. Set $\psi_i(x) = R(x, x_i)$, $i = 1, \dots, m$.
4. Set $\bar{\psi}_i(x) = \sum_{k=1}^i \beta_{ik} \psi_k(x)$, $i = 1, \dots, m$, (β_{ik} which obtained by the GS process).
5. Choose an initial function $u_0(x)$.
6. Set $n = 1$.
7. Set $B_n = \sum_{l=1}^n \beta_{nl} F(x_l, u_{n-1}(x_l))$.
8. Set $u_n(x) = \sum_{j=1}^n B_j \bar{\psi}_j(x)$.
9. If $n < m$, then set $n = n + 1$ and go to step 7, else stop.

Remark 3.2 *In comparison with Algorithm 2, Algorithm 1 needs not to use the GS orthogonalization process but in step 8 of it, we must to solve a linear system. The coefficient matrix of this system is positive definite because of the positive definiteness of the kernel. Therefore, it needs to decompose matrix B once using the QR decomposition and to solve a triangular system in step 8.*

4 Convergence analysis

In this section, we show that the approximate solution $u_{n,m}$ is converged to the analytical solution u uniformly. At first, the following lemma is given.

Lemma 4.1 *For a positive constant M , $A = \{u \mid \|u\|_{w_1} \leq M\}$ is a compact set in the space $C[0, 1]$ provided that $\|u'\|_{w_1} \leq c$, where c is a constant.*

Table 1: Results of Algorithms 1 and 2.

m	Example 5.1		Example 5.2	
	E_m	$E_{5,m}$	E_m	$E_{5,m}$
2	2.71828	0.09544	1.00000	0.22222
4	1.33040	0.01045	0.97862	0.02663
8	0.69767	0.00187	0.69761	0.00478
16	0.35213	0.00036	0.40235	0.00103
32	0.17439	0.00005	0.21231	0.00024

Table 2: Results of Algorithms 1 and 2.

m	Example 5.3		Example 5.4	
	E_m	$E_{5,m}$	E_m	$E_{5,m}$
2	1.00000	0.09178	0.84147	0.01545
4	0.88923	0.00904	0.51528	0.00139
8	0.50274	0.00163	0.18976	0.00025
16	0.25626	0.00035	0.07976	0.00005
32	0.12715	0.00008	0.03665	0.00001

Proof. It is enough to show that A is a bounded and equicontinuous set [5]. Since

$$\|R(x, y)\|_{w_1}^2 = \langle R(x, y), R(x, y) \rangle_{w_1} = R(x, x) < c_0,$$

where c_0 is a positive constant, there exists a constant c_1 such that

$$|u(x)| = |\langle u(y), R(x, y) \rangle_{w_1}| \leq \|u(y)\|_{w_1} \|R(x, y)\|_{w_1} \leq c_1 \|u(y)\|_{w_1}.$$

For any $u \in A$, we have $|u(x)| \leq c_1 M$. Hence A is a bounded set in the space $C[0, 1]$. On the other hand,

$$|u'(x)| = |\langle u(y), \frac{\partial R(x, y)}{\partial x} \rangle_{w_1}| \leq \|u(y)\|_{w_1} \|\frac{\partial R(x, y)}{\partial x}\|_{w_1} \leq c_2 \|u(y)\|_{w_1} \leq c_2 M$$

Then for any $u \in A$ and $\epsilon > 0$, we have $|u(x+h) - u(x)| \leq |u'(\eta)| |h| \leq c_2 M |h|$ where $\eta \in [x, x+h]$. So, there exists $\delta = \frac{\epsilon}{c_2 M}$ such that for $|h| < \delta$, we get $|u(x+h) - u(x)| < \epsilon$. Hence A is an equicontinuous set in the space $C[0, 1]$.

Theorem 4.1 *If L is an invertible bounded linear operator and $F(x, u(x))$ is a nonlinear bounded operator, it can be deduced that $\{u_{n,m}(x)\}_{n=1}^\infty$ is the bounded sequence of functions in $w_1[0, 1]$.*

Proof. We can write

$$\begin{aligned} & \|u_{n,m}(x)\|_{w_1}^2 \\ &= \langle u_{n,m}(x), u_{n,m}(x) \rangle_{w_1} \\ &= \langle \sum_{j=1}^m a_j \psi_j(x), \sum_{l=1}^m a_l \psi_l(x) \rangle_{w_1} \\ &= \sum_{j=1}^m a_j \langle \psi_j(x), \sum_{l=1}^m a_l \psi_l(x) \rangle_{w_1} \\ &= \sum_{j=1}^m a_j \langle \phi_j(x), \sum_{l=1}^m a_l \psi_l(x) \rangle_{w_1} \\ &= \sum_{j=1}^m a_j (\sum_{l=1}^m a_l \psi_l(x_j)) \\ &= a^T B a, \end{aligned}$$

where

$$a = [a_j], \quad j = 1, 2, \dots, m.$$

Now, since

$$B = [\psi_j(x_i)], \quad i, j = 1, 2, \dots, m,$$

is a positive definite matrix then we have $a = B^{-1} F$ and the assumed assumptions imply that

$$\|u_{n,m}(x)\|_{w_1} \leq M,$$

where M is a constant.

Theorem 4.2 *Assume that $\{x_i\}_{i=1}^\infty$ is dense in $[0, 1]$ and the assumptions of Theorem 4.1 and Lemma 4.1 hold. Then the approximate solution $u_{n,m}$ is converged to the analytical solution u .*

Proof. For $j = 1, 2, \dots, m$ and $n = 1, 2, \dots$, we have $Lu_{n,m}(x_j) = F(x_j, u_{n-1,m}(x_j))$.

According to Lemma 4.1, there exists a convergent subsequence $\{u_{n_l,m}(x)\}_{l=1}^\infty$ of $\{u_{n,m}(x)\}_{n=1}^\infty$ such that $u_{n_l,m}(x) \rightarrow u_{n,m}(x)$, uniformly as $l \rightarrow \infty$, $m \rightarrow \infty$. Then for $j = 1, 2, \dots, m$ and $n = 1, 2, \dots$, we derive

$$Lu_{n_l,m}(x_j) = F(x_j, u_{n_l-1,m}(x_j)). \tag{4.5}$$

Since the operators L and F are both continuous (according to the structure of L and assumption on F), after taking limit from both sides of (4.5), it can be inferred that u is the analytical solution of Eq. (1.1). So $u_{n_i,m}(x)$ is the approximate solution of Eq. (1.1).

5 Examples

In this section, we compare results of both Algorithms in solving four various problems using the following norms

$$\|u - u_{n,m}\|_{\infty} \simeq E_{n,m} = \max_{1 \leq i \leq m} |u(x_i) - u_{n,m}(x_i)|,$$

$$\|u - u_m\|_{\infty} \simeq E_m = \max_{1 \leq i \leq m} |u(x_i) - u_m(x_i)|,$$

where $u_{n,m}$ and u_m are approximate solutions obtained by Algorithms 1 and 2, respectively and u is the exact solution. The results of Tables 1, 2 (for $n = 5$) confirm the superiority of Algorithm 1.

Example 5.1 If $k(x, \xi) = x^2\xi$, $f(x) = \exp(x) - x^2 + x^2 \exp(x) - x^3 \exp(x)$ and $\mathcal{N}(u(x)) = u(x)$, then the Volterra integral equation (1.1) has the following exact solution $u(x) = \exp(x)$.

Example 5.2 If $k(x, \xi) = x^{\frac{7}{2}}\xi$, $f(x) = x^{\frac{7}{2}} - \frac{2}{11}x^9$ and $\mathcal{N}(u(x)) = u(x)$, then the Volterra integral equation (1.1) has the following exact solution $u(x) = x^{\frac{7}{2}}$.

Example 5.3 If $k(x, \xi) = x^2\xi$, $f(x) = \frac{1}{2}x^2 + \frac{1}{2}x^2 \cos(x^2)$ and $\mathcal{N}(u(x)) = \sin(u(x))$, then the Volterra integral equation (1.1) has the following exact solution $u(x) = x^2$.

Example 5.4 If $k(x, \xi) = \sin(x - \xi)$, $f(x) = \sin(x) + \frac{2}{3} \cos(x) - \frac{1}{6} \cos(2x) - \frac{1}{2}$ and $\mathcal{N}(u(x)) = u(x)^2$, then the Volterra integral equation (1.1) has the following exact solution $u(x) = \sin(x)$.

6 Conclusion

In this work, we proposed an iterative algorithm for solving nonlinear volterra integral equations on the basis of the reproducing kernel Hilbert space without using the Gram-Schmidt orthogonalization process. The results of some numerical examples show that the present method could be an accurate and reliable analytical-numerical

technique. Examples presented here belong to different categories such as linear or nonlinear problem with smooth or none-smooth kernel and solution. Nevertheless, our results only apply to the given examples; this, of course, does not mean that it holds in general. The advantage of the approach is that the method can be easily implemented. It seems that the method can be also applied for solving other nonlinear integral equations.

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