Numerical Solution of Stiff and Singularly Perturbed Boundary Value Problems With a Segmented-Adaptive Formulation of the Tau Method

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To Professor Y. L. Luke, in memoriam

Abstract. This paper concerns the application of Ortiz' recursive formulation of the Tau method to the construction of piecewise polynomial approximations to the solution of linear and nonlinear boundary value problems for ordinary differential equations. A practical error estimation technique, related to the concept of correction in Zadunaisky's sense, is considered and used in the design of an adaptive approach to the Tau method. It proves efficient in the numerical treatment of problems with rapid functional variations, stiff and singularly perturbed problems. A technique of increased accuracy at matching points of segmented Tau approximants is also discussed and successfully applied to several problems. Numerical examples show that, for a given degree of approximation, our segmented Tau approximant gives an accuracy comparable to that of the best segmented approximation of the exact solution by means of algebraic polynomials.

1. **Introduction.** We discuss the use of Ortiz' recursive formulation of the Tau method [23]–[25] in the numerical solution of boundary value problems for linear and nonlinear differential equations defined over an interval $a \leqslant x \leqslant b$. We consider *global* approximations over $[a, b]$, with a single polynomial expression, and *segmented* forms based on a step-by-step formulation of the Tau method considered by Ortiz in [26].

The Tau approximate solution of a differential problem defined by a differential operator D is represented in terms of the elements of a sequence Q of *canonical polynomials*. Such a sequence is uniquely determined by D, it is independent of the specific boundary conditions of the problem, and of the particular interval $[a, b]$ in which the solution is required. These properties make possible the use of segmentation within the framework and with the software [32] designed for the recursive formulation of the Tau method. The concept of correction, in Zadunaisky's sense [37] (see also Stetter [36]), is discussed in the context of the Tau method and related to a practical error estimation technique. This technique, based on *Tau estimators* introduced here, is systematically applied to all examples, linear or nonlinear. It is

also used in the control of the segmentation process: an adaptive form of the Tau method, based on this, is considered.

A technique of increased accuracy at matching points of segmented Tau approximations, introduced by Ortiz in [27] for initial value problems, is successfully used in linear and nonlinear boundary value problems. Numerical comparisons with other standard methods and, in particular, with an accurate technique of collocation with splines followed by a finite difference correction (see [18]) is definitely favorable to our approach. We find that, for a given degree of approximation, our *segmented* Tau approximants are close to the best segmented approximations by algebraic polynomials. This optimal result has been reported for the global case by Freilich and Ortiz in a recent paper [6].

Convergence results and error bounds for Tau approximations of the solution of differential equations are discussed by Luke [13], Lanczos [10]–[12], and Ortiz and Pham [30]–[31]. Systems of differential equations have been discussed by Freilich and Ortiz [6] and Crisci and Russo [3]. The effect of approximating the coefficients of a differential equation on the accuracy of the Tau approximant is discussed by Namasivayam and Ortiz in [17]. An interesting feature of the Tau method is the fact that no trial solutions, approximate quadratures or large matrix inversions are required. A summary of the results of this paper was presented to the Dundee Conference of Numerical Analysis, June 1981.

2. Recursive Formulation of the Tau Method: Some Basic Definitions and Notation. Let $\mathbf{P}_j$ be the class of polynomials of degree less than or equal to j. We shall consider the equation defined by the differential operator D:

$$(1) \quad Dy(x) := p_\nu(x)y^{(\nu)}(x) + \cdots + p_1(x)y^{(1)}(x) + p_0(x)y^{(0)}(x) = f(x),$$

where either $p_i(x) \in \mathbf{P}_{\alpha_i}$, $i = 0(1)\nu$, and $f(x) \in \mathbf{P}_F$, or they are close polynomial approximations of given functions. The symbol $y^{(i)}(x)$ stands for the ith derivative of $y(x)$, and $y^{(0)}(x) \equiv y(x)$. The solution $y(x)$ of (1) satisfies boundary conditions of the general form

$$(2) \quad \sum_{i=0}^{\nu-1} \left[a_{ri}y^{(i)}(a) + b_{ri}y^{(i)}(b) \right] = A_r, \qquad r = 1(1)\nu,$$

where some of the coefficients a_{ri}, b_{ri} may be equal to zero. We shall follow Ortiz' algebraic theory of the Tau method, of which we will give here some basic definitions and results. Further details and proofs can be found in [23]–[25], and [14].

A useful number associated with any differential operator D of the class $\mathscr{D}$, characterized by (1), is its height h,

$$h := \max_{0 \leqslant i \leqslant \nu} (\alpha_i - i).$$

Each differential operator $D \in \mathscr{D}$ is uniquely associated with a sequence Q of *canonical polynomials* $Q_n(x)$ defined for all indices $n \in \mathbf{N} - S$. The set $\mathbf{N} := \{0, 1, 2, 3, \ldots\}$; S is finite, usually very small, and the number s of its elements is bounded by $\nu + h$. The polynomials $Q_n(x)$ satisfy the functional equation

$$(3) \quad DQ_n(x) = x^n + r_n(x), \qquad n \in \mathbf{N} - S,$$

where $r_n(x) \in \mathbf{R}_S = \operatorname{span}_{i \in S}\{x^i\}$ is called the *residual* of $Q_n(x)$. If all powers of x can be obtained as images of polynomials under the differential operator D, then S will be empty. If D is applied to the monomial x^m, we obtain the *generating polynomials*

$$(4) \qquad Dx^m = \sum_{i=0}^{n} a_i x^i, \qquad m \in \mathbf{N},$$

from which a recursive relation, involving at most $\nu + h$ (assuming $p_0(x) \neq 0$) canonical polynomials, can immediately be deduced if $a_n \neq 0$ (see Ortiz' Theorem 3.3 in [24]). Otherwise, (4) is used to find the minimal set S of indices of undefined canonical polynomials and $\operatorname{Ker}(D)$, the algebraic kernel of D, which contains all the exact polynomial solutions of (1). The recursive relation for the $Q_n(x)$'s is self starting and the polynomials themselves depend neither on the supplementary conditions ((2) in our case) imposed on $y(x)$, nor on the interval in which the solution is required. These properties will be used in the design of a segmented formulation of the Tau method of [26] for boundary value problems.

Let $\mathbf{v} = \{v_i(x)\} = V\mathbf{x}$ be a polynomial basis defined by a lower triangular matrix $V = ((v_{ij}))$, $i, j \in \mathbf{N}$, acting on $\mathbf{x} = (1, x, x^2, \ldots)'$. Clearly $\hat{Q} = \{\hat{Q}_n(x)\}$, $n \in \mathbf{N} - S$, is such that $D\hat{Q}_n(x) = v_n(x) + \hat{r}_n(x)$, $\hat{r}_n(x) \in \mathbf{R}_S$, if

$$(5) \qquad \hat{Q}_n(x) = \sum_{j=0}^{n} v_{nj} Q_j(x), \quad \text{with } j \notin S.$$

Definition 1. A polynomial $g(x)$ is called a *Tau approximant* of order n of $y(x)$ if it satisfies exactly the differential equation (1) with a polynomial perturbation term $H_n(x) \in \mathbf{P}_n$, and if it satisfies exactly boundary conditions (2).

Clearly $g(x) \in \mathbf{P}_{n-h}$. Let

$$(6) \qquad H_n(x) = \sum_{i=0}^{m} \tau_i^{(n)} v_{n-i}(x) \in \mathbf{P}_n$$

be the perturbation term, expressed in the basis $\mathbf{v}$, and let $\tau_i^{(n)}$, $i = 0(1)m$, be free parameters. Let us assume further that

$$f(x) = \sum_{i=0}^{F} f_i v_i(x),$$

and that n is chosen sufficiently large for $F \leqslant n$. Then,

$$(7) \qquad y_n(x) := \sum_{i=0}^{m} \tau_i^{(n)} \hat{Q}_{n-i}(x) + \sum_{i=0}^{F} f_i \hat{Q}_i(x), \quad \text{with } i \notin S,$$

satisfies the equation $Dy_n(x) = H_n(x) + f(x)$ if s parameters $\tau_i^{(n)}$ are chosen in such a way that the residuals of $Dy_n(x)$ match the components of $H_n(x) + f(x)$ belonging to $\mathbf{R}_S$. If ν further parameters $\tau_i^{(n)}$ are fixed for $y_n(x)$ to satisfy exactly the boundary conditions (2), then $y_n(x)$ is a Tau approximant of order n of $y(x)$. If there exist t exact polynomial solutions of (1), then $m = s + \nu - t - 1$.

Finally, if (1) is given in an integrated form, through an integral operator I, a sequence $\overline{Q} = \{\overline{Q}_n(x)\}$, $n \in \mathbf{N} - S_I$, such that $I\overline{Q}_n(x) = x^n + \bar{r}_n(x)$, $\bar{r}_n(x) \in \mathbf{R}_{S_I}$ can be constructed immediately in terms of the sequence Q associated with D:

$$(8) \qquad \overline{Q}_n(x) = \frac{n!}{(n - \nu)!} Q_{n-\nu}(x), \quad \text{with } S_I = \{ n \in \mathbf{N} : n < \nu, n - \nu \in S \}.$$

A variety of choices of basis $\mathbf{v}$ is possible in the Tau method. If $\mathbf{v}$ is $\{x^n\}$, $n \in \mathbf{N}$, the Tau method realizes the power series expansion method, where a high accuracy is to be expected near the point of expansion. Lanczos [11]–[12] suggested the choice of Chebyshev and Legendre polynomials to obtain a better distribution of errors in the equation over the interval in which the approximate solution is required. The error in the equation, namely $H_n(x)$, is related to the error of approximation

$$e_n(x) = y_n(x) - y(x)$$

through the inverse operator of D: $D^{-1}H_n(x) = e_n(x)$, on account of the linearity of D. The choice of $H_n(x)$ close to a best uniform approximation of the function identically equal to zero in $[a, b]$ by means of algebraic polynomials is an attempt to take advantage of the relation between $H_n(x)$ and $e_n(x)$. However, that relation requires a careful analysis if quantitative results are required. The behavior of $e_n(x)$ relative to the perturbation term $H_n(x)$ is discussed by Namasivayam and Ortiz [16].

Remark 1. If the solution $y(x)$ of (1)–(2) is a polynomial of degree k, any Tau approximate solution of degree $\geq k$ will detect it exactly. In this sense we say that the Tau method is *exact of degree k*.

3. Correction of Tau Approximants. We now consider the correction technique of Zadunaisky [37] and Stetter [36] in the context of the Tau method.

THEOREM 1. *Correction of a Tau approximant leads to a Tau approximant of a higher order.*

Proof. Let us consider the error function $e_n(x)$. From (1) it follows that $e_n(x) \in C^{(\nu)}[a, b]$ and that it satisfies the differential equation $De_n(x) = H_n(x)$, where all τ-parameters in $H_n(x)$ are fixed. It also satisfies boundary conditions (2) with $A_r = 0$, for $r = 1(1)\nu$.

Let $[e_n(x)]_m$ be an mth order ($m > n$) Tau approximant of the error function. Then the polynomial $[e_n(x)]_m$ satisfies the differential equation

$$(9) \qquad D[e_n(x)]_m = H_n(x) + H_m(x) = f(x) + H_n(x) - [f(x) - H_m(x)],$$

with homogeneous boundary conditions (2). Let $y_m(x)$ be the mth order Tau approximant of $y(x)$. Then, $e_n(x) - e_m(x) = y_n(x) - y_m(x)$. From the uniqueness of the Tau approximant for a given order m (see Ortiz [24]) and (9), it follows that $[e_n(x)]_m = e_n(x) - e_m(x)$. Thus, $y_n(x)$ plus the correction term $[e_n(x)]_m$ equals $y_m(x)$.

Definition 2. $[e_n(x)]_m$ defined by (9) is called the *Tau estimator* of order n, m. Examples of Tau approximants corrected with the Tau estimator, for $m \geq n + 1$, are given in the last section of this paper.

4. Implicit Matching of Tau Approximants. Let

$$\Pi = \left\{ x_0 = a < x_1 < x_2 < \cdots < x_{p-1} < x_p = b \right\}$$

be a partition of the interval $[a, b]$ into subintervals $[x_{j-1}, x_j], j = 1(1)p$,

Definition 3. The vector

$$(10) \qquad \mathbf{y}_n(x) = \left[y_{n1}(x), y_{n2}(x), \dots, y_{np}(x) \right],$$

with $y_{nj}(x)$ defined for $x_{j-1} \leqslant x \leqslant x_j, j = 1(1)p$, is a piecewise Tau approximant of order n of the solution $y(x)$ of the boundary value problem (1)–(2) if each of the $y_{nj}(x), j = 1(1)p$, satisfies (1) with a polynomial perturbation term $H_n^j(x)$ defined for $x_{j-1} \leqslant x \leqslant x_j$, and

$$(11\text{-a}) \qquad \sum_{i=0}^{\nu-1} \left[a_{ri} y_{n1}^{(i)}(a) + b_{ri} y_{np}^{(i)}(b) \right] = A_r \quad \text{for } r = 1(1)\nu,$$

$$(11\text{-b}) \qquad y_{nj-1}^{(i)}(x_{j-1}) = y_{nj}^{(i)}(x_{j-1}) \quad \text{for } j = 2(1)p, i = 0(1)\nu - 1.$$

Remark 2. Condition (11-a) imposes on $\mathbf{y}_n(x)$ the ν boundary conditions given by (2), while (11-b) is a continuity requirement for $y_{nj}(x)$ and its $\nu - 1$ derivatives at the $p - 1$ interior partition points of Π. If $S \neq \varnothing$, matching coefficients of terms belonging to $\mathbf{R}_S$ in the right-hand side of the differential equation with the residuals of $Dy_{nj}(x)$ provides us with s conditions in each of the p subintervals defined by the partition. Thus, a total of $p(s + \nu)$ conditions.

THEOREM 2. *The construction of an nth order piecewise Tau approximation* $\mathbf{y}_n(x)$ *of the solution* $y(x)$ *of boundary value problem* (1)–(2) *depends on only one matrix V and one canonical sequence Q.*

Proof. Let us consider the family of mappings $\mathscr{S}_j, j = 1(1)p$ (see Ortiz [26]) acting on $y_n(x)$ of (7) and such that

$$\mathscr{S}_j y_n(x) = y_{nj}(x), \qquad x_{j-1} \leqslant x \leqslant x_j,$$

and let us call $\tau_{ij}^{(n)}$ the free parameters corresponding to $y_{nj}(x)$. Without loss of generality, let us assume that $a = 0$, and let $E_j = 1/(x_j - x_{j-1}), j = 1(1)p$. The mappings $\mathscr{S}_j$ admit a trivial realization: replacing v_{nj} by $v_{ni}^j = v_{ni}(E_j)$, we find that $((v_{ni}^j))$ defines the basis $\mathbf{v}^j$ corresponding to the interval $x_{j-1} \leqslant x \leqslant x_j$, for $j = 1(1)p$. Taking into account (7) the result is proved.

We now discuss the computational implications of Theorem 2. Let $\tau_j^{(n)} = (\tau_{ij}^{(n)})'$, $i = 0(1)s + \nu - 1$, be a vector, the elements of which are the Tau parameters of the component $y_{nj}(x)$ of $y_n(x)$, for $j = 1(1)p$, and let $\tau^{(n)} = (\tau_1^{(n)\prime}, \dots, \tau_p^{(n)\prime})'$ be a vector with $p(s + \nu)$ components. Let $a = x_{j-1}, b = x_j$. Then (7) gives us $y_{nj}(x)$ for $j = 1(1)p$. The piecewise polynomial Tau approximant $\mathbf{y}_n(x)$ of $y(x)$ will be determined if $\tau^{(n)}$ is found.

From Remark 2 it follows that $\tau^{(n)}$ is implicitly defined by a system of linear algebraic equations of the form

$$(12) \qquad Z_n \tau^{(n)} = \mathbf{W}_n.$$

Let us split $y_{nj}(x)$ (see (7), with $a = x_{j-1}, b = x_j$) into two parts: one dependent on the Tau parameters, and then related to Z_n, and the other independent of them, thus

related to $\mathbf{W}_n$. We shall write $\phi_{nj}(x)\tau_j^{(n)}$ for the vector with ν rows, containing as elements that first part of $y_{nj}(x)$ and its $\nu - 1$ derivatives. Let $\xi_{nj}(x)$ be a vector containing in each of its rows the second part of $y_{nj}(x)$ and its $\nu - 1$ derivatives. The *residual* $Dy_{nj}(x)$ of $y_{nj}(x)$ belongs to $\mathbf{R}_S$, which is a subspace generated by s linearly independent basis elements. We will split the residual of $y_{nj}(x)$ in the same way as before, placing the s components in $\mathbf{R}_S$ of the first part, dependent on the Tau terms, in successive rows of a vector $\rho_{nj}\tau_j^{(n)}$ and the rest of it in the successive rows of a vector η_{nj}.

With the help of matrices ϕ_{nj} and ρ_{nj}, of orders $\nu \times (s + \nu)$ and $s \times (s + \nu)$, respectively, we can discuss the structure of matrix Z_n of (12). Z_n is made up of blocks $\alpha_j^{(n)}$, each one of them related to one component of $\mathbf{y}_n$, for $j = 1(1)p$. These blocks have the form

$$\alpha_j^{(n)} = \begin{array}{|c|}\hline \phi_{nj}(x_{j-1}) \\ \hline \rho_{nj} \\ \hline -\phi_{nj}(x_j) \\ \hline \end{array} \quad \text{for } j = 2(1)p - 1;$$

in $\alpha_p^{(n)}$ the third subblock is missing, and $\alpha_1^{(n)}$ has the same second and third subblocks, but a different first subblock γ_{n1}. Let

$$\mathscr{A} = ((a_{ri})), \quad \mathscr{B} = ((b_{ri})), \quad r = 1(1)\nu, i = 0(1)\nu - 1, \text{ and } \mathbf{A} = (A_1, \ldots, A_\nu)',$$

where a_{ri}, b_{ri}, and A_i are the coefficients of the boundary conditions (11-a); then γ_{n1} is such that

$$\gamma_{n1}\tau^{(n)} = \mathscr{A}\phi_{n1}(a)\tau_1^{(n)} + \mathscr{B}\phi_{np}(b)\tau_p^{(n)}.$$

Matrix Z_n is constructed by linking the blocks $\alpha_j^{(n)}$ in such a way that the third subblock of $\alpha_j^{(n)}$ faces the first subblock of $\alpha_{j+1}^{(n)}$ for $j = 1(1)p - 1$. The first ν components of $\mathbf{W}_n$ have the following form:

$$\mathbf{A} - \mathscr{A}\xi_{n1}(a) - \mathscr{B}\xi_{np}(b);$$

with this choice the boundary conditions (11-a) are incorporated into (12). The remaining elements of $\mathbf{W}_n$ follow if ξ_{nj} and η_{nj} are assembled following the same rules as for the corresponding Tau-dependent elements in Z_n. The rows of Z_n with ϕ's realize the continuity conditions (11-b), while the rows with ρ's make sure that the sum of residuals of each component y_{nj} match the terms in $\mathbf{R}_S$ on the right-hand side of the Tau problem for each $x_{j-1} \leqslant x \leqslant x_j$, $j = 1(1)p$. We have assumed that $t = 0$, if this is not the case, only $s + \nu - t$ Tau parameters are required, as t free parameters are provided by the exact polynomial solutions of $Dy(x) = 0$.

Remark 3. The following observation enables us to introduce a considerable simplification in the computational procedure. Let us assume for simplicity that $|x_{j-1} - x_j| = |b - a|/p = $ constant for $j = 1(1)p$; otherwise a scaling factor E_j will be required. Let us single out one of the subintervals defined by the partition Π, say the first one $[x_0, x_1]$. On account of Theorem 2, all components $y_{nj}(x)$ of $\mathbf{y}_n(x)$ can be determined as shifts of a *Tau master element* (see [26]) $y_{n1}(x)$ if we only replace x_{j-1} by x_1 in the left-hand side of (11-b), and x_{j-1} by x_0 in its right-hand side. Then, the only basis required will be $\mathbf{v}$, defined for $x_0 \leqslant x \leqslant x_1$, and the same sequence of

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canonical polynomials (see (5)) will be sufficient to construct all components of the piecewise Tau approximant $y_n(x)$. The procedure for the construction of $y_{n1}(x)$ could be regarded as an *integration formula* based on the Tau method of [25] and specifically designed for each differential equation (1)–(2). The graphs of successive segmented Tau approximants $y_{nj}(x)$, $j = 1(1)p$, will all be in the interval $[x_0, x_1]$. The ordinate at the endpoint of the graph of $y_{nj}(x)$ will be the same as that of $y_{nj+1}(x_0)$. Sliding these graphs in the direction of $\overline{x_0 x_p}$ we obtain the graph of $y_n(x)$.

If Remark 3 is taken into account the structure of Z_n is greatly simplified: except for the subblock representing the boundary conditions, all blocks $\alpha_j^{(n)}$ are similar.

As early as 1956 Lanczos observed a significant increase in accuracy at the endpoint of the approximation range of a Tau approximant when Chebyshev polynomials are replaced by Legendre polynomials (see Lanczos [11], [12] and Luke [13]). Ortiz [26] used this property in the design of a segmented formulation of the Tau method for initial value problems, with increased accuracy at matching points. Examples given in Section 7 of this paper show that the accuracy of segmented Tau approximants for linear and nonlinear boundary value problems improves by the use of the technique of increased accuracy at matching points. Theoretical results in this direction have been recently reported by Freilich and Ortiz [7] and by Namasivayam and Ortiz [16]. As collocation is a special realization of the Tau method, where the collocation nodes are the zeros of $H_n(x)$, it is clear that Lanczos' observation, the results reported in this paper, and those of [24] and [26] apply to the collocation method.

6. Nonlinear Boundary Value Problems and the Tau Method. The numerical solution of nonlinear boundary value problems with the recursive formulation of the Tau method [24] is based on the approximation of the solution of the nonlinear problem by a sequence of Tau approximants $y_{n,k}(x)$ of linear boundary value problems with variable coefficients. Each of the $y_{n,k}(x)$ is used, in an iterative cycle, to represent the nonlinear terms of the given differential equation, and gives a new $y_{n,k+1}(x)$. The fixed point of such a sequence is, under convergence conditions depending on the linearization scheme used, the function $y(x)$, solution of the initial problem. Details of this procedure can be found in Ortiz [28]. In practice such a process is only repeated a small number, N, of times and stopped when the maximum difference between the coefficients of two successive approximations is smaller than a given tolerance parameter T, specified beforehand. A test is also made on the size of the perturbation term corresponding to that approximation. If the sequence of approximants does not reach the tolerance parameter T after N cycles, either the degree of the Tau approximants is increased, the initial approximation redefined or the interval segmented by using the technique sketched in Section 5. The initial approximation is usually chosen to be a polynomial satisfying (2). In some special cases the differential equation allows for the immediate determination of an algebraic curve with a contact of order $\nu - 1$ with the solution $y(x)$ which is equally effective. For instance, a tangent to $y(x)$ in the case of a second order differential equation. The technique of increased convergence at matching points improves the accuracy of numerical results, as will be shown in Section 7.

Remark 4. In the case of nonlinear boundary value problems the piecewise Tau approximants $y_{nj,k}(x)$, at stage k of the process, satisfy different differential equations in each subinterval $[x_{j-1}, x_j], j = 1(1)p$.

7. Numerical Examples.

(1) *Dirichlet and Neumann Linear Boundary Value Problems.* Let us consider the problem (see de Boor and Swartz [4]):

$$(14) \quad \begin{cases} Dy(x) := y''(x) - 4y(x) = 4\cosh 1, \\ y(0) = y(1) = 0, \qquad 0 \leqslant x \leqslant 1, \end{cases}$$

for which results obtained by using finite difference residual correction of a collocation solution constructed with four cubic splines centered at the points 0, 1/3, 2/3, and 1, have been reported recently by Oliveira [18]. In Table 1 we are only concerned with *global* Tau approximants defined for $0 \leqslant x \leqslant 1$. For both the differential form (14) and the corresponding integrated form we compare the first Tau correction, defined by $[e_n(x)]_{n+1}$ for $n = 3(2)9$, with the exact error. We wish to remark that the integrated form is never constructed, its Tau approximant is computed by using the result of (8).

TABLE 1

Global Tau approximants and their corrections

Degree	DIFFERENTIAL FORM		INTEGRATED FORM	
	First Tau Correction	Exact Error	First Tau Correction	Exact Error
3	7.369×10^{-2}	7.415×10^{-2}	9.866×10^{-3}	9.940×10^{-3}
5	4.576×10^{-4}	4.589×10^{-4}	7.358×10^{-5}	7.390×10^{-5}
7	1.275×10^{-6}	1.278×10^{-6}	3.172×10^{-7}	3.181×10^{-7}
9	3.378×10^{-9}	3.384×10^{-9}	9.048×10^{-10}	9.064×10^{-10}

In Table 2 we present the same information when segmentation is used: four Tau approximants are constructed over subintervals of $[0, 1]$ of equal length. We remark that for $n \geqslant 4$ the approximation of the differential form is more accurate than that of the integrated form.

TABLE 2

Four piecewise Tau approximants over the equally

segmented interval $0 \leqslant x \leqslant 1$ and their corrections

Degree	DIFFERENTIAL FORM		INTEGRATED FORM	
	First Tau Correction	Exact Error	First Tau Correction	Exact Error
3	2.072×10^{-3}	2.073×10^{-3}	2.016×10^{-3}	2.032×10^{-3}
4	1.141×10^{-5}	1.196×10^{-5}	1.405×10^{-5}	1.644×10^{-5}
5	5.448×10^{-7}	5.470×10^{-7}	2.373×10^{-6}	2.386×10^{-6}
6	2.115×10^{-9}	2.235×10^{-9}	1.163×10^{-8}	1.282×10^{-8}
7	1.192×10^{-10}	1.196×10^{-10}	1.181×10^{-9}	1.186×10^{-9}
8	3.766×10^{-13}	3.908×10^{-13}	4.440×10^{-12}	4.750×10^{-12}

In Table 3 we attempt a ranking of our global and segmented approximations before and after correction with the Tau estimator $[e_n(x)]_{n+1}$. After correction, a *global* Tau approximant of the solution of (14) is more accurate than the piecewise polynomial approximation obtained by Oliveira [18] by using collocation with four cubic splines and a finite difference correction.

If four cubic Tau approximants are corrected with the Tau estimator, the accuracy over Oliveira's approximation increases by a factor of 3.169×10^{-3}. By using the technique of increased accuracy at matching points, which accounts for a switch from Chebyshev to Legendre polynomials in the basis **v**, that factor, now 1.619×10^{-4}, would show an even higher accuracy.

TABLE 3

A hierarchy of global and segmented approximants

of the boundary value problem (14)

Numerical technique	Maximum Abs. Error	Type of Approximant
Corrected piecewise cubic differential Tau with the use of the technique of increased accuracy at matching points	5.8×10^{-7}	Segmented
Corrected piecewise cubic differential Tau	1.1×10^{-5}	Segmented
Corrected piecewise cubic integrated Tau	1.4×10^{-5}	Segmented
Corrected cubic global integrated Tau	7.4×10^{-5}	Non-segmented
Piecewise cubic differential Tau with the use of the technique of increased accuracy at matching points	1.6×10^{-4}	Segmented
Corrected cubic global differential Tau	4.6×10^{-4}	Non-segmented
Finite differences corrected collocation with four cubic splines [18]	3.6×10^{-3}	Segmented
Piecewise cubic differential Tau	4.0×10^{-3}	Segmented
Cubic global integrated Tau	1.0×10^{-2}	Non-segmented
Collocation with cubic splines [18]	1.5×10^{-2}	Segmented
Cubic global differential Tau	7.4×10^{-2}	Non-segmented

Remark 4. The accuracy obtained with the Tau approximant at the top of Table 3 (5.827×10^{-7}) is almost identical to the upper bound for the best uniform segmented approximation of the exact solution $y(x)$ by means of four algebraic polynomials of degree four, which is the degree of the corrected Tau pieces. The upper bound was estimated on the basis of Lagrange's error formula (see Meinardus [15]). Therefore, it seems possible to use segmented Tau approximants as an initial guess for the construction of segmented best approximations.

A similar linear boundary value problem, now with Neumann conditions, and over a longer interval is:

$$\begin{cases} Dy(x):= \ y''(x) - y(x) = 1, \\ y(0) = 0, \quad y'(20) = 1, \qquad 0 \leqslant x \leqslant 20, \end{cases}$$

considered in detail by Scott [35] with finite difference techniques.

Table 4 displays the maximum of the value of the first Tau estimator and of the exact absolute error for some of the Tau approximants considered before. They have been constructed either over [0, 20], or over four equal and consecutive subintervals of [0, 20]. In this, as in many other Neumann problems, the differential form gives consistently better results than the integrated one, even for small values of n (see Onumanyi [20]).

TABLE 4

Tau approximation of a Neumann problem over a large interval

Approximant	$n = 7$		$n = 8$	
	Exact error	Tau estimator	Exact error	Tau estimator
Global integral Tau	2.1×10^{-1}	2.1×10^{-1}	5.4×10^{-2}	5.3×10^{-2}
Global differential Tau	1.1×10^{-1}	1.3×10^{-1}	2.1×10^{-2}	2.3×10^{-2}
Piecewise integral Tau	8.8×10^{-4}	8.7×10^{-4}	1.5×10^{-4}	1.4×10^{-4}
Piecewise differential Tau	1.4×10^{-4}	1.3×10^{-4}	2.4×10^{-5}	2.4×10^{-5}

(2) *A Stiff Boundary Value Problem With Nonpolynomial Coefficients.* If coefficients or the right-hand side of a differential equation are not polynomials, the Tau method can be used to find polynomial approximations of them. The same applies to transcendental nonlinear terms (see for example Ortiz and Samara [33]).

Let us consider the singular perturbation boundary value problem

(15) $$\begin{cases} Dy(x):= \ y''(x) - Py(x) = \cos x, \\ y(0) = y(\pi/2) = 1, \qquad 0 \leqslant x \leqslant \pi/2 \end{cases}$$

(see Guderley [8]). The nonpolynomial term will be replaced by a polynomial approximation of degree 14 which, in this case, we generate with the present Tau method technique.

The graph of the solution $y(x)$ of the singularly perturbed boundary value problem (15) is, for large P, close to $y = 0$ inside the interval $(0, \pi/2)$, and jumps to $y = 1$ when x approaches either $x = 0$ or $x = \pi/2$ (see Figure 1).

We have computed Tau approximations of $y(x)$ for $P = 1000$ and for $P = 10000$. They are global Tau approximations of a moderately large degree ($n = 27$) and segmented Tau approximants of a lower degree ($n = 14$). The ability of these approximations to follow the rapid variation of $y(x)$ near the endpoints of the interval $[0, \pi/2]$ is reported in Table 5 for both Chebyshev and Legendre perturbation terms. The graphs of the two segmented Tau approximants for $P = 1000$ and $P = 10000$ are reproduced in Figure 1.

TABLE 5

Tau approximation of a stiff boundary value problem

with a nonpolynomial term (15)

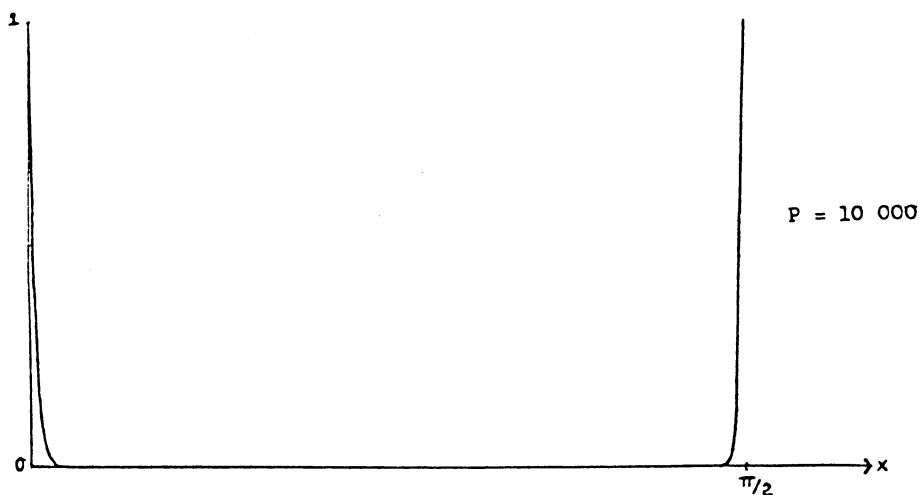| Stiffness	Degree of approximant	Perturbation term used	Nr. of equal subintervals	Max. Abs. Error
10^3	27	Chebyshev	1, no segm.	1.0×10^{-7}
10^3	14	Chebyshev	8	2.0×10^{-10}
10^3	14	Legendre	8	1.1×10^{-10}
10^4	28	Chebyshev	1, no segm.	1.0×10^{-2}
10^4	14	Chebyshev	12	7.8×10^{-7}
10^4	14	Legendre	12	4.4×10^{-7}

FIGURE 1

Graphs of the exact solution of the singularly perturbed boundary value problem (15), *for* $P = 10000$, *and of a segmented Tau approximant of degree* 14, *in the construction of which the technique of increased accuracy at matching points has been used. The maximum deviation between the two curves, seen as one in the graph, is* 4.4×10^{-7} (*see Table* 5).

(3) *Control of the Segmentation Process With* $[e_n(x)]_m$ *in an Adaptive Formulation of the Tau Method*. We now consider the singular perturbation boundary value problem with a boundary layer at $x = 1$ defined by the differential equation

(16)
$$\begin{cases} Dy(x) := -y''(x) + \frac{d}{dx}[P(1-cx)y(x)] = 0, \\ y(0) = 1, \quad y'(1) = 0, \qquad 0 \le x \le 1, \end{cases}$$

where $c = 0.98$ and $P = 1000$. The graph of the solution is a spike, symmetric about $x = 1$, extending in the y-axis direction from 1 up to about 50 when x is in $[0, 2]$. Problem (16) has been discussed numerically in a recent paper by Barrett and Morton [1], with a quasi-symmetrization technique based on Galerkin's method with a piecewise linear trial space; further references can be found in that paper.

We have constructed for this problem global Tau approximants of degrees 27 and 28, then a segmented piecewise Tau approximant of degree 7 defined over four equal

subintervals of $[0, 1]$. The Tau estimator $[e_n(x)]_m$, with $m = n + 1$, was used to detect the region of rapid variation of the last segmented approximant, and then a new nonuniform segmentation was introduced, with nodes in $x = 0.0$; 0.7; 0.8; 0.9 and 1.0. A segmented Tau piecewise approximant, of degree 7 over each subinterval, was constructed. In Table 6 we reproduce our numerical results; we have also included the values of the exact solution up to 2D. Agreement to 2D is provided by the Tau approximant.

TABLE 6

Use of the Tau estimator $[e_n(x)]_m$ for the generation of a nonuniform segmentation strategy, better adapted to the singular nature of the solution of the boundary layer problem (16); $m = n + 1$.

Type of approximant	Nr. of pieces	$x = 0.7$	$x = 0.8$	$x = 0.9$	$x = 1.0$
Global Tau, $n = 27$	1	3.10	4.73	9.32	49.99
Global Tau, $n = 28$	1	3.21	4.73	9.32	49.94
Segmented piecewise Tau, uniform segmentation, $n = 7$	4	3.22	4.48	8.40	48.21
Segmented piecewise Tau, adapted-nonuniform segmentation, controlled by the Tau estimator $[e_n(x)]_m$, $n = 7$	4	3.21	4.73	9.31	49.95
EXACT RESULTS		3.21	4.73	9.31	49.95

Results reported in [1] with quasi-symmetrization lead to large errors: $+0.14$ and -0.52, near $x = 1$ ($x = 0.8$ and 0.9 respectively). A comparison between the upwinding technique of [2] and Tau method approximants is also favorable to the recursive formulation of the Tau method (see Ortiz [29]).

(4) *Tau Approximate Solution and Error Estimation of a Nonlinear Boundary Value Problem With Global and Segmented Approximants.* The numerical solution of nonlinear differential equations with the recursive formulation of the Tau method is reduced to the approximate solution of a sequence of linear problems, as indicated in Section 6, that is, to problems of the type considered before. However, we shall discuss explicitly a nonlinear problem with a singularity of the first kind proposed by Russell and Shampine [34], for which results with the finite difference techniques of de Hoog and Weiss [5] are available,

$$(17) \qquad \frac{d}{dx}\begin{bmatrix} y_1(x) \\ -y_2(x) \end{bmatrix} = \frac{1}{x}\begin{bmatrix} 0 & 0 \\ 0 & 2 \end{bmatrix}\begin{bmatrix} y_1(x) \\ y_2(x) \end{bmatrix} + \begin{bmatrix} y_2(x) \\ y_1^5(x) \end{bmatrix};$$

$$\begin{bmatrix} y_1(1) \\ y_2(0) \end{bmatrix} = \begin{bmatrix} \sqrt{.75} \\ 0 \end{bmatrix}, \qquad 0 \leqslant x \leqslant 1,$$

which we treat as the second order boundary value problem

$$(18) \qquad \begin{cases} Dy(x) := y''(x) + (2/x)y'(x) + y^5(x) = 0, \\ y'(0) = 0, \quad y(1) = \sqrt{0.75}, \qquad 0 \leqslant x \leqslant 1. \end{cases}$$

A direct approximation of systems of differential equations with the recursive formulation of the Tau method is also possible if *vector* canonical polynomials are used; see [6] for details on this approach. The exact solution of (17)–(18) is known to be $y(x) = (1 + x^2/3)^{-1/2}$, which makes it possible to compare the error of our Tau approximations of $y(x)$ with the estimation of that error provided by the Tau estimator $[e_n(x)]_m$.

We fix the tolerance parameter T (see Section 3) equal to 10^{-6} and the maximum permissible number of cycles N equal to 5 which turns out to be sufficiently large. Table 7 displays the errors of approximation of a global and a segmented Tau approximant, the latter over the subintervals $[0, 0.5]$ and $[0.5, 1]$.

TABLE 7

Maximum absolute error of approximation for the nonlinear boundary value problem (18) when approximated with global and segmented Tau approximants.

Degree of the Tau approximant	Global Approximation		Segmented approximation	
	Max. Abs. Error	Nr. of cycles required	Max. Abs. Error	Nr. of cycles required
4	8.7×10^{-4}	3	7.8×10^{-5}	3.3
5	6.0×10^{-5}	3	1.6×10^{-6}	3.3
6	6.6×10^{-6}	3	1.0×10^{-6}	3.3

The Tau error estimator $[e_n(x)]_m$, applied to both global and segmented approximants, leads to remarkably accurate estimations of the exact error. Table 8 reproduces results for $n = 4$, $m = n + 1$ and $n + 2$, with global and segmented Tau approximants. For $m = n + 1$ the order of the error is correctly estimated in both cases.

TABLE 8

Error estimation of global and segmented Tau approximants of the nonlinear singular boundary value problem (18), given by the Tau estimator $[e_n(x)]_m$ for $n = 4$.

Global Tau approximant		Segmented Tau approximants	
m	Difference between the exact and estimated errors	m, m	Difference between the exact and estimated errors
$n + 1$	6.01×10^{-5}	$n + 1, n + 1$	1.57×10^{-6}
$n + 2$	6.50×10^{-6}	$n + 2, n + 2$	1.05×10^{-6}

The nonlinear system of differential equations

$$\frac{d}{dx}\begin{bmatrix} y_1(x) \\ y_2(x) \end{bmatrix} = \begin{bmatrix} y_2(x) \\ (1/\varepsilon)[y_1^2(x) - x^2]\,y_2(x) \end{bmatrix}, \qquad \begin{bmatrix} y_1(-1) \\ y_1(0) \end{bmatrix} = \begin{bmatrix} 0.96 \\ 0.001 \end{bmatrix},$$

$-1 \leqslant x \leqslant 0$, is a nonlinear model of a singularly perturbed differential equation where the leading coefficient of the reduced equation has turning points, the location of which depends on the unknown function. It is a stiff problem with a multiplicity

of solutions which has been treated by Kedem [9] by collocation ($\varepsilon = 15$), and by Ortiz [29] with the Tau method (global approximant); further references can be found in these papers. High-order boundary value problems for differential equations are considered in [21] (global approximant). Further numerical results on a large variety of problems where the technique discussed in this paper has been successfully applied are reported in Onumanyi [19]-[20].

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