

Numerical Study on Phase-Fitted and Amplification-Fitted Diagonally Implicit Two Derivative Runge-Kutta Method for Periodic IVPs

(Kajian Berangka ke atas Suai-Fasa dan Suai-Pembesaran Kaedah Dua-terbitan Pepenjuru Tersirat Kaedah Runge-Kutta untuk MNA Berkala)

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ABSTRACT

A fourth-order two stage Phase-fitted and Amplification-fitted Diagonally Implicit Two Derivative Runge-Kutta method (PFAFDITDRK) for the numerical integration of first-order Initial Value Problems (IVPs) which exhibits periodic solutions are constructed. The Phase-Fitted and Amplification-Fitted property are discussed thoroughly in this paper. The stability of the method proposed are also given herewith. Runge-Kutta (RK) methods of the similar property are chosen in the literature for the purpose of comparison by carrying out numerical experiments to justify the accuracy and the effectiveness of the derived method.

Keywords: Diagonally implicit methods; initial values problems; ordinary differential equations; phase-fitted and amplification-fitted; stability region; two derivative Runge-Kutta method

ABSTRAK

Kaedah Runge-Kutta Dua Terbitan Pepenjuru Tersirat Suai-Fasa dan Suai-Pembesaran (RKDTPTSFSP) tahap dua peringkat empat untuk penyelesaian pengamiran berangka Masalah Nilai Awal (MNA) peringkat pertama yang mengandungi penyelesaian berkala dibina. Sifat suai-fasa dan suai-pembesaran dibincangkan secara menyeluruh dalam kertas kajian ini. Kestabilan kaedah yang dicadangkan adalah seperti berikut. Kaedah Runge-Kutta (RK) dengan sifat yang sama dipilih di dalam kajian sorotan untuk tujuan perbandingan dengan menjalankan uji kaji berangka untuk memastikan kejituhan dan keberkesanan kaedah yang diterbitkan.

Kata kunci: Kaedah pepenjuru tersirat; kaedah Runge-Kutta dua terbitan; masalah nilai awal; persamaan pembezaan biasa; rantau kestabilan; suai-fasa dan suai-pembesaran

INTRODUCTION

The Ordinary Differential Equations (ODEs) of first-order for the numerical solution of the IVPs are considered

$$q' = f(t, q), \text{ given the initial condition, } q(t) = q_0, \quad (1)$$

where their solutions show periodically or oscillatory behavior in which the eigenvalue is in complex form. This type of problems appears throughout certain fields of applied sciences, for instance, mechanics, electronics, circuit simulation, orbital mechanics, astrophysics, and

molecular dynamics. In general, periodically or oscillatory behavior problems are mostly known with second or higher order. It is therefore essential to perform order reduction to solve the ODEs (1) by reducing them to first-order problems.

Anastassi and Simos (2012), Chen et al. (2012), and Kosti et al. (2012a) efficiently solved the Schrödinger equation and related periodically problems by designing a new explicit phase-fitted and amplification-fitted for the optimization of the method.

RK methods for solving oscillatory problems using several techniques, for instance, phase-fitted and amplification-fitted, trigonometrically-fitted and exponentially-fitted techniques have been developed and expanded by several famous authors such as Simos (1998) in his written paper. Simos (1998) designed a Runge-Kutta method with exponentially-fitted properties for the numerical integration of IVPs of order five. Konguetsof and Simos (2003) introduced explicit symmetric multistep method which is exponentially-fitted and trigonometrically-fitted of eighth-order.

Recently, Adel et al. (2016) and Fawzi et al. (2015) derived two fourth-order modified RK and classical RK method with phase-fitted and amplification-fitted property, respectively. Meanwhile, Demba et al. (2016a, 2016b) suggested Runge-Kutta-Nyström (RKN) methods with trigonometrically-fitted property to solve second-order IVPs with periodic solutions in nature derived on Simos' RKN method. Two Derivative Runge-Kutta (TDRK) methods which are explicit in nature given by Chan and Tsai (2010) in which they include the second derivative in its general formula making it special. Just one evaluation of function f is involved along with a several number of function g to be evaluated at every step. With this finding, they managed to derive methods up to order seven with five stages as well as some embedded pairs.

The numerical integration of radial Schrödinger equation and periodic problems are constructed by Zhang et al. (2013) using a TDRK method with trigonometrically-fitted of order five. Other than that, Fang et al. (2013) and Chen et al. (2015) constructed two TDRK methods of order four and three practical TDRK methods with exponentially-fitted, respectively. The newly derived methods are compared with some widely-known optimized codes as well as conventional RK methods with exponentially-fitted property mentioned in the literature.

In this recent year, there are no findings of research associated with phase-fitting and amplification-fitting in DITDRK methods. The benefits or drawbacks of applying phase-fitted and amplification-fitted property to DITDRK methods have not yet discussed thoroughly by researchers especially mathematicians. A two stage fourth-order DITDRK method with phase-fitted and amplification-fitted property is therefore derived in this paper. A summary of the TDRK method is discussed in Section 2. The next section considered the conditions for the phase-fitted and amplification-fitted property. The construction of the phase-fitted and amplification-fitted

DITDRK method is defined in Section 4. A description on the stability property is discussed briefly in Section 5. The numerical results, discussion, and conclusion are presented briefly in Sections 6, 7, and 8, respectively.

TWO DERIVATIVE RUNGE-KUTTA METHODS

The scalar ODEs (1) is considered with $g: \Re^N \rightarrow \Re^N$. It is assumed, in this case, the second derivative is known where

$$q'' = g(q) := f'(q)f(q), g: \Re^N \rightarrow \Re^N. \quad (2)$$

The numerical integration of IVPs (1) for a TDRK method is given by

$$q_{n+1} = q_n + \Delta t \sum_{i=1}^s b_i f(q_i) + \Delta t^2 \sum_{i=1}^s \hat{b}_i g(Q_i), \quad (3)$$

$$Q_i = q_n + \Delta t \sum_{j=1}^s a_{ij} f(q_j) + \Delta t^2 \sum_{j=1}^s \hat{a}_{ij} g(Q_j), \quad (4)$$

The lowest number of function evaluations for diagonally implicit methods can be established by considering the methods in the following manner where $i = 1, \dots, s$.

$$q_{n+1} = q_n + \Delta t f(t_n, q_n) + \Delta t^2 \sum_{i=1}^s \hat{b}_i g(t_n + \Delta t c_i, Q_i), \quad (5)$$

$$Q_i = q_n + \Delta t c_i f(t_n, q_n) + \Delta t^2 \sum_{j=1}^i \hat{a}_{ij} g(t_n + \Delta t c_j, Q_j), \quad (6)$$

where $i = 1, \dots, s$.

Assume that all of the following DITDRK parameters a_{ij} , \hat{a}_{ij} , b_i , \hat{b}_i and c_i are real and s is method's stages number. We introduced the -dimensional vectors $b = [b_1, b_2, \dots, b_s]^T$, $\hat{b} = [\hat{b}_1, \hat{b}_2, \dots, \hat{b}_s]^T$, $c = [c_1, c_2, \dots, c_s]^T$ and $s \times s$ matrices $A = [a_{ij}]$ and $\hat{A} = [\hat{a}_{ij}]$ where $1 \leq i, j \leq s$. We use the following simplifying assumption,

$$\sum_{j=1}^s \hat{a}_{ij} = \frac{1}{2} c_i^2, \quad (7)$$

Table 1 shows the order conditions for unique DITDRK methods given in Chan and Tsai (2010).

TABLE 1. Order conditions for unique DITDRK methods

Order	Conditions
1	$b^T e = 1$
2	$\hat{b}^T e = \frac{1}{2}$
3	$\hat{b}^T c = \frac{1}{6}$
4	$\hat{b}^T c^2 = \frac{1}{12}$
5	$\hat{b}^T c^3 = \frac{1}{20}$
	$\hat{b}^T \hat{A}c = \frac{1}{120}$
6	$\hat{b}^T c^4 = \frac{1}{30}$
	$\hat{b}^T c \hat{A}c = \frac{1}{180}$
	$\hat{b}^T \hat{A}c^2 = \frac{1}{360}$
7	$\hat{b}^T c^5 = \frac{1}{42}$
	$\hat{b}^T c^2 \hat{A}c = \frac{1}{252}$
	$\hat{b}^T c \hat{A}c^2 = \frac{1}{504}$
	$\hat{b}^T \hat{A}c^3 = \frac{1}{840}$
	$\hat{b}^T \hat{A}^2 c = \frac{1}{5040}$

The method described herewith is identified as a unique DITDRK method. The remarkable aspect of this method is it requires just one evaluation of function f and a few evaluations of function g per step compared to a number of evaluations of function f per step in the conventional RK methods. The following Butcher tableau illustrates the significant difference between the DITDRK method and the unique DITDRK method.

$$\begin{array}{c|cc||c|c} c & A & \hat{A} \\ \hline b^T & \hat{b}^T \end{array} \longrightarrow \begin{array}{c||c} c & \hat{A} \\ \hline \hat{b}^T & \end{array}$$

PHASE-FITTED AND AMPLIFICATION-FITTED PROPERTY
The following linear scalar equation is considered,

$$q' = i\lambda q. \quad (8)$$

The exact solution with initial value $q(t_0) = q_0$ of this equation satisfies

$$q(t_0 + \Delta t) = H_0(z)q_0, \quad (9)$$

where $H_0(z) = \exp(z)$, $z = iv$. A phase advance $v = \lambda\Delta t$ is experienced by the exact solution whereby the amplification appears to remain stable and secure after a cycle of time Δt .

The DITDRK method is adapted to the test equation (8) to yield

$$q_1 = H(z)q_0, \quad (10)$$

where

$$H(z) = \left(1 + v^2 \hat{b}(I - v^2 \hat{A})^{-1} e\right) + i \left(v + v^3 \hat{b}(I - v^2 \hat{A})^{-1} c\right), \quad (11)$$

where $e = [1, \dots, 1]^T$.

The stability function of the DITDRK method is presented by $H(z)$ which in term of complex number. The function is split in terms of the real (denoted as $U(v)$) and imaginary (denoted as $V(v)$) part of $H(z)$. Further, we have the argument of $H(z)$ or simply $\arg H(z) = \tan^{-1}\left(\frac{V(v)}{U(v)}\right)$ and the magnitude of $H(z)$ or $|H(z)| = \sqrt{U^2(v) + V^2(v)}$ for small Δt . According to the analysis above, the following definition arises.

Definition 1 (van der Houwen & Sommeijer (1987))
The quantities

$$\tilde{P}(v) = v - \arg H(z), \tilde{D}(v) = 1 - |H(z)|, \quad (12)$$

are defined as the phase lag (or dispersion) and the error of amplification factor (or dissipation) of the method, respectively. If

$$\tilde{P}(v) = c_\phi v^{\alpha+1} + \mathcal{O}(v^{\alpha+3}), \tilde{D}(v) = c_d v^{\beta+1} + \mathcal{O}(v^{\beta+3}), \quad (13)$$

then, the method is defined as dispersive of order α and dissipative of order β , respectively.

If

$$\tilde{P}(v) = 0, \tilde{D}(v) = 0, \quad (14)$$

the method is defined as phase-fitted (or zero-dispersive) and amplification-fitted (or zero dissipative), respectively.

Theorem 2 (Chen et al. 2012)

The method is justified as phase-fitted and amplification-fitted if and only if

$$U(v) = \cos(v), V(v) = \sin(v). \quad (15)$$

The local truncation error (LTE), $LTE = q(t_0 + \Delta t) = \mathcal{O}(\Delta t^{\zeta+1})$, for any $(\zeta + 1)$ -th differentiable function $g(q)$, when equations (5) and (6) are applied to the first-order ODEs (1). Thence, the method is said to have a (algebraic) order ζ . Define

$$EC_{\zeta+1}(v) = \left(\sum_{i=1}^j (\tau_j^{(\zeta+1)})^2 \right)^{\frac{1}{2}} \quad (16)$$

where $\tau_j^{(\zeta+1)}$ is the error coefficient of the method. The non-negative number

$$EC_{\zeta+1} = \lim_{v \rightarrow 0} EC_{\zeta+1}(v), \quad (17)$$

is known as the method's error constant.

DERIVATION OF THE NEW PHASE-FITTED AND AMPLIFICATION-FITTED METHOD

If and only if Theorem 2 is satisfied, then only a DITDRK method appeared to be phase-fitted and amplification-fitted. Thus, the proposed method is derived by combining the DITDRK method with the phase-fitted and amplification-fitted property proposed in this section.

First, a fourth-order two stages DITDRK method will be derived. Referring to the order conditions in Table 1 up to fourth-order, we have

$$\hat{b}_1 + \hat{b}_2 - \frac{1}{2} = 0, \quad (18)$$

$$\hat{b}_1 c_1 + \hat{b}_2 c_2 - \frac{1}{6} = 0, \quad (19)$$

$$\hat{b}_1 c_1^2 + \hat{b}_2 c_2^2 - \frac{1}{12} = 0. \quad (20)$$

Solving equation (18)-(20) we obtain \hat{b}_1, \hat{b}_2 and c_1 in term of c_2

$$\hat{b}_1 = \frac{1}{(36c_1^2 - 24c_1 + 6)}, \quad (21)$$

$$\hat{b}_2 = \frac{1}{3} \left(\frac{9c_1^2 - 6c_1 + 1}{6c_1^2 - 4c_1 + 1} \right), \quad (22)$$

$$c_2 = \frac{1}{2} \left(\frac{2c_1 - 1}{3c_1 - 1} \right). \quad (23)$$

Our main focus is to choose c_1 in such a way that a very small value of the principal local truncation error coefficient, $\|\tau^{(5)}\|_2$ might be achieved. There will be a significant global error difference with an inaccurate choice of c_1 . The graph of $\|\tau^{(5)}\|_2$ against c_1 is plotted in Figure 1 where a small value of c_1 is chosen within the range of $[0.0, 1.0]$. Therefore, the value of c_1 is between $[0.1, 0.3]$ with the help of Maple software where we use the minimisation command for non-linear functions. For simplicity, we have chosen $c_1 = \frac{1}{5}$ for an ideal optimized pair by running empirical experiment.

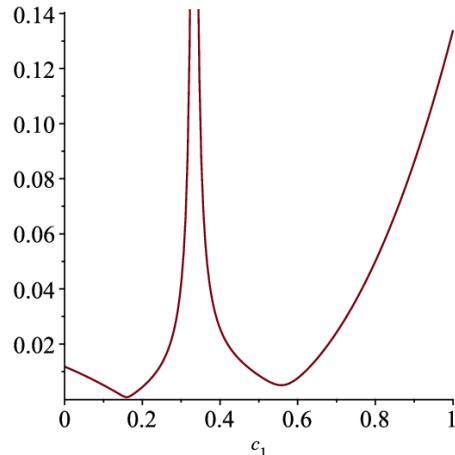


FIGURE 1. The graph of $\|\tau^{(5)}\|_2$ against c_1

The following Butcher tableau represents the coefficients

of the method and are referred to as DITDRK(2,4).

TABLE 2. Butcher Tableau for DITDRK(2,4) Method

1	1
5	$\frac{50}{209}$
3	$\frac{1}{209}$
$\frac{1}{4}$	$\frac{1}{50}$
	$\frac{25}{800}$
	$\frac{4}{50}$
	$\frac{25}{66}$
	$\frac{4}{33}$

The stability function (11) for two stages fourth-order DITDRK method is considered. Therefore, by choosing \hat{a}_{21} and c_1 as free parameters, we have

$$I = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix}, e = \begin{bmatrix} 1 \\ 1 \end{bmatrix}, \hat{b} = \begin{bmatrix} \frac{25}{66} \\ \frac{4}{33} \end{bmatrix}, c = \begin{bmatrix} c_1 \\ \frac{1}{4} \end{bmatrix}, \hat{A} = \begin{bmatrix} \frac{1}{50} & \frac{1}{50} \\ \hat{a}_{21} & \frac{1}{50} \end{bmatrix}. \quad (24)$$

We substituted the matrices (24) into $H(z)$ given by equation (11) and splitted the complex number of $H(z)$ into real and imaginary as mentioned. The free parameters, \hat{a}_{21} and c_1 are taken as the ideal combination for the optimized value of the maximum global error. By implementing Theorem 2, (15) is solved to get the coefficients of \hat{b}_1 and c_4 and this resulting in

$$\cos(v) = \frac{1}{33} \frac{10000 v^4 \hat{a}_{21} - 792 v^4 - 37950 v^2 + 82500}{(v^2 + 50)^2}, \quad (25)$$

$\sin(v)$

$$= \frac{1}{33} \frac{v(10000 v^4 \hat{a}_{21} c_1 - 625 v^4 c_1 - 117 v^4 - 31250 v^2 c_1 - 4200 v^2 + 82500)}{(v^2 + 50)^2}. \quad (26)$$

By solving (25) and (26) we will obtain the following

$$\hat{a}_{21} = \frac{(33 \cos(v) + 792)v^4 + (3300 \cos(v) + 37950)v^2 + 82500 \cos(v) - 82500}{10000 v^4}, \quad (27)$$

$$c_1 = \frac{33 \sin(v) v^2 + 117 v^3 + 1650 \sin(v) - 1650 v}{33 \cos(v) v^3 + 167 v^3 + 1650 \cos(v) v - 1650 v}. \quad (28)$$

The following Taylor expansions as $v \rightarrow 0$ are obtained as follows

$$\begin{aligned} \hat{a}_{21} &= \frac{209}{800} + \frac{77 v^2}{120000} - \frac{781 v^4}{6720000} + \frac{803 v^6}{604800000} + \frac{59 v^8}{7257600000} - \frac{2081 v^{10}}{6604416000000} \\ &\quad + \dots, \\ c_1 &= \frac{1}{5} + \frac{11 v^2}{3125} + \frac{3476 v^4}{41015625} + \frac{13111549 v^6}{3691406250000} + \frac{1679796241 v^8}{9228515625000000} \\ &\quad + \frac{66016889468987 v^{10}}{6298461914062500000000} + \dots. \end{aligned}$$

The following expansions shall be obtained by direct calculation:

$$\begin{aligned} \hat{b}^T e &= \frac{1}{2}, \\ \hat{b}^T c &= -\frac{5}{66} + \frac{825 \sin(v) v^2 + 2925 v^3 + 41250 \sin(v) - 41250 v}{2178 \cos(v) v^3 + 11022 v^3 + 108900 \cos(v) v - 108900 v} \quad (29) \\ &= \frac{1}{6} + \mathcal{O}(v), \\ \hat{b}^T c^2 &= -\frac{1}{66} + \frac{25}{66} \left(\frac{M}{N}\right)^2 = \frac{1}{12} + \mathcal{O}(v). \end{aligned}$$

Subsequently, all of the order conditions till order four are satisfied by the coefficients shown in Table 2. But the condition for order five was not satisfied. For instance,

$$\hat{b}^T c^3 = 0 \neq \frac{1}{20} + \mathcal{O}(v). \quad (30)$$

Therefore, it is a method of order four. The coefficients of error of the DITDRK(2,4) for order five are given by

$$\tau_1^{(5)} = \frac{1}{240}, \quad \tau_2^{(5)} = \frac{1}{750}. \quad (31)$$

Therefore, for DITDRK(2,4), we obtain the following

$$EC_5 = \frac{1}{6000} \sqrt{689}. \quad (32)$$

As we have proven that this newly derived method is fourth-order, it is therefore known as PFAFDITDRK(2,4). The error coefficients of PFAFDITDRK(2,4) are given by

$$\begin{aligned} \tau_1^{(5)} &= \frac{1}{880} + \frac{25}{66} \left(\frac{M}{N} \right)^3, \\ \tau_2^{(5)} &= \left(\frac{1}{132} + \frac{1}{82500 v^4} \right) \left(\frac{MP}{N} \right) - \frac{43}{6600}. \end{aligned} \quad (33)$$

For PFAFDITDRK(2,4), we have

$$\begin{aligned} EC_5(v) &= \left(\frac{1}{108900000000 N^6} (15625000000 M^6 + 93750000 (MN)^3 \right. \\ &\quad + 6250000 (MN^2)^2 - 10750000 MN^5 + 4763125 N^6) \\ &\quad \left. + \frac{1}{108900000000 N^6 v^4} (20000 (MN^2)^2 P - 17200 MN^5 P) \right. \\ &\quad \left. + \frac{1}{6806250000 v^8} \left(\frac{MP}{N} \right)^{\frac{1}{2}} \right), \end{aligned} \quad (34)$$

where $M = 33(v) v^2 + 117 v^3 + 1650 \sin(v) - 1650 v$, $N = 33 \cos(v) v^3 + 167 v^3 + 1650 \cos(v) v - 1650 v$, $P = 33 \cos(v) v^4 + 792 v^4 + 3300 \cos(v) v^2 + 37950 v^2 + 82500 \cos(v) - 82500$.

PFAFDITDRK(2,4) will reduce to its actual method, DITDRK(2,4) as $v \rightarrow 0$. Apart from that, PFAFDITDRK(2,4)

will have the identical error constant as DITDRK(2,4) as $v \rightarrow 0$.

STABILITY AND CONVERGENCE OF THE NEW METHOD

The linear stability of the method being developed is analysed in this section. Applying equation (8) to the DITDRK method produces the difference equation

$$q_{n+1} = H(z)q_n, z = iv, i^2 = -1, \quad (35)$$

where $H(z)$ is given as (11).

Definition 3 A DITDRK method is said to be absolutely stable if $|H(z)| < 1$ for all $z \in (-v, 0)$. The stability polynomial of the PFAFDITDRK(2,4) method is shown as follows.

$$\begin{aligned} H(v) &= \frac{1}{49883818359375000000000 (v^2 - 50)^2} (56936760886197823 v^{15} \\ &\quad - 4763031005859375 v^{14} + 1022777450921523750 v^{13} + \\ &\quad 122886657714843750 v^{12} + 17921737015284375000 v^{11} + \\ &\quad 20070098876953125000 v^{10} + 105179969425781250000 v^9 \\ &\quad - 1756820983886718750000 v^8 + \\ &\quad 9699631347656250000000 v^6 + 590291850585937500000000 v^5 + \\ &\quad 2751923979492187500000000 v^4 + 15796542480468750000000000 v^3 + \\ &\quad 5736639111328125000000000 v^2 + 124709545898437500000000000 v + \\ &\quad 124709545898437500000000000 + \dots). \end{aligned} \quad (36)$$

We plot and compare the region of stability of the PFAFDITDRK(2,4) method up to $v^i, i = 6, 8, 14$ and its actual method as in Figure 2.

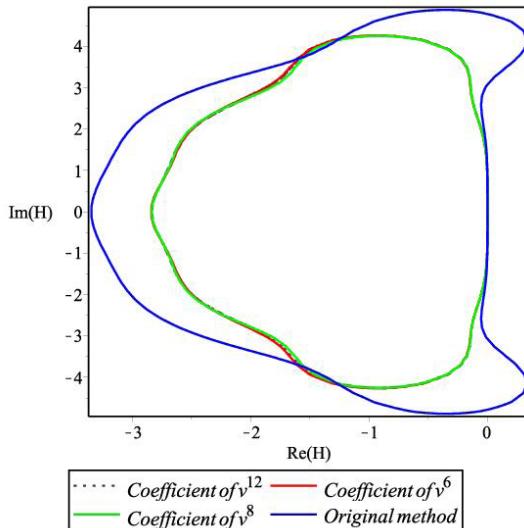


FIGURE 2. Stability region of PFAFDITDRK(2,4) method for different order

The stability interval with the coefficients v^6 , v^8 and v^{12} of this method are (-2.843,0.000), (-2.837,0.000) and (-2.833,0.000), respectively. The stability regions in Figure 2 is observed and as the coefficients order tends to infinity, the stability interval becomes further away from the original method where it is given by (-3.347,0.000).

Through the stability interval, we can literally consider the largest value of Δt the method could take to ensure it will remain stable. $v = \lambda \Delta t$ is mentioned earlier and the test problems represents the value of λ . Therefore, the value of Δt is obtained by dividing v with λ . The stability test as following would illustrate on how the regions of stability are used for practical purposes. We have

$$q' = \lambda(q - \varphi) + \varphi', q(0) = \varphi(0), Re(\lambda) < 0, t \in [0, 2000],$$

given that $\varphi(t)$ is a smooth function. Letting $\lambda = -1$, $\varphi(t) = \sin(t)$ and $q(t) = \varphi(t)$ is the exact solution.

Stability of the method can be achieved once the maximum global error is small enough and therefore converging to its exact solution. Instead of that, a larger maximum global error indicates that the method is unstable, meaning that they are actually diverging from its exact solution. The stability test will be conducted to demonstrate the connection between Δt , λ and $|H(z)|$. When $\Delta t = 4.15$, the stability is achieved whereby this is the largest value of Δt can be used to ensure the method remain stable in this particular test for stability. Table 3 represents the global error for a variety of Δt values.

TABLE 3. Stability test for PFAFDITDRK(2,4) using coefficient of v^8 with $\lambda = -1$ for variable Δt

Δt	$ H(z) $	Global Error
3.20	2.399948118	1.729241×10^{236}
3.00	1.496894148	5.112085×10^{114}
2.83	0.9808945284	4.720623×10^0
1.00	0.3650531765	$1.836955 \times 10^{(-3)}$
0.15	0.8607077753	$5.005154 \times 10^{(-7)}$
0.01	0.9900498337	$9.473644 \times 10^{(-12)}$

Definition 4 (Henrici, 1962)

The numerical method with order p is zero stable if numerical solutions remain bounded in the limit $\Delta t \rightarrow 0$, with the modulus of roots for the first characteristic polynomial are less than or equal to zero.

In studying the zero stability of the DITDRK method, the characteristic polynomial of method (5)-(6) is:

$$p(\xi) = (\xi - 1) \quad (37)$$

Hence, the method is zero stable since the roots, $\xi = 1$ are less than or equal to one.

Definition 5 (Suli & Mayers 2003)

The method is consistent with the order at least p if and only if local truncational error, $T_{p+1} = \mathcal{O}(\Delta t^{p+1})$ as $\Delta t \rightarrow 0$. Consider DITDRK methods in the class as follow:

$$\sum_{j=0}^s \delta_j q_{n+j} = \Delta t \gamma_j f_{n+j} + \Delta t^2 \phi g(q_{n+k}, q_{n+k-1}, \dots, q_n, t_n; \Delta t^2). \quad (38)$$

On putting $s = 1$, then

$$\delta_1 = 1, \delta_0 = -1, \gamma_0 = 1, \quad \phi g(q_n, t_n; \Delta t^2) = \sum_{i=1}^s \hat{b}_i Q_i, \quad (39)$$

$$Q_i = q_n + \Delta t c_i f(t_n, q_n) + \Delta t^2 \sum_{j=1}^i \hat{a}_{ij} g(t_n + \Delta t c_j, Q_j),$$

where $i = 1, \dots, s$.

The condition for (39) to be consistent are

$$\begin{aligned} \sum_{j=0}^s \delta_j &= 0, \quad \sum_{j=0}^s (j\delta_j - \gamma_j) = 0, \\ \frac{\phi g(q(t_n), q(t_n), \dots, q(t_n), t_n; 0)}{\sum_{j=0}^s j\delta_j} &= g(t_n, q(t_n)). \end{aligned} \quad (40)$$

Applying the conditions (40), the necessary and sufficient condition for DITDRK methods to acquire consistency is

$$\phi g(q(t_n), t_n; 0) = g(t_n, q(t_n)) \Leftrightarrow \sum_{i=0}^s \hat{b}_i = \frac{1}{2}. \quad (41)$$

Here, local truncation error, T_{n+1} at t_{n+1} is expressed as the residual when q_{n+j} is replaced by $q(t_{n+j})$ which is

$$\begin{aligned} T_{n+1} &= q(t_{n+1}) + \Delta t f(t_n, q(t_n)) - [q(t_n) + \Delta t f(t_{n+1}, q(t_{n+1}))] \\ &\quad - \Delta t^2 \phi g(q(t_n), t_n; \Delta t^2), \end{aligned} \quad (42)$$

where ϕg is defined in (39). Assuming that p is the largest integer whereby $T_{n+1} = \mathcal{O}(\Delta t^{p+1})$, then the method has order p (Lambert 1991). We denote by \tilde{q}_{n+1} the value at t_{n+1} generated by DITDRK method when the localising assumption, $q_n = q(t_n)$ is made. Since

$$\tilde{q}_{n+1} = q_n + \Delta t f(t_n, q_n) + \Delta t^2 \phi g(q_n, t_n; \Delta t^2). \quad (43)$$

Then we have

$$q(t_{n+1}) - \tilde{q}_{n+1} = T_{n+1}. \quad (43)$$

DITDRK method is consistent if they follow (41) such that

$$\begin{aligned} q(t_{n+1}) + \Delta t f(t_n, q(t_n)) - [q(t_n) + \Delta t f(t_{n+1}, q(t_{n+1}))] \\ = \frac{\Delta t^2}{2} f'(t_n) - \frac{\Delta t^2}{2} g(t_n, q(t_n)) + \mathcal{O}(\Delta t^3) \end{aligned} \quad (44)$$

By reason of $f'(t_n) = g(t_n, q(t_n))$, T_{n+1} for DITDRK method is equal to $\mathcal{O}(\Delta t^3)$, it shows that DITDRK method is consistent if their order is at least 2, which is in line with our definitions of order for linear multistep methods. Since the order of DITDRK method is at least 2, and hence, this method is consistent.

Convergence is a property of numerical method related to truncation errors that ensures the numerical solution converges onto the exact solution and the global truncation error goes to zero at all step size indices in

the limit $\Delta t \rightarrow 0$ (Atkinson 2009). Maximum absolute global truncation error between the analytical solution and numerical solution the gets smaller as the step size becomes lesser.

Definition 6 (Lambert 1991)

The numerical method is convergent if acquiring the properties of zero stability and consistency.

Since DITDRK method is zero-stable and consistent, implies that DITDRK method is convergent.

PROBLEMS TESTED AND NUMERICAL RESULTS

The derived method PFAFDITDRK(2,4) are compared in term of their numerical performances with some famous existing RK and TDRK methods by considering Problems 1-5 as follows. C Programming codes are used for solving differential equations where all the problems chosen are having oscillatory solutions.

Problem 1 (Harmonic Oscillator)

$$\begin{aligned} q_1'(t) &= q_2(t), \quad q_1(0) = q_{0_1}, \quad t \in [0, t_{end}], \\ q_2'(t) &= -\omega^2 q_1(t), \quad q_2(0) = q_{0_2}. \end{aligned}$$

Exact solution is

$$q_1(t) = \bar{c}_1 \sin(\omega t) + \bar{c}_2 \cos(\omega t),$$

$$q_2(t) = \bar{c}_3 \omega \cos(\omega t) - \bar{c}_4 \omega \sin(\omega t).$$

Total energy as given in Pokorny (2009)

$$E(q_1, q_2) = \frac{q_1^2}{2} + \frac{q_2^2}{2} = \frac{\Psi^2}{2},$$

where Ψ depends on the initial conditions.

Problem 2 (Inhomogeneous problem (Van de Vyver 2007))

$$q_1' = q_2, \quad q_1(0) = 1, \quad t \in [0, 1000],$$

$$q_2' = -100q_1 + 99 \sin(t), \quad q_2(0) = 11.$$

Exact solution is

$$q_1(t) = \cos(10t) + \sin(10t) + \sin(t),$$

$$q_2(t) = -10 \sin(10t) + 10 \cos(10t) + \cos(t).$$

Problem 3 (An almost Periodic Orbit problem (Stiefel & Bettis 1969))

$$\begin{aligned} q_1' &= q_2, \quad q_1(0) = 1, \quad t \in [0, 1000], \\ q_2' &= -q_1 + 0.001 \cos(t), \quad q_2(0) = 0, \\ q_3' &= q_4, \quad q_3(0) = 0, \\ q_4' &= -q_3 + 0.001 \sin(t), \quad q_4(0) = 0.9995. \end{aligned}$$

Exact solution is

$$\begin{aligned} q_1(t) &= \cos(t) + 0.0005t \sin(t), \quad q_2(t) = -\sin(t) + 0.0005t \cos(t) + 0.0005t \sin(t), \\ q_3(t) &= \sin(t) - 0.0005t \cos(t), \quad q_4(t) = \cos(t) + 0.0005t \sin(t) - 0.0005 \cos(t). \end{aligned}$$

Problem 4 (Duffing problem (Kosti et al. 2012b)

$$\begin{aligned} q_1' &= q_2, \quad q_1(0) = 0.200426728067, \\ q_2' &= -q_1 - q_1^3 + 0.002 \cos(1.01t), \quad q_2(0) = 0, \quad t \in [0, 1000]. \end{aligned}$$

Exact solution is

$$q_1(t) = 0.200179477536 \cos(1.01t) + 2.46946143 \times 10^{-4} \cos(3.03t) + 3.04014 \times 10^{-7} \cos(5.05t) + 3.74 \times 10^{-10} \cos(7.07t),$$

$$\begin{aligned} q_2(t) &= -0.2021812723 \sin(1.01t) - 7.482468133 \times 10^{-4} \sin(3.03t) - 1.53527070 \times 10^{-6} \sin(5.05t) - 2.64418 \times 10^{-9} \sin(7.07t). \end{aligned}$$

Problem 5 (Prothero-Robinson problem Chan & Tsai 2010)

$$q' = \lambda(q - \varphi) + \varphi', \quad q(0) = \varphi(0), \quad Re(\lambda) < 0, \quad t \in [0, 1000],$$

where $\varphi(t)$ is a smooth function and taking $\lambda = -1$, $\varphi(t) = \sin(t)$.

Exact solution is $q(t) = \varphi(t)$.

Figures 3-18 used the following abbreviations.

PFAFDITDRK(2,4): Fourth-order two stages phase-fitted and amplification-fitted DITDRK method proposed in this paper. TFDIRKK(3,4): Fourth-order three stages trigonometrically-fitted DIRK method developed in Kalogiratou (2013). PFAFDIRKA(3,4): Fourth-order three stages phase-fitted and amplification-fitted DIRK method given by Ahmad et al. (2016). EFDIRKE(3,4): Fourth-order three stages exponentially-fitted DIRK method given in Ehigie et al. (2018).

Figures 3-18 represents the behaviour of these numerical results in graphics form.

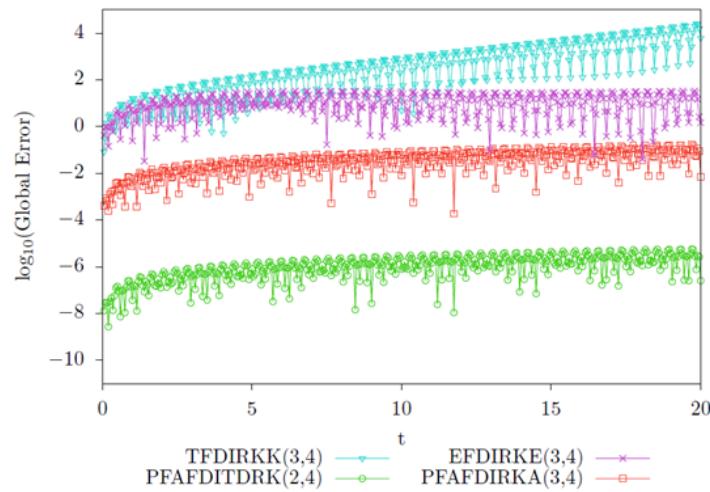


FIGURE 3. (Conservation of Energy). The logarithm error of energy (Global Error) when solving the harmonic oscillator (Problem 1) at each integration point for $\omega = 8$, $q_{01} = 1$, $q_{02} = -2$ and $\Delta t = 1/20$

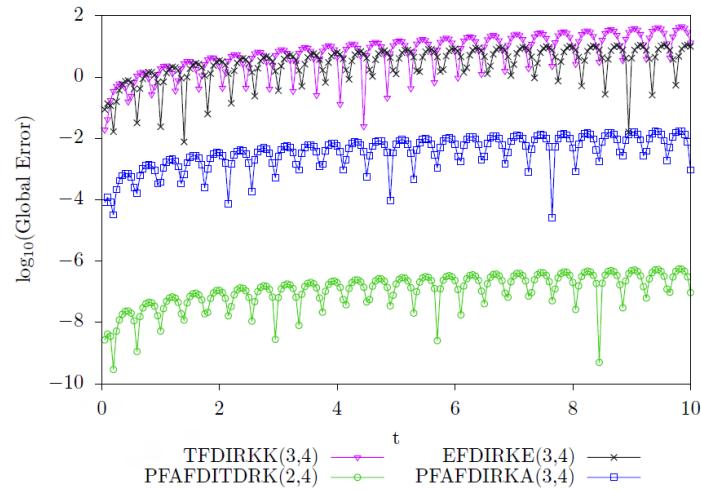


FIGURE 4. The error when solving the harmonic oscillator (Problem 1) at each integration point where $\omega = 8, q_{01} = 1, q_{02} = -2$ and $\Delta t = 1/20$

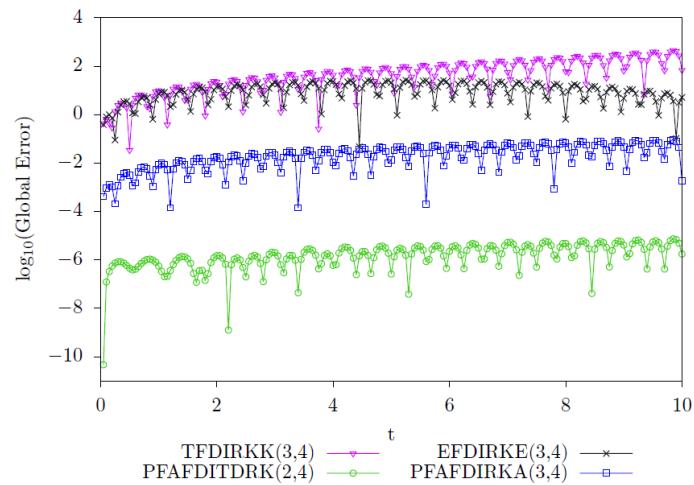


FIGURE 5. The global error when solving the inhomogeneous problem (Problem 2) at each integration point where $\Delta t = 1/20$

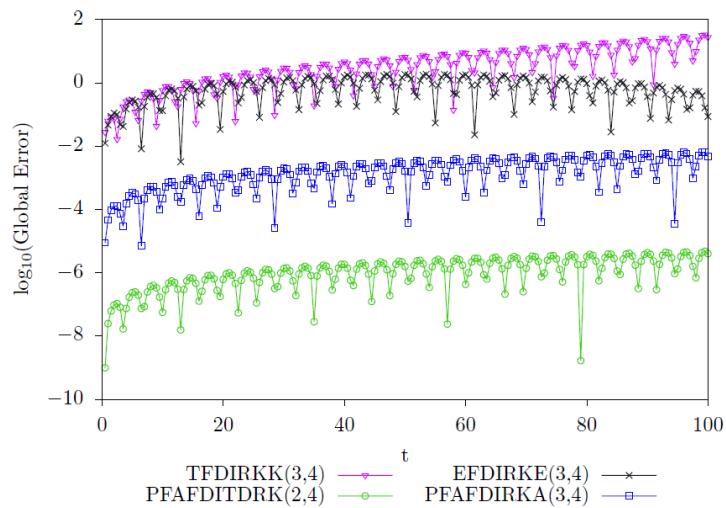


FIGURE 6. The global error when solving the almost periodic problem (Problem 3) at each integration point where $\Delta t=1/2$

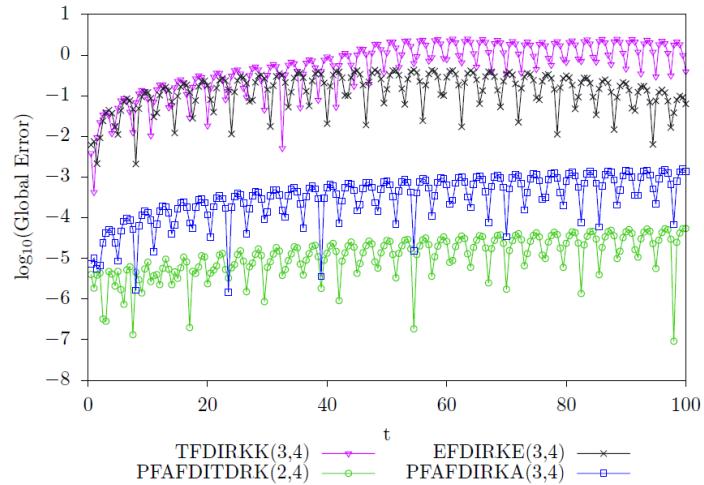


FIGURE 7. The global error when solving the Duffing problem
(Problem 4) at each integration point where $\Delta t = 1/2$

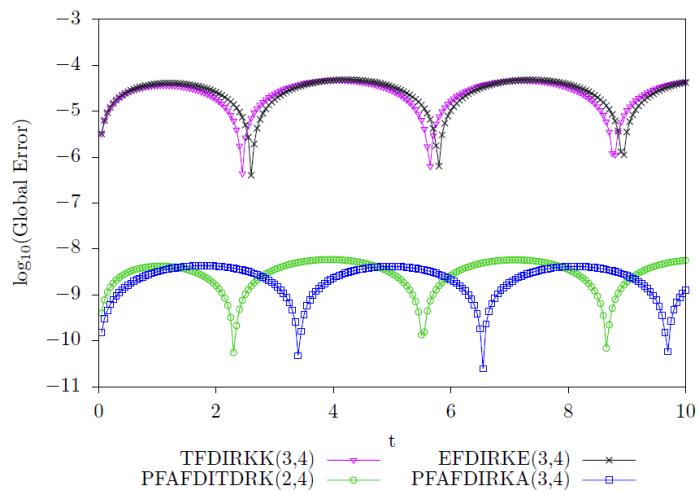


FIGURE 8. The global error when solving the Prothero-Robinson
problem (Problem 5) at each integration point where $\Delta t = 1/20$

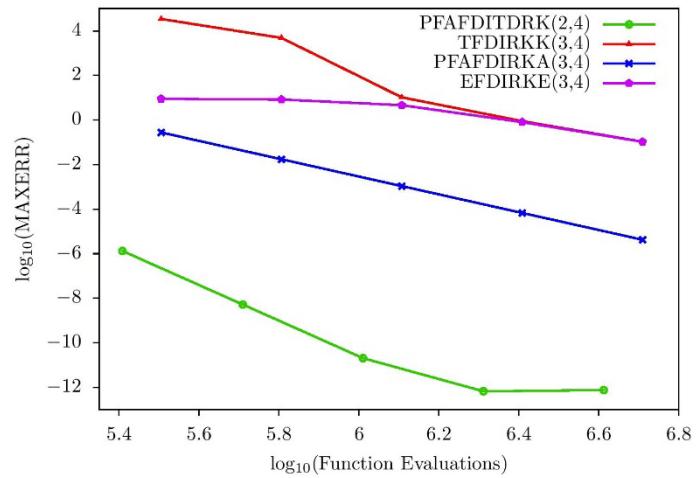


FIGURE 9. The curve for the harmonic oscillator (Problem 1) with
 $\lambda = 8, \Delta t = 1.0/2^i, i = 5, \dots, 9$ with $t_{\text{end}} = 1000$

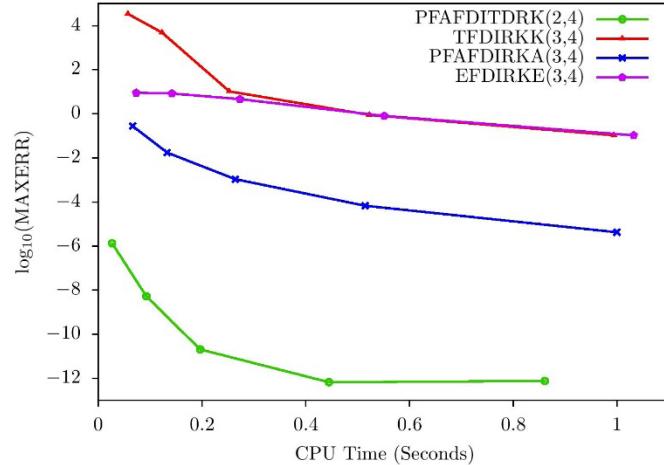


FIGURE 10. The curve for the harmonic oscillator (Problem 1) with $\lambda = 8, \Delta t = 1.0/2^i, i = 5, \dots, 9$ with $t_{\text{end}} = 1000$

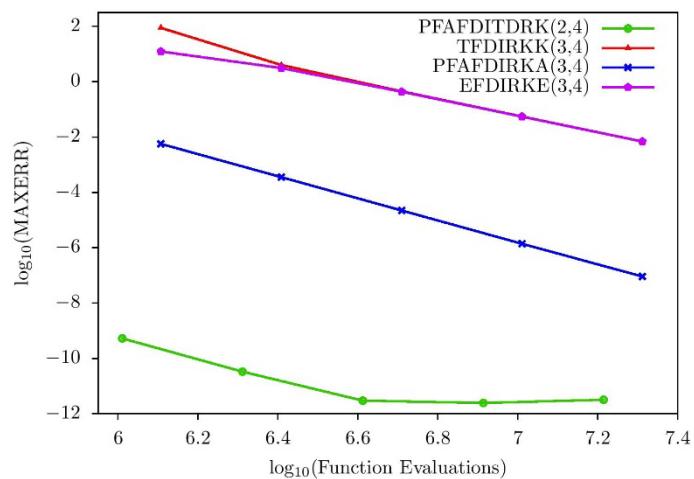


FIGURE 11. The curve for the inhomogeneous problem (Problem 2) with time step $\Delta t = 1.0/2^i, i = 7, \dots, 11$

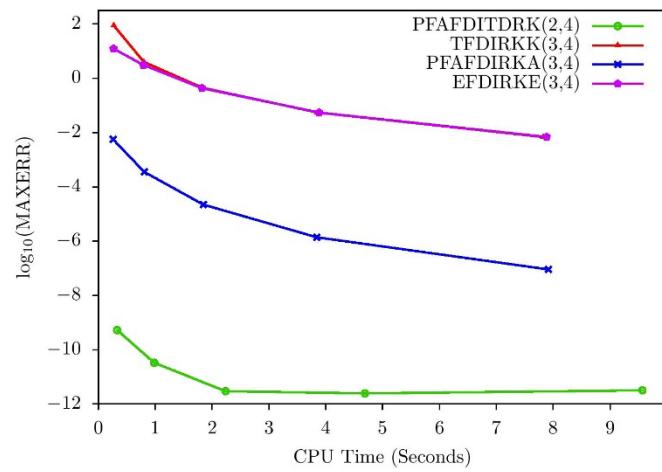


FIGURE 12. The curve for the inhomogeneous problem (Problem 2) with time step $\Delta t = 1.0/2^i, i = 7, \dots, 11$

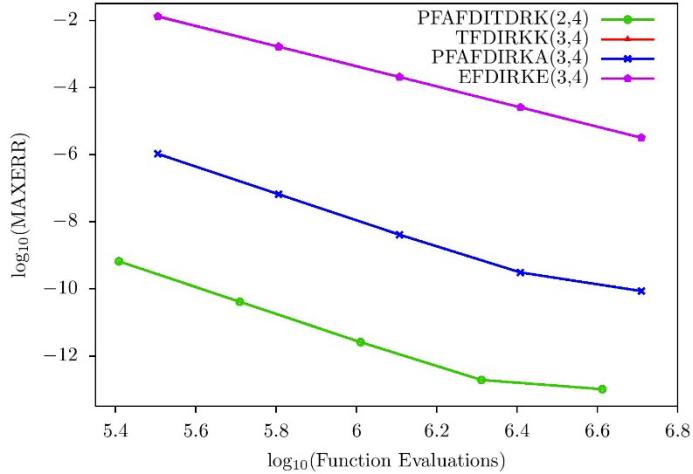


FIGURE 13. The curve for the almost periodic problem (Problem 3)
with time step $\Delta t = 1.0/2^i, i = 5, \dots, 9$

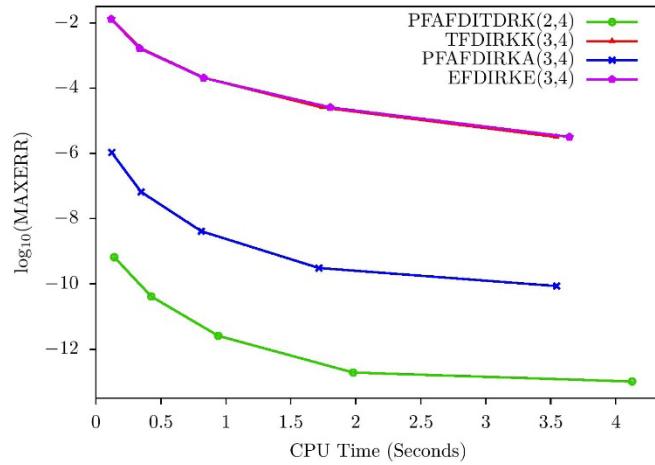


FIGURE 14. The curve for the almost periodic problem (Problem 3)
with time step $\Delta t = 1.0/2^i, i = 5, \dots, 9$

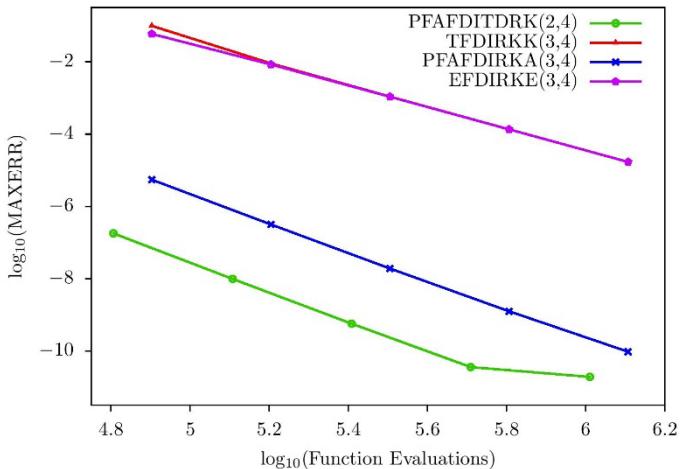


FIGURE 15. The curve for the Duffing problem (Problem 4) with
time step $\Delta t = 1.0/2^i, i = 3, \dots, 7$

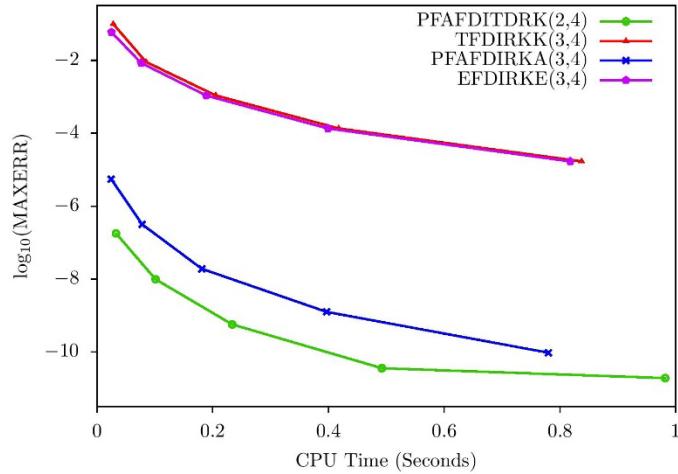


FIGURE 16. The curve for the Duffing problem (Problem 4) with time step $\Delta t = 1.0/2^i, i = 3, \dots, 7$

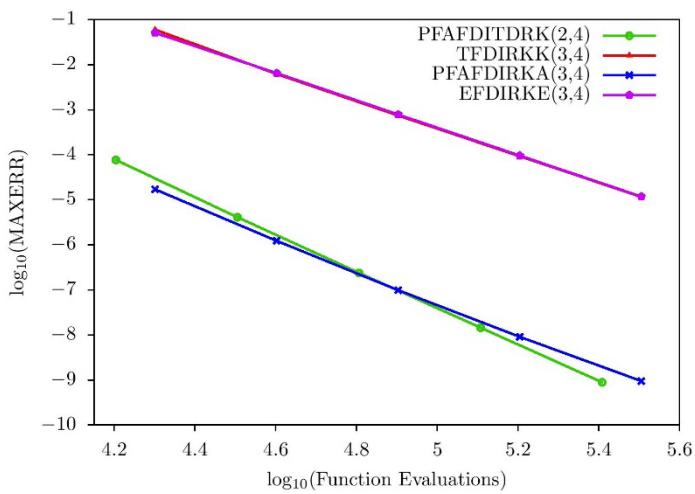


FIGURE 17. The curve for the Prothero-Robinson problem (Problem 5) with time step $\Delta t = 1.0/2^i, i = 1, \dots, 5$

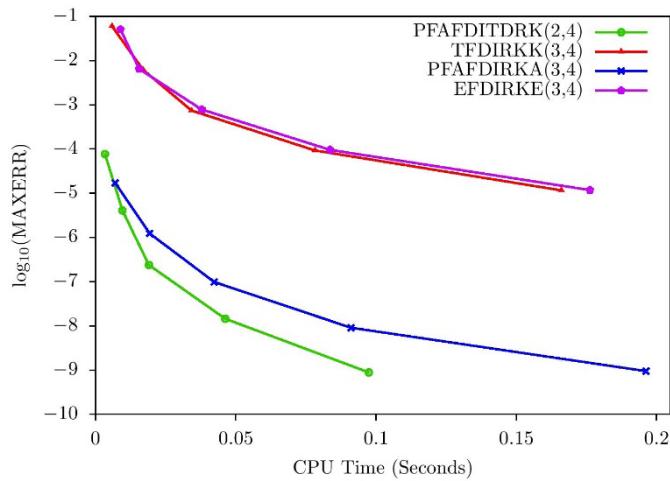


FIGURE 18. The curve for the Prothero-Robinson problem (Problem 5) with time step $\Delta t = 1.0/2^i, i = 1, \dots, 5$

DISCUSSION

The numerical results has shown the standard properties of the proposed phase-fitted and amplification-fitted DITDRK method, PFAFDITDRK(2,4) which was obtained earlier. Several well-known existing RK methods are chosen as the comparison with the proposed method. The energy error at every integration point can be seen in Figure 3. Conservation of energy is succeeded by the phase-fitted and amplification-fitted DITDRK method when it experienced smaller amount of energy error compared to TFDIRKK(3,4), PFAFDIRKA(3,4) and EFDIRKE(3,4). The log number of global error against the time of integration for different time step, are plotted for distinct problems as shown in Figures 4-8. From Figures 4-7, it is identified that global error developed by the PFAFDITDRK(2,4) method is smaller compared to TFDIRKK(3,4), PFAFDIRKA(3,4) and EFDIRKE(3,4). Meanwhile in Figure 8, the global error between PFAFDITDRK(2,4) and PFAFDIRKA(3,4) are rather close between one another but still the proposed method has the smallest global error.

Next, a long period of integration of the global error and the efficiency of the method are plotted. The log of the maximum global error versus the logarithm number of function evaluations and CPU time is plotted as given in Figures 9-18 to show the accuracy of the designed method. From Figures 9-18, the global error produced by PFAFDITDRK(2,4) method is smaller compared to the same order existing RK methods. In Figures 12, 14 and 16, PFAFDITDRK(2,4) takes longer CPU time compared to other existing RK methods due to its method complexity which is caused by the existence of the extra g to be evaluated at every step. In Figure 17, at the beginning of the graph, PFAFDITDRK(2,4) has slightly bigger maximum global error compared to PFAFDIRKA(3,4). As the value of Δt decreases, PFAFDITDRK(2,4) has smaller maximum global error compared to PFAFDIRKA(3,4). From Figures 9-18, it can be seen that PFAFDITDRK(2,4) method has the smallest maximum global error and the least amount of function evaluations per step.

One of the disadvantages of the derived method is that it is not suitable for solving stiff oscillatory or highly oscillatory problems which required the need of P-stable or strongly stable method. Therefore, we suggested that in the future work, the derivation of P-stable PFAFDITDRK is considered when one tries to solve stiff oscillatory or highly oscillatory problems.

Based on the phase-fitted and amplification-fitted property, the fitted property works well in solving linear problems but is not suitable in solving non-linear problems.

Hence, we did not include non-linear problem in the problems tested.

CONCLUSION

In this area of study, a fourth-order phase-fitted and amplification-fitted DITDRK method of has been proposed. Based on the numerical experiments, we can simplified that the proposed PFAFDITDRK(2,4) method is more promising than any of the other well-known existing DIRK methods with trigonometrically-fitted and phase-fitted and amplification-fitted property in terms of efficiency and accuracy as well as the number of function evaluations per step.

ACKNOWLEDGEMENTS

We are grateful to the Institute of Mathematical Research (INSPEM) and the Department of Mathematics and Statistics, Universiti Putra Malaysia for their continuous support and guidance throughout the research work. This study was supported financially by the Fundamental Research Grant Scheme (Ref. No. FRGS/1/2018/STG06/UPM/02/2) awarded by the Malaysia Ministry of Education and MyBrainSc.

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Received: 26 June 2020

Accepted: 27 October 2020