

# Spectral Deferred Correction Methods for Ordinary Differential Equations

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## Introduction to Scientific Computing

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### ✍ Problem Description

Consider an initial value problem:

$$\begin{cases} \varphi'(t) = F(t, \varphi(t)), & t \in [a, b] \\ \varphi(a) = \phi_a \end{cases} \quad -(1.1)$$

The goal is to find the function  $\varphi(t)$  satisfy these condition. IVP generally fall into two cases, Non-stiff problem and stiff problem. We will focus our attention on solving stiff problem.

For a stiff problem, standard numerical methods have the following issue:

- **Runge-Kutta:** These method are stable, but can become computationally expensive when requiring very high order
- **Implicit multistep algorithm:** the method can reach very high order convergence, but it have poor stability properties.

This paper aim to construct a method that can solve **stiff problem** with Arbitrary Order Accuracy. The strategy involves combining Picard Integral with a Gauss-Legendre grid.

## 1. Notation

### 1.1 The Picard Integral equation

Integrating equation (1.1) with respect to t, then we have

$$\varphi(t) = \varphi(a) + \int_a^t F(\tau, \varphi(\tau)) d\tau. \quad -(1.2)$$

which is call Picard Integral equation.

### 1.2 Residual function

Given an approximate solution  $\varphi^0(t)$  to (1.2), then we have

$$\epsilon(t) = \varphi(a) + \int_a^t F(s, \varphi^0(s)) ds - \varphi^0(t). \quad -(1.3)$$

### ✍ Remark

### Why we choose Picard Integral

Integration is numerically stable, whereas differentiation is numerically unstable.

Now we define the error  $\delta(t) = \varphi(t) - \varphi^0(t) - (\star)$ , then

$$\begin{aligned} \text{substitue } (\star) \text{ into (1.2)} \Rightarrow \quad \varphi(t) &= \varphi(a) + \int_a^t F(s, \delta(s) + \varphi^0(s)) ds - (1.5) \\ (1.5)-(1.3) \Rightarrow \quad \delta(t) &= \int_a^t F(s, \delta(s) + \varphi^0(s)) - F(s, \varphi^0(s)) ds + \epsilon(t) \\ \delta(t) &= \int_a^t G(s, \delta(s)) ds + \epsilon(t) \quad - (1.6) \end{aligned}$$

where  $G : \mathbb{R} \times \mathbb{C} \rightarrow \mathbb{C}$  given by  $G(t, \delta) = F(t, \varphi^0(t) + \delta(t)) - F(t, \varphi^0(t))$ . The form of (1.4) is called Picard-type integral.

### **Remark**

The structure of equation (1.6) is identical to the original Picard integral. Therefore, we can use the same numerical methods to solve for  $\delta(t)$ , eliminating the need to design a new method for our error terms.

### **Application to Euler's method**

We present an example for how to apply this notation to Euler's method. For the standard forward Euler method applied to  $\varphi$ , we have

$$\varphi_{i+1} = \varphi_i + h_i \cdot F(t_i, \varphi_i), \quad h_i = t_{i+1} - t_i.$$

Then, the forward Euler method for the solution (1.6) is given by

$$\delta_{i+1} = \delta_i + h_i \cdot G(t_i, \delta_i) + (\epsilon(t_{i+1}) - \epsilon(t_i))$$

### 1.3 Nodes selection

Given  $m \in \mathbb{N}$ , let  $r_1, r_2, \dots, r_m$  be the Gauss-Legendre nodes on the interval  $[-1, 1]$ . For an interval  $[a, b] \subset \mathbb{R}$ , let  $s_1, s_2, \dots, s_m$  be the  $m$  Gaussian nodes on the interval  $[a, b]$ , given by the formula.

$$s_i = \frac{b-a}{2} \cdot r_i + \frac{b+a}{2}$$

### **Remark**

### Why Gauss-Legendre Nodes

- **Spectral Accuracy:** Gaussian quadrature provides the highest possible degree of precision for numerical integration, which is essential for the high order convergence rates targeted by the SDC method.
- **Avoiding the Runge Phenomenon:**

### 1.4 Lagrange Interpolation

Given the nodes  $t_1, t_2, \dots, t_m$  and function values  $\varphi = (\varphi_1, \varphi_2, \dots, \varphi_m)$ , the Lagrange interpolant  $L^m(\varphi, t)$  is the unique polynomial of degree  $m - 1$  defined by

$$L^m(\varphi, t) = \sum_{i=1}^m c_i(t) \cdot \varphi_i \quad (1.7), \quad \text{where } c_i(t) = \prod_{j \neq i} \frac{t - t_j}{t_i - t_j}$$

### 1.5 Spectral Integration and Differential Matrices

Let  $f = (f_1, f_2, \dots, f_m)$  be the vector of function values at the nodes defined by

$$f = \begin{bmatrix} F(t_1) \\ F(t_2) \\ \vdots \\ F(t_m) \end{bmatrix}.$$

Then, the integration matrix  $S^m$  is the linear operator that maps the function values  $f$  to the integrals of the interpolant from start of the interval to each grid node

$$(S^m(f))_i = \int_a^{t_i} L^m(f, t) dt.$$

Also, there can be written in:

$$S^m = \begin{pmatrix} \int_a^{t_1} L_1(f, t) dt & \cdots & \int_a^{t_m} L_1(f, t) dt \\ \vdots & & \vdots \\ \int_a^{t_1} L_m(f, t) dt & \cdots & \int_a^{t_m} L_m(f, t) dt \end{pmatrix}$$

#### Remark

- If the underlying function  $F$  is a polynomial of degree  $m - 1$ , the operations  $S^m$  are exact. For smooth functions, the error decays exponentially with  $m$ .
- The matrix  $S^m$  depends only on the nodes and is **fixed** for all iterations.

## 2. Spectral deferred correction(SDC).

Bases on the derived error equation and spectral matrices, we will construct the full iterative algorithm. The method follows a prediction-correction paradigm

## 2.1 Step of algorithm

1. Assume we have an approximate solution  $\varphi^{[0]}$  on the grip  $[a, b]$ , By using the standard Gauss-Legendre nodes on  $[a, b]$ ,  $\varphi^{[j]}$  will be used to denote the  $j$ -th approximate solution.

$$\varphi^{[j]} = (\varphi_1^{[j]}, \varphi_2^{[j]}, \dots, \varphi_m^{[j]}) = (\varphi(s_1), \varphi(s_2), \dots, \varphi(s_m))$$

2. The residual function  $\epsilon(t)$  will be approximate by the vector  $\varphi^{[j]}$  defined by

$$\sigma(\varphi^{[j]}) = S^m \bar{F}(\varphi^{[j]}) - \varphi^{[j]} + \bar{\varphi}_a$$

pseudocode

```
1:Input:Initial value, number of nodes m, number of sweeps J
2:Prediction(Initialization):
3:Compute an initial approximate solution on the Gauss-Legendre grid using a
low-order method
4:\Correlation loop
5:for j=1 to J do
    ##Compute Residual
    Calculate the residual vector using the spectral integration matrix S^m
    \sigma(\varphi^{[j]})=S^m\bar{F}(\varphi^{[j]})-
    \varphi^{[j]}+\bar{\varphi}_a

    ##Solve error Equation
    Solve for the correction using a first-order method on the error
equation
    \delta^{[j]}=C_{imp}(G,\sigma(\varphi^{[j-1]}))

    ##Update Solute
    \varphi^{[j]}=\varphi^{[j-1]}+\delta
6:Output:High-order solution \varphi^{[J]}
```

## 3.Analysis of method

### 3.1 Behavior at large $|\lambda|$ and extrapolation

- **Explicit Euler:** For stiff problem, the explicit SDC schemes are unstable
- **Implicit Euler:** The implicit SDC schemes are generally stable. However, when analyzing the amplification factor,  $A_m(\lambda)$  as  $|\lambda| \rightarrow \infty$ , we observe

$$\lim_{|\lambda| \rightarrow \infty} = A_m(\lambda) = \mu(m, J) \neq 0$$

This means that the method is Not L-stable.

Therefore, we construct a combinations method.

### Composite Schemes

To achieve L-stability, we compute two solutions by using two different number of nodes  $m_1, m_2$  with asymptotic limits  $\mu_1, \mu_2$ . The linear combination

$$\varphi_{comb} = \frac{\mu_2 \varphi_1 - \mu_1 \varphi_2}{\mu_2 - \mu_1}.$$

### 3.2 Order of Accuracy

For any sufficiently smooth function F, the approximation convergence with an order of accuracy given by

$$Order = \min(m, J + 1)$$

where m is the number of Gauss-Legendre nodes and J is the number of correction sweeps.

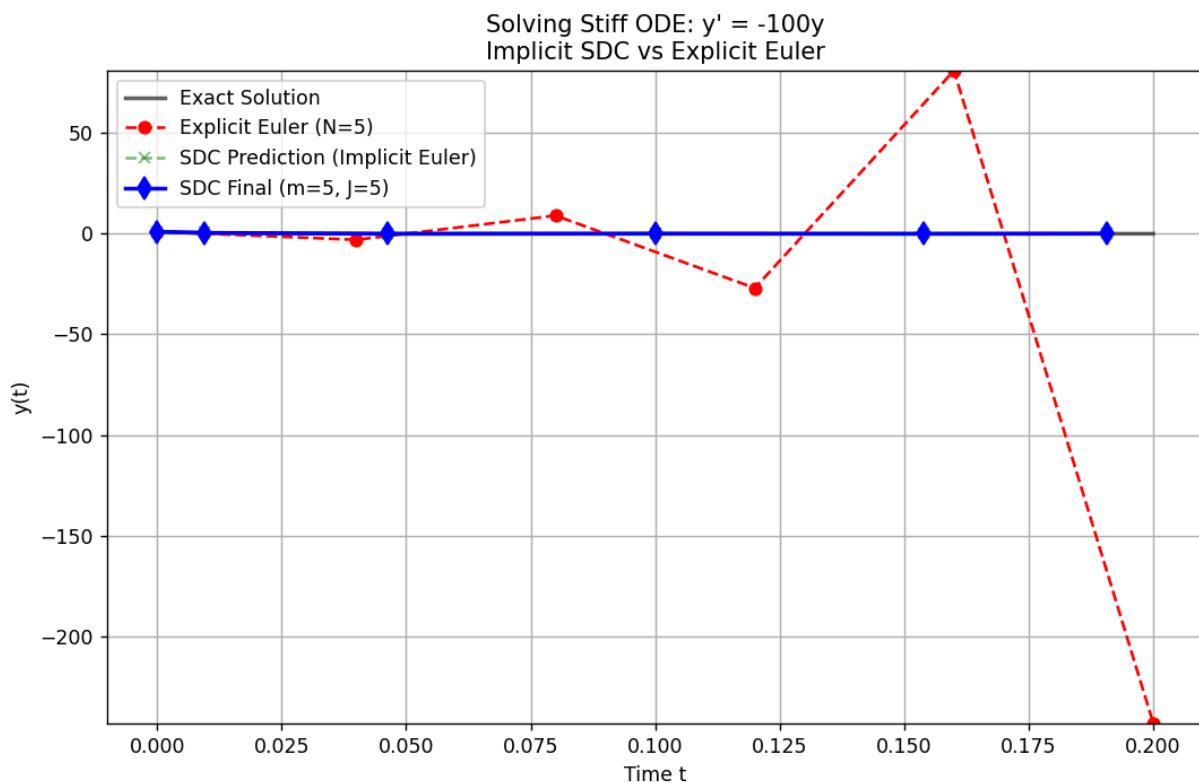
## 4. Numerical example

Consider

$$\begin{cases} \varphi' = -100\varphi \\ \varphi(0) = 1 \end{cases}$$

### 4.1 Analysis of result

- **Explicit Euler method:** The Explicit Euler solution (red dashed line) violates the stability condition ( $0.04 > 0.02$ ). Consequently, it is unstable, oscillating and diverging to values near -250 at  $t = 0.2$ .
- **SDC:** The SDC result remains bounded and stable despite the large step size. Furthermore, the iterative spectral deferred correction sweeps successfully boost the accuracy, resulting in a solution that is visually indistinguishable from the exact solution



## 5. Conclusion

We explored Spectral Deferred Correction (SDC) methods for solving stiff ordinary differential equations with arbitrary order accuracy. By using Picard integral equation and Gauss-Legendre quadrature, the SDC approach iteratively corrects a low-order approximation to achieve high-precision results. Consequently, SDC proves to be a powerful and flexible alternative to traditional methods for high-accuracy scientific computing.