

AN EXTENDED TECHNIQUE FOR COMPUTATION OF LAPLACE TRANSFORMED DYNAMIC FUNDAMENTAL SOLUTIONS FOR 3D ANISOTROPIC ELASTIC SOLIDS

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Abstract. This paper presents an extended procedure for computation of integral representations of regular parts of Laplace domain three-dimensional dynamic anisotropic elastic full space displacement fundamental solutions and their spatial derivatives. The problem is that under specific conditions these integrals become highly oscillatory. For the modified integral expressions, we present a technique that utilizes specialized quadrature rule which in turn is a variation of well-known Levin's method for solving highly oscillatory integrals. Results of numerical investigations suggest improved performance (regarding number of integration points) compared to using the Gauss-Legendre quadrature.

1 INTRODUCTION

Elastodynamic fundamental solutions (or Green's functions) play significant role in formulation and solution of initial boundary value problems. Besides that, fundamental solutions are essential for the development of diverse formulations of Boundary Element Method.

For general homogeneous anisotropic linearly elastic media full-space fundamental solutions are not available in the explicit closed-form expressions. Frequency domain anisotropic fundamental solutions are represented as a sum of static (fundamental solutions of corresponding static problem) and dynamic parts.

Pan and Chen [1] presented an excellent work dedicated to evaluation of static anisotropic fundamental solutions. Concerning the dynamic parts of elastodynamic fundamental solutions for anisotropic solids practically important results were provided by Wang and Achenbach [2, 3]. They suggested using Radon transform and obtained dynamic parts of fundamental solutions in the form of surface integrals over a half of a unit sphere. Resulting

expressions have rather complicated integrands that become rapidly oscillating for large distances between source and observation points or high frequencies. For these specific cases we propose an extended yet simple computational strategy based on a quadrature rule developed by Evans and Webster [4]. This quadrature is a variation of the well-known Levin's collocation method [5]. Results of numerical experiments indicate significantly improved performance of suggested procedure when compared to using the Gauss-Legendre quadrature.

2 LAPLACE DOMAIN ELASTODYNAMIC FUNDAMENTAL SOLUTIONS

Laplace domain full space anisotropic elastic dynamic fundamental solutions for displacements and tractions are represented as sums of respective singular (static) and regular (dynamic) parts as

$$\bar{g}_{ij}(\mathbf{y}, \mathbf{x}, s) = \bar{g}_{ij}(\mathbf{r}, s) = g_{ij}^S(\mathbf{r}) + \bar{g}_{ij}^R(\mathbf{r}, s), \quad i, j = \overline{1, 3}, \quad (1)$$

$$\bar{h}_{mi}(\mathbf{y}, \mathbf{x}, s) = \bar{h}_{mi}(\mathbf{r}, s) = C_{ijkl} g_{mk,l}^S(\mathbf{r}) n_j(\mathbf{y}) + C_{ijkl} \bar{g}_{mk,l}^R(\mathbf{r}, s) n_j(\mathbf{y}) = h_{mi}^S(\mathbf{r}) + \bar{h}_{mi}^R(\mathbf{r}, s), \quad k, l, m = \overline{1, 3}, \quad (2)$$

$$\mathbf{r} = \mathbf{y} - \mathbf{x}, \quad r = |\mathbf{r}|, \quad (3)$$

where g_{ij}^S and h_{mi}^S are the static parts, \bar{g}_{ij}^R is the dynamic part, s is the complex frequency, \mathbf{y} and \mathbf{x} are the position vectors of the observation point source point, respectively, n_j is the unit vector at \mathbf{y} , C_{ijkl} is the fourth order elasticity tensor.

In this paper, we concern ourselves only with the aspects of computing the dynamic part of displacement fundamental solution \bar{g}_{ij}^R and its spatial derivative $\bar{g}_{ij,l}^R$. Their modified and expanded expressions of those obtained in [2, 3] can be written as follows

$$\bar{g}_{ij}^R(\mathbf{r}, s) = -\frac{s\sqrt{\rho}}{8\pi^2} \int_0^{2\pi} \int_0^{\pi/2} \sum_{m=1}^3 \frac{E_{im}(\varphi, \psi) E_{jm}(\varphi, \psi) \sin(\psi)}{\lambda_m^{3/2}(\varphi, \psi)} e^{sr\sqrt{\rho} \left(-\frac{\cos(\psi)}{\sqrt{\lambda_m(\varphi, \psi)}} \right)} d\psi d\varphi, \quad (4)$$

$$\bar{g}_{ij,l}^R(\mathbf{r}, s) = \frac{s^2 \rho}{8\pi^2} \int_0^{2\pi} \int_0^{\pi/2} \sum_{m=1}^3 \frac{n_l(\varphi, \psi) E_{im}(\varphi, \psi) E_{jm}(\varphi, \psi) \sin(\psi)}{\lambda_m^2(\varphi, \psi)} e^{sr\sqrt{\rho} \left(-\frac{\cos(\psi)}{\sqrt{\lambda_m(\varphi, \psi)}} \right)} d\psi d\varphi, \quad (5)$$

$$c_m(\varphi, \psi) = \sqrt{\frac{\lambda_m(\varphi, \psi)}{\rho}}, \quad k_m(\varphi, \psi) = \frac{s}{c_m(\varphi, \psi)}, \quad (6)$$

where $\lambda_m(\varphi, \psi)$ and $E_{jm}(\varphi, \psi)$ are the eigenvalues and the corresponding eigenvectors of the matrix $\Gamma_{ij}(\mathbf{n}(\varphi, \psi)) = C_{kijl} n_k(\varphi, \psi) n_l(\varphi, \psi)$. Vector $\mathbf{n}(\varphi, \psi)$ is defined as (see Figures 1, 2)

$$\mathbf{n}(\varphi, \psi) = \mathbf{d}(\varphi) \sin \psi + \mathbf{e} \cos \psi = [n_1, n_2, n_3]^T, \quad (7)$$

$$\mathbf{d}(\varphi) \cdot \mathbf{r} = 0, \quad \mathbf{e} = \frac{\mathbf{r}}{r}, \quad \mathbf{e} = [e_1, e_2, e_3]^T, \quad (8)$$

$$\mathbf{p} = \frac{[e_2, -e_1, 0]^T}{\sqrt{1-e_3^2}}, \quad \mathbf{q} = \mathbf{e} \times \mathbf{p} = \frac{[e_1 e_3, e_2 e_3, -(1-e_3^2)]^T}{\sqrt{1-e_3^2}}, \quad (9)$$

$$\mathbf{p} \cdot \mathbf{e} = 0, \quad \mathbf{q} \cdot \mathbf{e} = 0, \quad \mathbf{p} \cdot \mathbf{q} = 0, \quad \mathbf{d}(\varphi) = \mathbf{p} \cos \varphi + \mathbf{q} \sin \varphi, \quad 0 \leq \varphi \leq 2\pi, \quad (10)$$

$$\mathbf{d}(\varphi) = \frac{[e_2 \cos \varphi + e_1 e_3 \sin \varphi, -e_1 \cos \varphi + e_2 e_3 \sin \varphi, -(1-e_3^2) \sin \varphi]^T}{\sqrt{1-e_3^2}}. \quad (11)$$

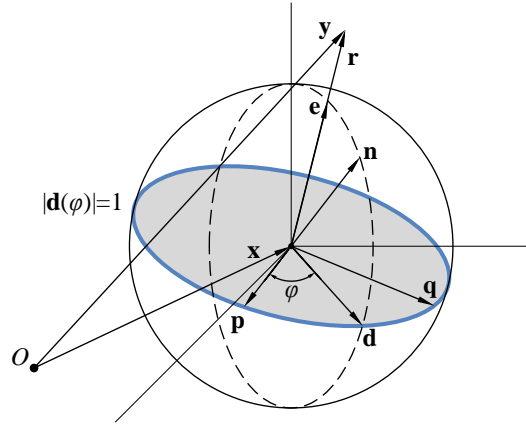


Figure 1: Overview of quantities required for integration of dynamic part of fundamental solution and its derivative

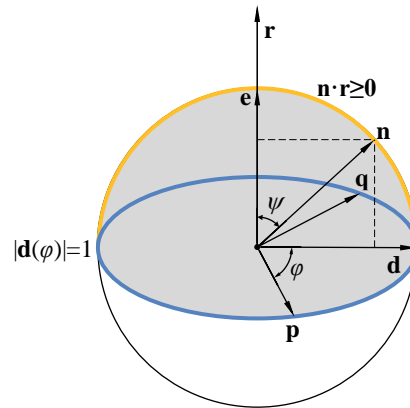


Figure 2: Quantities involved in defining vector \mathbf{n}

Close inspection of integrals defined in equations (4) and (5) reveals that for numerical integration of these integrals an eigenvalue problem for Christoffel matrix $\Gamma_{ij}(\mathbf{n}(\varphi, \psi))$ has to be also solved numerically for every value of φ and ψ which is very computationally expensive. For practically feasible implementation of Laplace domain anisotropic elastodynamic fundamental solutions in any numerical method it is important to minimize

number of integration points required for numerical evaluation of integrals in (4) and (5).

Let value of φ be fixed and consider integrals with the respect to ψ

$$I_1^{ij}[\varphi, r, s] = I_1[\varphi, r, s] = \int_0^{\pi/2} \sum_{m=1}^3 \frac{\sin(\psi) E_{im}(\psi) E_{jm}(\psi)}{\lambda_m^{3/2}(\psi)} e^{sr\sqrt{\rho} \left(-\frac{\cos(\psi)}{\sqrt{\lambda_m(\psi)}} \right)} d\psi, \quad (12)$$

$$I_2^{ij,k}[\varphi, r, s] = I_2[\varphi, r, s] = \int_0^{\pi/2} \sum_{m=1}^3 \frac{n_k(\psi) \sin(\psi) E_{im}(\psi) E_{jm}(\psi)}{\lambda_m^2(\psi)} e^{sr\sqrt{\rho} \left(-\frac{\cos(\psi)}{\sqrt{\lambda_m(\psi)}} \right)} d\psi. \quad (13)$$

We denote complex frequency as

$$\tau = sr\sqrt{\rho} = \langle s = \alpha + i\omega \rangle = \alpha r\sqrt{\rho} + i\omega r\sqrt{\rho}. \quad (14)$$

With (14) we have

$$I_1^{ij}[\varphi, \tau] = I_1[\varphi, \tau] = \int_0^{\pi/2} \sum_{m=1}^3 \frac{\sin(\psi) E_{im}(\psi) E_{jm}(\psi)}{\lambda_m^{3/2}(\psi)} e^{\tau \left(-\frac{\cos(\psi)}{\sqrt{\lambda_m(\psi)}} \right)} d\psi, \quad (15)$$

$$I_2^{ij,k}[\varphi, \tau] = I_2[\varphi, \tau] = \int_0^{\pi/2} \sum_{m=1}^3 \frac{n_k(\psi) \sin(\psi) E_{im}(\psi) E_{jm}(\psi)}{\lambda_m^2(\psi)} e^{\tau \left(-\frac{\cos(\psi)}{\sqrt{\lambda_m(\psi)}} \right)} d\psi. \quad (16)$$

Exponential term in the integrands makes these integrals highly oscillatory for the large values of the imaginary part $\text{Im}(\tau) = \omega r\sqrt{\rho}$ of the complex frequency τ .

For further convenience we rewrite integrals $I_1^{ij}[\varphi, \tau]$ and $I_2^{ij,k}[\varphi, \tau]$ as follows

$$I_1^{ij}[\varphi, \tau] = I_1[\varphi, \tau] = \int_0^{\pi/2} \sum_{m=1}^3 \frac{\sin(\psi) E_{im}(\psi) E_{jm}(\psi)}{\lambda_m^{3/2}(\psi)} e^{\tau \left(-\frac{\cos(\psi)}{\sqrt{\lambda_m(\psi)}} \right)} d\psi = \sum_{m=1}^3 \int_0^{\pi/2} f_{ij}^m(\psi) e^{\tau q_m(\psi)} d\psi, \quad (17)$$

$$\begin{aligned} I_2^{ij,k}[\varphi, \tau] &= I_2[\varphi, \tau] = \int_0^{\pi/2} \sum_{m=1}^3 \frac{n_k(\psi) \sin(\psi) E_{im}(\psi) E_{jm}(\psi)}{\lambda_m^2(\psi)} e^{\tau \left(-\frac{\cos(\psi)}{\sqrt{\lambda_m(\psi)}} \right)} d\psi \\ &= \sum_{m=1}^3 \int_0^{\pi/2} g_{ij,k}^m(\psi) e^{\tau q_m(\psi)} d\psi, \end{aligned} \quad (18)$$

$$f_{ij}^m(\psi) = \frac{\sin(\psi) E_{im}(\psi) E_{jm}(\psi)}{\lambda_m^{3/2}(\psi)}, \quad g_{ij,k}^m(\psi) = \frac{n_k(\psi) \sin(\psi) E_{im}(\psi) E_{jm}(\psi)}{\lambda_m^2(\psi)}, \quad (19)$$

$$q_m(\psi) = -\frac{\cos(\psi)}{\sqrt{\lambda_m(\psi)}}, \quad (20)$$

where $f_{ij}^m(\psi)$ and $g_{ij,k}^m(\psi)$ are amplitude functions, and $q_m(\psi)$ are corresponding to phase functions.

3 COMPUTATION OF INTEGRALS

Specialized procedures [6] are required to efficiently evaluate highly oscillatory integrals defined in equations (17) and (18).

Amplitude functions $f_{ij}^m(\psi)$ and $g_{ij,k}^m(\psi)$ and phase functions $q_m(\psi)$ are not available in an explicit closed-form and for every value of integration variable ψ they need to be evaluated numerically. This prevents from using any Filon-type method to compute integrals (17) and (18).

For a complicated phase functions such as in our case, Levin's collocation method offers a suitable framework for evaluating highly oscillatory integrals. In Levin's approach an integration problem of form

$$I = \int_{-1}^1 f(x) e^{i\gamma q(x)} dx, \quad (21)$$

is transformed into an ordinary differential equation (ODE)

$$F'(x) + i\gamma q'(x)F(x) = f(x). \quad (22)$$

After ODE (22) is solved, integral (21) is obtained by substitution

$$I = \int_{-1}^1 (F'(x) + i\gamma q'(x)F(x)) e^{i\gamma q(x)} dx = \int_{-1}^1 \frac{d}{dx} [F(x) e^{i\gamma q(x)}] dx = F(1) e^{i\gamma q(1)} - F(-1) e^{i\gamma q(-1)}. \quad (23)$$

To find $F(x)$ it is approximated as

$$F(x) \approx F_n(x) = \sum_{k=1}^n \beta_k v_k(x), \quad (24)$$

where $v_k(x)$ are linearly independent basis functions.

Therefore integral (21) can be approximated substituting (24) into (23)

$$I \approx I_n = \sum_{k=1}^n \beta_k v_k(1) e^{i\gamma q(1)} - \sum_{k=1}^n \beta_k v_k(-1) e^{i\gamma q(-1)}. \quad (25)$$

In order to obtain coefficients β_k , a following collocation system is assembled on a set of points $\{x_j\}$, $j = \overline{1, n}$, and then solved

$$\sum_{k=1}^n \beta_k v_k'(x_j) + i\gamma q'(x_j) \sum_{k=1}^n \beta_k v_k(x_j) = f(x_j). \quad (26)$$

To evaluate integrals (17) and (18) using Levin's method it would require to solve three complex valued collocation systems with 24 different right-hand sides each. We employ a variation of Levin's method proposed by Evans and Webster [4]. Consider a quadrature rule

$$I = \int_{-1}^1 f(x) e^{i\gamma q(x)} dx \approx \sum_{j=0}^N w_j f(x_j), \quad (27)$$

which is exact for the functions

$$h_k(x) = i\tau q'(x)p_k(x) + p'_k(x), \quad k = \overline{0, N}. \quad (28)$$

Functions defined in (28) are substituted on a set of points $\{x_j\}$ into (27) to obtain linear system for the weights w_j

$$a_{kj}w_j = b_k, \quad (29)$$

$$a_{kj} = i\tau q'(x_j)p_k(x_j) + p'_k(x_j), \quad k, j = \overline{0, N}, \quad (30)$$

$$b_k = p_k(1)e^{i\tau q(1)} - p_k(-1)e^{i\tau q(-1)}, \quad k = \overline{0, N}. \quad (31)$$

Following original paper by Evans and Webster [4]

$$x_j = \cos(j\pi/N), \quad p_j(x) = T_j(x), \quad j = \overline{0, N}, \quad (32)$$

where $T_j(x)$ is the j -th Chebyshev polynomial of the first kind.

In our case, to assemble systems defined by (29) we need to define derivatives of the phase functions

$$q'_m(\psi) = \frac{1}{2} \frac{2\lambda'_m(\psi)\sin(\psi) + \lambda''_m(\psi)\cos(\psi)}{\lambda_m^{3/2}(\psi)}, \quad \lambda'_m(\psi) = \sum_{i=1}^3 \sum_{j=1}^3 E_{im}(\mathbf{n})E_{jm}(\mathbf{n})\Lambda_{ij}(\mathbf{n}, \boldsymbol{\beta}), \quad (33)$$

$$\boldsymbol{\beta}(\varphi, \psi) = \mathbf{d}(\varphi)\cos(\psi) - \mathbf{e}\sin(\psi) = [\beta_1, \beta_2, \beta_3]^T, \quad \Lambda_{ijk}(\mathbf{n}, \boldsymbol{\beta}) = C_{ijkl}(\beta_i n_l + n_i \beta_l), \quad i, j, k, l = \overline{1, 3}. \quad (34)$$

To evaluate integrals (17) and (18) domain of integration $[0, \pi/2]$ over ψ is subdivided into a sufficient number of intervals and Evans-Webster quadrature is applied on each interval.

4 NUMERICAL EXAMPLE

Consider evaluation of integrals $I_1^{ij}[\varphi, \tau]$ and $I_2^{ij,k}[\varphi, \tau]$ for the following parameters

$$\rho = 2216 \text{ kg/m}^3, \quad \mathbf{C} = \begin{bmatrix} 17.77 & 3.78 & 3.76 & 0.24 & -0.28 & 0.03 \\ & 19.45 & 4.13 & -0.41 & 0.07 & 1.13 \\ & & 21.79 & -0.12 & 0.01 & 0.38 \\ & & & 8.30 & 0.66 & 0.06 \\ \text{symm.} & & & & 7.62 & 0.52 \\ & & & & & 7.77 \end{bmatrix} \text{GPa}, \quad \varphi = \pi/5. \quad (35)$$

Dimensionless quantities are introduced

$$\tilde{\mathbf{C}} = \mathbf{C}/p_{\max}, \quad \tilde{\rho} = \rho l_{\max}^2/p_{\max}/t_{\max}^2, \quad \tilde{s} = 2.0 + i\omega, \quad \tau = \tilde{s}\tilde{r}\sqrt{\tilde{\rho}}, \quad (36)$$

$$t_{\max} = 0.005 \text{ s}, \quad p_{\max} = 1.0 \cdot 10^{10} \text{ Pa}, \quad l_{\max} = 200 \text{ m}. \quad (37)$$

$$\tilde{\mathbf{x}} = [0, 0, 0]^T, \quad \tilde{\mathbf{y}} = [\sin(\pi/3)\cos(\pi/4), \sin(\pi/3)\sin(\pi/4), \cos(\pi/3)]^T, \quad \tilde{r} = |\tilde{\mathbf{y}} - \tilde{\mathbf{x}}| = 1. \quad (38)$$

The errors between results obtained with Evans-Webster (EW) based procedure and converged values obtained with high-order Gauss-Legendre (GL) rule denoted with

superscript “*ref*” are measured as

$$err(I_1) = \frac{\|I_1^{ref} - I_1\|_2}{\|I_1^{ref}\|_2}, \quad err(I_2) = \max_{1 \leq k \leq 3} \left(\frac{\|I_2^{ij,k;ref} - I_2^{ij,k}\|_2}{\|I_2^{ij,k;ref}\|_2} \right), \quad (39)$$

where $\|\cdot\|_2$ denotes the spectral norm of a matrix.

Different orders of EW method on each interval are considered for $\omega=500$. Figures 3 and 4 present obtained errors depending on the total number of evaluations of integrands. Results for EW orders 5, 6, and 7 are depicted in Figure 3 and for EW orders 8, 9, and 10 are depicted in Figure 4.

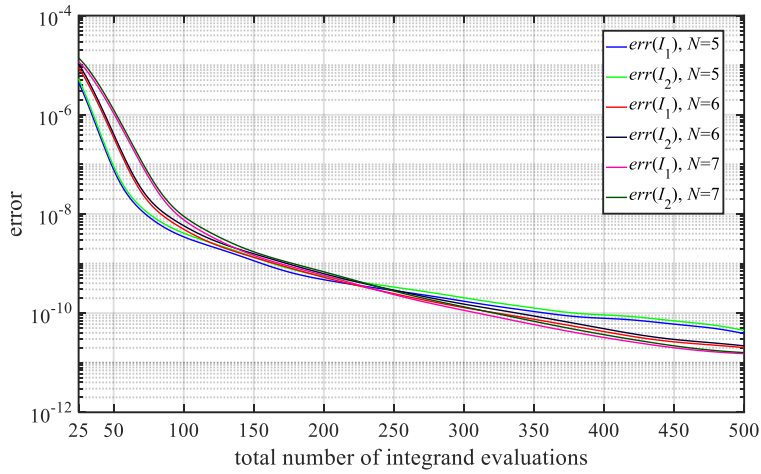


Figure 3: Errors $err(I_1)$ and $err(I_2)$ for $\omega = 500$, $N = 5, 6, 7$

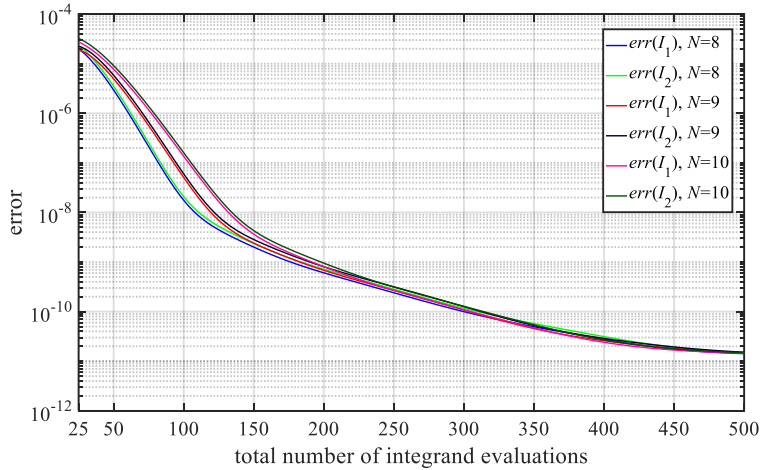


Figure 4: Errors $err(I_1)$ and $err(I_2)$ for $\omega = 500$, $N = 8, 9, 10$

Now we fix maximum allowed error for evaluating I_2 . For frequencies in interval $1 \leq \omega \leq 1000$, Figure 5 presents the number of integrand evaluation required for

$err(I_2) \leq 1 \cdot 10^{-10}$ for Gauss-Legendre quadrature and proposed technique with EW method orders $N = 5, 6, 7, 8, 9$ and 10 .

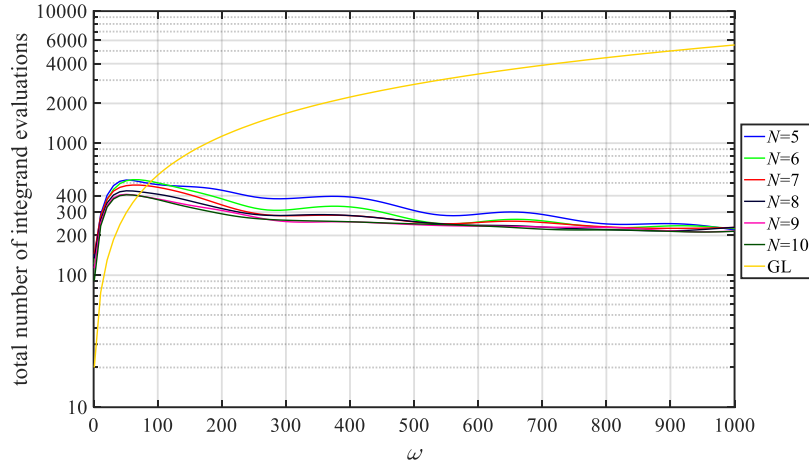


Figure 5: Number of integrand evaluations required for $err(I_2) \leq 1 \cdot 10^{-10}$ and $1 \leq \omega \leq 1000$

For lower frequencies $\omega < 100$ GL predictably performs better. Obtained results indicate that required numbers of integrand evaluations for all considered EW orders are close. Therefore, it is computationally beneficial to use lower orders in practice.

5 CONCLUSIONS

- A simple technique based on a variation of well-known Levin's method is presented for evaluation of highly oscillatory integrals appearing in Laplace domain anisotropic elastodynamic solutions.
- For high frequencies presented approach performs favorably in terms of required number of integrand evaluations compared to using Gauss-Legendre quadrature rule.
- Obtained results of numerical example indicate that errors decrease with increasing frequency which suggest that proposed procedure enjoys asymptotic property of Levin's method.

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