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A hybrid analytical-numerical method for solving advection-dispersion problems on a half-line



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ABSTRACT

This paper employs the unified transform, also known as the Fokas method, to solve the advectiondispersion equation on the half-line. This method combines complex analysis with numerics. Compared to classical approaches used to solve linear partial differential equations (PDEs), the unified transform avoids the solution of ordinary differential equations and, more importantly, constructs an integral representation of the solution in the complex plane which is uniformly convergent at the boundaries. As a consequence, such solutions are well suited for numerical computations. Indeed, the numerical evaluation of the solution requires only the computation of a single contour integral involving an integrand which decays exponentially fast for large values of the integration variable. A novel contribution of this paper, with respect to the solution of linear evolution PDEs in general, and the implementation of the unified transform in particular, is the following: using the advection-dispersion equation as a generic example, it is shown that if the transforms of the given data can be computed analytically, then the unified transform yields a fast and accurate method that converges exponentially with the number of evaluations *N* yet only has complexity O(N). Furthermore, if the transforms are computed numerically using *M* evaluations, the unified transform gives rise to a method with complexity O(NM). Results are successfully compared to other existing solutions.

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1. Introduction

Several problems in physics and engineering are modelled by means of the advection-dispersion equation (ADE). Starting from the pioneering work of Taylor [1], the ADE has been used to predict transport of passive or reactive scalars in natural (e.g. atmospheric, subsurface and surface flows) as well as in confined (e.g., channels and tubes) environments. Given its importance in many fields in science and engineering, it is not surprising that there exists a plethora of analytical solutions for the ADE under different physical configurations. In this work, we focus on the use of the ADE in environmental flows such as those occurring in natural porous formations (e.g., groundwater) and surface water bodies (e.g., rivers and channels).

There have been significant efforts in developing analytical solutions for the ADE in the environmental engineering and hydrological communities. For example, Ogata and Banks [2], van

Genuchten and Alves [3], Kreft and Zuber [4], Lee [5] and Guerrero et al. [6,7], provide a compendium of analytical solutions for ADE that are applicable to fully saturated porous media. Other works [8–10] have derived analytical and semi-analytical solutions in order to study solute transport in a fractured porous medium. With the goal of predicting transport in geological formations displaying heterogeneities, analytical solutions have also been applied to estimate the ensemble mean concentration (see [11,12] for details). Similarly, analytical solutions exist for contaminant mixing in rivers [13–16] and for bed-load sediment transport [17].

All of the above works rely on classical methods for solving linear partial differential equations, namely separation of variables and specific integral transforms [18–20]. However, as explained in [21], the traditional transforms have several disadvantages which will be discussed below using the diffusion equation as an illustrative example.

Consider the diffusion equation on the half-line, with Dirichlet boundary conditions:

$$\begin{aligned} &\frac{\partial \phi(\mathbf{x},t)}{\partial t} = \frac{\partial^2 \phi(\mathbf{x},t)}{\partial x^2}, \ \mathbf{0} < \mathbf{x} < \infty; \\ &\phi(\mathbf{x},\mathbf{0}) = \phi_o(\mathbf{x}), \ \mathbf{0} < \mathbf{x} < \infty; \ \phi(\mathbf{0},t) = g_o(t), \ t > \mathbf{0}, \end{aligned}$$

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where $\phi(x, t)$ denotes a scalar (e.g. temperature or concentration), $\phi_o(x)$ and $g_o(t)$ are given functions and $\phi_o(x) \rightarrow 0$ as $x \rightarrow \infty$. Using the sine transform, one finds the solution representation

$$\phi(\mathbf{x},t) = \frac{2}{\pi} \int_0^\infty \sin(\lambda \mathbf{x}) e^{-\lambda^2 t} \left[\int_0^\infty \sin(\lambda \xi) \phi_o(\xi) d\xi - \int_0^t e^{\lambda^2 \tau} g_o(\tau) d\tau \right] d\lambda.$$
(2)

- 1. The traditional approaches construct solutions which are not uniformly convergent at the boundaries unless the boundary conditions are homogeneous. In the particular case of Eq. (2), one cannot take the limit $x \rightarrow 0$ inside the integral otherwise the solution vanishes since $\sin(\lambda x)$ vanishes for x = 0.
- 2. Because of 1, it is not straightforward to verify that the function $\phi(x, t)$ defined by the right-hand side of (2) satisfies the boundary condition $\phi(0, t) = g_a(t)$.
- 3. Because of 1, the solution obtained from traditional methods is not directly suitable for numerical computations.
- 4. The traditional approaches require separability of the partial differential equation, domain and boundary conditions.
- 5. It is in general difficult to derive the proper transform.
- 6. Traditional transforms exist only for certain boundary value problems. For example, there does not exist an *x*-transform for the equation $\frac{\partial \phi(x,t)}{\partial t} + \frac{\partial^3 \phi(x,t)}{\partial x^2} = 0$, see [22,23].

The unified transform [21], also known as the Fokas method, has been used to obtain solutions for the diffusion equation [24–26], the reactive-diffusion equation [27], and recently has been applied to a physical problem with non-separable boundary conditions [28], as well as to interface problems [23,22]. However, the Fokas method has not been implemented to advective-dispersive (or convective-diffusive) problems.

In this work, we focus on solving the ADE via the unified transform. This work is structured as follows. Section 2 provides the formulation of the relevant physical-mathematical problem. The methodology is presented in Section 3 and specific semianalytical solutions are obtained in Section 4. Computational results are illustrated and discussed in Section 5 with particular emphasis on the speed and order of convergence of the method. Finally, a summary is provided in Section 6.

2. Problem formulation

Consider a one-dimensional semi-infinite medium with space and time variables *x* and *t*. The concentration of an inert solute is assumed to be governed by the ADE

$$\frac{\partial C(x,t)}{\partial t} + u \frac{\partial C(x,t)}{\partial x} = D \frac{\partial^2 C(x,t)}{\partial x^2}, \qquad 0 < x < \infty,$$
(3)

subject to the following boundary and initial conditions:

$$C(0,t) = g_o(t),\tag{4}$$

$$C(x,t) \to 0, \quad \text{as } x \to \infty, \quad t \ge 0$$
 (5)

$$C(x,0) = C_o(x), \tag{6}$$

where $C [M/L^3]$ denotes the concentration of an inert solute, u [L/T] represents the constant velocity and $D [L^2/T]$ is the dispersion coefficient. The problem defined by Eqs. (3)–(6) can represent solute transport in a river for which a time variable concentration $g_o(t)$ is injected at the origin at time zero. It can also represent transport of a non-reactive substance through a fully saturated soil column.

3. Methodology

Our starting point is the definition of the Fourier transform pair,

$$\widehat{C}(\lambda,t) = \int_{0}^{\infty} e^{-i\lambda x} C(x,t) dx, \qquad \operatorname{Im}\{\lambda\} \le 0,$$
(7)

$$C(x,t) = \frac{1}{2\pi} \int_{-\infty}^{\infty} e^{i\lambda x} \widehat{C}(\lambda,t) d\lambda \qquad 0 < x < \infty,$$
(8)

with ι denoting the imaginary unit and λ is a complex variable.

In order to obtain a solution to Eq. (3), we will proceed with the following methodological procedure proposed by Fokas [29,21].

- 1. Employ the Fourier transform to obtain the *global relation* (GR). This equation provides an expression that couples the function C(x, t) with its derivatives on the boundary of the domain.
- 2. Obtain an integral representation of the solution C(x, t) in the complex λ -plane.
- 3. By using the symmetry properties of the GR, eliminate from the integral representation of C(x, t) (obtained in Step 2) the transforms of the unknown boundary values.

Next, we employ the above steps to construct a solution for the ADE (3) subject to the initial and boundary conditions (4)–(6).

Employing Eq. (7) into (3) and integrating by parts we find the following GR:

$$e^{\left(\lambda^{2}D+\iota\lambda u\right)t}\widehat{C}(\lambda,t)=\widehat{C}_{o}(\lambda)-\widetilde{g}_{1}(\lambda,t)-\iota\lambda\widetilde{g}_{o}(\lambda,t),\qquad \mathrm{Im}\{\lambda\}\leq0,\quad(9)$$

where,

$$\widehat{C}_{o}(\lambda) = \int_{0}^{\infty} e^{-i\lambda x} C_{o}(x) dx, \quad \operatorname{Im}\{\lambda\} \le 0,$$
(10)

$$\widetilde{g}_{1}(\lambda,t) = \int_{0}^{t} e^{\left(\lambda^{2}D + i\lambda u\right)\tau} \left[D\left(\frac{\partial C}{\partial x}\right) \Big|_{x=0} - ug_{o}(\tau) \right] d\tau, \ \lambda \in \mathbb{C}, \ t > 0,$$
(11)

$$\widetilde{g}_{o}(\lambda,t) = D \int_{0}^{t} e^{\left(\lambda^{2}D + \iota\lambda u\right)\tau} g_{o}(\tau) d\tau, \quad \lambda \in \mathbb{C}, \, t > 0.$$
(12)

The derivation of (9) follows precisely the same steps used in the usual transform methods; details can be found in Appendix A.

Solving the GR (9) for $\hat{C}(\lambda, t)$ and employing the inverse Fourier transform (8), we find

$$C(x,t) = \frac{1}{2\pi} \int_{-\infty}^{\infty} e^{\imath \lambda x - (\lambda^2 D + \imath \lambda u)t} \widehat{C}_o(\lambda) d\lambda - \frac{1}{2\pi} \int_{-\infty}^{\infty} e^{\imath \lambda x - (\lambda^2 D + \imath \lambda u)t} [\widetilde{g}_1(\lambda, t) + \imath \lambda \widetilde{g}_o(\lambda, t)] d\lambda.$$
(13)

In summary, using the transform pair (7) and (8) and following the traditional transform methodology, we obtain an integral representation for the scalar quantity C(x, t). However, the representation (13) involves the *t*-transform of the unknown boundary value $\partial C/\partial x|_{x=0}$ defined by the function $\tilde{g}_1(\lambda, t)$ in (11). In order to eliminate this transform, we will adopt the two *novel* steps described below.

First, we deform the contour of the second integral in the righthand-side of Eq. (13) from the real axis to the upper half of the complex plane. In this connection, we need to determine the domain in the λ -complex plane where $\exp[\iota\lambda x - (\lambda^2 D + \iota\lambda u)(t - \tau)]$ is bounded (for $x \ge 0$ and $(t - \tau) \ge 0$). This can be achieved by analyzing the exponentials $\exp[\iota\lambda]$ and $\exp[-(\lambda^2 D + \iota\lambda u)]$: the exponential $\exp[\iota\lambda]$ is bounded for $\operatorname{Im}\{\lambda\} \ge 0$. We also note that $\lambda^2 D + \iota\lambda u = D(\lambda_R^2 - \lambda_I^2 + 2\iota\lambda_R\lambda_I) + \iotau(\lambda_R + \iota\lambda_I)$, where $\lambda = \lambda_R + \iota\lambda_I$. Thus, $\operatorname{Re}\{\lambda^2 D + \iota\lambda u\} = D(\lambda_R^2 - \lambda_I^2) - u\lambda_I$. Examining the curve

$$D(\lambda_R^2 - \lambda_I^2) - u\lambda_I = 0, \tag{14}$$

we note that if $\lambda_I = 0$, then $\lambda_R = 0$. Therefore, if $\lambda_R = 0$ then $\lambda_I = 0$ or $\lambda_I = -u/D$. Also, we observe that $\lambda_I^2 + (u/D)\lambda_I \ge 0$ or $\lambda_I(\lambda_I + u/D) = 0$, hence, either $\lambda_I \ge 0$ or $\lambda_I < -u/D$. Thus, we can replace the contour in the second term of the right-hand-side of (13) with the contour ∂D^+ in Fig. 1, where the curve ∂D^+ is defined by

$$D(\lambda_R^2 - \lambda_I^2) - u\lambda_I = 0, \quad \text{with } \lambda = \lambda_R + \iota\lambda_I.$$
 (15)

Both functions $\tilde{g}_1(\lambda, t)$ and $\tilde{g}_o(\lambda, t)$ (see Eqs. (11) and (12)) depend on λ only through the combination $\lambda^2 D + \iota \lambda u$. This fact motivates the search for a transformation $\lambda \rightarrow v$ such that the above combination is invariant:

$$\lambda^2 D + \iota \lambda u = v^2 D + \iota v u, \tag{16}$$

or

$$(\lambda^2 - \nu^2)D + \iota u(\lambda - \nu) = 0.$$
(17)

Hence,

$$(\lambda - \nu)[(\lambda + \nu)D + \iota u] = 0.$$
(18)

Thus, $v = \lambda$ or $v = -\lambda - u/D$. Replacing λ by $-\lambda - u/D$ in the global relation (9) we obtain:

$$e^{\left(\lambda^{2}D+\iota\lambda u\right)t}\widehat{C}\left(-\lambda-\iota\frac{u}{D},t\right)=\widehat{C}_{o}\left(-\lambda-\iota\frac{u}{D}\right)-\widetilde{g}_{1}(\lambda,t)+\left(\iota\lambda-\frac{u}{D}\right)\widetilde{g}_{o}(\lambda,t).$$
(19)

Solving (19) for \tilde{g}_1 , substituting the resulting equation in (13) with the integration in the second term replaced by the integration along the contour ∂D^+ , we obtain the equation

. (.2- .) ^

$$C(\mathbf{x},t) = \frac{1}{2\pi} \int_{-\infty}^{\infty} e^{i\lambda x - (\lambda^2 D + i\lambda u)t} C_o(\lambda) d\lambda$$

$$- \frac{1}{2\pi} \int_{\partial D^+} e^{i\lambda x - (\lambda^2 D + i\lambda u)t} \times$$

$$\left\{ \widehat{C}_o \left(-\lambda - \imath \frac{u}{D} \right) - \widehat{C} \left(-\lambda - \imath \frac{u}{D}, t \right) e^{(\lambda^2 D + i\lambda u)t} + \left(2\imath\lambda - \frac{u}{D} \right) \widetilde{g}_o(\lambda,t) \right\} d\lambda.$$
(20)

It is important to note that the unknown term $\widehat{C}(-\lambda - uu/D, t)$ does not contribute to the final expression, since both this term and exp $(\iota\lambda x)$ are bounded and analytic in D^+ . Furthermore, $\widehat{C}(-\lambda - uu/D, t)$ decays as $\lambda \to \infty$. Thus, according to Cauchy's theorem, the integral

$$\int_{\partial D^+} e^{\imath \lambda x} \widehat{C}\left(-\lambda - \imath \frac{u}{D}, t\right) d\lambda$$



Fig. 1. Boundaries ∂D^+ and ∂D^- of the domains D^+ and D^- which are defined by $D^{\pm} = \{\lambda = \lambda_R + \iota \lambda_I, \ \lambda_R \in \mathbb{R}, \ \lambda_I \in \mathbb{R}, \ \text{Re}\{\lambda^2 D + \iota u \lambda\} \leq 0\}.$

vanishes. Therefore, the general solution of the ADE (3) subject to the conditions (4)–(6) is given by

$$C(x,t) = \frac{1}{2\pi} \int_{-\infty}^{\infty} e^{\iota \lambda x - \left(\lambda^2 D + \iota \lambda u\right)t} \widehat{C}_o(\lambda) d\lambda - \frac{1}{2\pi} \int_{\partial D^+} e^{\iota \lambda x - \left(\lambda^2 D + \iota \lambda u\right)t} \left[\widehat{C}_o(-\lambda - \iota \frac{u}{D}) + \left(2\iota \lambda - \frac{u}{D}\right) \widetilde{g}_o(\lambda,t) \right] d\lambda,$$
⁽²¹⁾

where the contour ∂D^+ is depicted in Fig. 1 and defined by (15).

4. Specific solutions

We illustrate several features of the analytical expression derived in Section 3. Using Eq. (21) as a starting point, we develop solutions for two specific cases of increasing level of complexity.

4.1. Test Case 1 - Continuously injected solute at constant concentration

For our first demonstration, we set in Eqs. (4) and $(6)g_o(t) = C_i$ and $C_o(x) = 0$. This situation represents, for example, advectivedispersive transport of a solute through a homogeneous, fully saturated, porous medium where an inactive tracer is continuously injected at the origin with inlet concentration C_i , i.e., these conditions are met in laboratory scale tracer experiments in soil columns. Under the above conditions, Eq. (21) becomes

$$C(\mathbf{x},t) = -\frac{1}{2\pi} \int_{\partial D^+} e^{i\lambda \mathbf{x} - \left(\lambda^2 D + i\lambda u\right)t} \left(2i\lambda - \frac{u}{D}\right) \widetilde{g}_o(\lambda,t) d\lambda, \tag{22}$$

with

$$\widetilde{g}_{o}(\lambda, t) = C_{i}D \int_{0}^{t} e^{\left(\lambda^{2}D + i\lambda u\right)\tau} d\tau$$

$$= C_{i}D \left[\frac{e^{\left(m\lambda + D\lambda^{2}\right)t} - 1}{\lambda(m + D\lambda)}\right].$$
(23)

Using (22) and (23), the equation satisfied by the scalar C(x,t) becomes

$$C(\mathbf{x},t) = -\frac{C_i D}{2\pi} \int_{\partial D^+} e^{i\lambda x - \left(\lambda^2 D + i\lambda u\right)t} \frac{\left(2i\lambda - \frac{u}{D}\right)}{\lambda(iu + D\lambda)} \left[e^{\left(\lambda^2 D + i\lambda u\right)t} - 1 \right] d\lambda, \quad (24)$$

which can be further simplified to

$$C(x,t) = -\frac{iC_i}{2\pi} \int_{\partial D^+} \left(\frac{2D\lambda + iu}{D\lambda + iu} \right) e^{i\lambda x} \frac{\left[1 - e^{-\left(\lambda^2 D + i\lambda u\right)t} \right]}{\lambda} d\lambda,$$
(25)

where the integration over ∂D^+ will be carried out numerically. In this connection, it is convenient (using Cauchy's theorem) to replace ∂D^+ by the contour *h* depicted below (see Fig. 2)). This contour has the advantage that both $\exp[\iota\lambda x]$ and $\exp[-(\lambda^2 D + \iota\lambda u)t]$ (which oscillates on the contour ∂D^+) decay on the contour *h*.

In summary, Eq. (25) is the solution for the initial boundary value problem subject to $g_o(t) = C_i$ and $C_o(x) = 0$ and will be used



Fig. 2. Schematic of the contour h.

to compute the concentration field. Additional details regarding the numerical integration are provided in Section 5. It is shown in Appendix B that Eq. (25) can be simplified further.

The problem under consideration, i.e. $g_o(t) = C_i$ and C(x, 0) = 0, has been tackled in the past by Ogata and Banks [2]. In order to obtain a solution for this problem, Ogata and Banks [2] mapped the ADE into a diffusion equation and then made use of Duhamel's theorem [18] to derive the following *approximate* analytical expression:

$$C(x,t) \approx \frac{C_i}{2} \left\{ \text{Erfc}\left[\frac{x-ut}{2\sqrt{Dt}}\right] + e^{\frac{ux}{D}} \text{Erfc}\left[\frac{x+ut}{2\sqrt{Dt}}\right] \right\}.$$
(26)

We will utilize Eq. (26), which represents an approximate solution for the concentration field, to benchmark the results obtained through the unified transform (computed by using Eq. (25)).

4.2. Test Case 2 – Time variable boundary condition and spatially distributed initial condition

Next, we derive the solution for the following case: $g_o(t) = \mathcal{A}_i[\cos(\omega t) + 1]$ and $C_o(x) = \exp[-x]$ (see Eqs. (4) and (6)), where \mathcal{A}_i and ω are amplitude and angular frequency respectively. Under this condition, both $\hat{C}_o(\lambda)$ and $g_o(t)$ (see Eq. (21)) can be computed analytically. Explicit expressions for $\hat{C}_o(\lambda)$, $\hat{C}_o(-\lambda - iu/D)$ and $g_o(t)$ are provided below:

$$\widehat{C}_{o}(\lambda) = -\frac{1}{\lambda - 1},\tag{27}$$

$$\widetilde{C}_{o}\left(-\lambda - \iota \frac{u}{D}\right) = \frac{iD}{\iota u + D(\iota + \lambda)}$$

$$\widetilde{g}_{o}(\lambda, t) = D\mathcal{A}_{i} \left\{ \frac{e^{\lambda t(u+D\lambda)} - 1}{u - 1} \right\} +$$
(28)

$$\left\{\frac{e^{\lambda t(u+D\lambda)}[\lambda(u+D\lambda)]}{\lambda^2(u+D\lambda)^2+\omega^2}\right\}$$
(2)
$$\left\{\frac{e^{\lambda t(u+D\lambda)}[\lambda(u+D\lambda)\cos(\omega t)+\omega\sin(\omega t)]-\lambda(u+D\lambda)}{\lambda^2(u+D\lambda)^2+\omega^2}\right\}$$

Substituting these into (21) leads to

$$C(x,t) = \frac{1}{2\pi} [\mathcal{P}_1(x,t) - \mathcal{P}_2(x,t)],$$
(30)

where

$$\mathcal{P}_{1}(\mathbf{x},t) = -\int_{-\infty}^{\infty} e^{i\lambda x - \left(\lambda^{2}D + i\lambda u\right)t} \left(\frac{1}{\lambda - 1}\right) d\lambda; \qquad (31)$$

$$\mathcal{P}_{2}(\mathbf{x},t) = \left\{ \mathcal{A}_{i} \int_{\partial D^{+}} e^{i\lambda x - \left(\lambda^{2}D + i\lambda u\right)t} \left(2i\lambda D - u\right) \left[\frac{e^{it(u+D\lambda)} - 1}{\lambda(u+D\lambda)}\right] d\lambda + \int_{\partial D^{+}} e^{i\lambda x - \left(\lambda^{2}D + i\lambda u\right)t} \left[\frac{Di}{|u+D(t+\lambda)} + \mathcal{A}_{i}(2i\lambda D - u) \times \frac{e^{it(u+D\lambda)} \left[\lambda(u+D\lambda)\cos(\omega t) + \cos\sin(\omega t)\right] - \lambda(u+D\lambda)}{\lambda^{2}(u+D\lambda)^{2} + \omega^{2}} \right] d\lambda \right\}.$$

In summary, Eq. (30) provides the solution of the initial boundary value problem with $g_o(t) = \mathcal{A}_i[\cos(\omega t) + 1]$ and $C_o(x) = \exp[-x]$. Similar to Test Case 1, the contours in (31) and (32) can be replaced by the contour *h* depicted in Fig. 2 which is used in the numerical evaluations and defined in the next section.

5. Computational results

For the upcoming computational illustrations, we plot the results for Test Cases 1 and 2 in dimensionless form. The following dimensionless quantities were adopted for the concentration C, distance x and time t respectively:

$$\chi = \frac{C}{C_i}; \quad \xi = \frac{xu}{D}; \quad \eta = \frac{u^2 t}{D}, \tag{33}$$

with C_i denoting a characteristic concentration value at the inlet boundary.

The analytic solution of the solute concentration (after a suitable parametrization) can be represented in terms of integrals of the following two forms:

$$L_1(f,\alpha) = \int_0^\infty f(x)e^{-\alpha x}dx, \quad L_2(f,\alpha) = \int_{-\infty}^\infty f(x)e^{-\alpha x^2}dx, \quad (34)$$

where the functions *f* are well behaved (smooth and bounded). In the case that α is not small, we shall evaluate these integrals using Gauss-Laguerre quadrature for L_1 and Gauss-Hermite quadrature for L_2 . For the functions considered in this paper these quadrature rules will converge exponentially (we refer the reader to [30–32] for the theory of these quadrature rules) with the number of evaluations, which we denote throughout by N. These quadrature rules (the weights and nodes) can also be computed extremely fast in $\mathcal{O}(N)$ operations [33–35]. Similarly, if α is small (as in the case of (25) for small x) we shall truncate the domain of integration (typically to $|\lambda| < 100$) and apply standard Gaussian quadrature which will also converge exponentially and have $\mathcal{O}(N)$ cost. In all cases (assuming the functions \hat{C}_0 and \tilde{g}_0 are known), we recover a quadrature method which can be implemented in $\mathcal{O}(N)$ operations per point (x, t) which will also converge pointwise exponentially in N. Note that this is a key advantage of the unified transform over classical methods as mentioned in the introduction. We will also consider the case where \hat{C}_o and \tilde{g}_o themselves must be computed in Section 5.3.

All computational results were obtained using MATLAB on a standard desktop computer and to compare we have computed accurate reference solutions using adaptive Gauss-Kronrod quadrature to evaluate the integrals in the unified transform (with tolerance set to machine precision). For the numerical evaluations of the concentration field through Eq. (21), the numerical computation of the integrals in (25) and (30) is achieved by deforming ∂D^+ to a contour *h* between the real line and ∂D^+ (see Fig. 2)). A convenient choice for *h* is the union of the two line segments

$$\{\ell[\cos\left(\pi-\theta\right)+\iota\sin\left(\pi-\theta\right)]:\ell\ge 0\},\tag{35}$$

$$\{\ell[\cos\left(\theta\right) + \iota\sin\left(\theta\right)] : \ell \ge 0\},\tag{36}$$

where θ is the angle between the line segment and the real line. For our illustrations, we set $\theta = \pi/6$.

5.1. Results for Test Case 1

As a first example, we consider the computation of the solution for $\eta = 16$. Fig. 3 shows the computed solution using Eq. (B11) with N = 50 quadrature points, denoted as Approach 1 in the plot, and the approximate analytical solution (26) obtained through traditional methods. For completeness, we also include a comparison with a numerical solution based on the Crank-Nicolson and backward Euler finite-difference methods with $\Delta t = \Delta x = 0.5$. In addition, Fig. 3 compares with a solution, also based via the unified transform, obtained by mapping the ADE into a diffusion equation (denoted as Approach 2 in the plot). The unified transform-based solution for the diffusion equation is well documented in the literature [24,21] and for the sake of brevity, it is not reproduced here. All of the methods agree well.

Fig. 4 shows the convergence of the method for different η as we increase *N*, where we have approximated the relative error measured in L^2 norm (i.e. normalised by the L^2 norm of the reference solution computed using adaptive quadrature). As expected, we see exponential convergence in *N*, with the simplified version requiring smaller *N* (see Fig. 4, left). Interestingly, whilst we need larger *N* for larger times to achieve modest errors, there seems to be a cut-off beyond which the errors depend similarly with respect to *N* independent of η . To verify our solution, we have also shown the error when using the Crank-Nicolson and backward Euler



Fig. 3. Dimensionless concentration $\chi(\xi, \eta)$ versus distance ξ at $\eta = 16$. Comparison between the proposed analytical solution with other analytical and numerical solutions. Approach 1 refers to the solution given by (B11) (which can also be obtained via (25)) for N = 50. Approach 2 refers to the solution obtained through the unified transform method by mapping the ADE into a diffusion equation. We have also shown the solutions obtained using Crank-Nicolson and backward Euler finite-difference methods with $\Delta x = 0.5$ (note the increased accuracy in using Crank-Nicolson).

finite-difference methods (see Fig. 4, right). For the finitedifference methods, we note that one needs to solve a triangular $M \times M$ system at each step where $M \approx \mathcal{O}(\Delta x^{-1})$. This can be done in $\mathcal{O}(M)$ operations as opposed to $\mathcal{O}(M^3)$ for a full system but still gives $\mathcal{O}(\Delta x^{-1}\Delta t^{-1})$ operations to compute the solution at any fixed time *t* and the methods only converge algebraically. This shows a significant advantage of the unified transform: as well as obtaining the solution at any time without time-stepping, the integral representation lends itself to extremely efficient numerical integration.

Next, we consider the computation of the solution over a range of values of ξ and η . Fig. 5 (top) shows the concentration profile computed using the unified transform. As time increases, the front of the solute cloud advances through the flow domain and becomes more diffused with time. We have also shown the absolute error of the method with N = 250 using both (25) and (B11) (see Fig. 5 bottom left and bottom right respectively). The method is very fast, taking approximately 0.29 s using (25) and 0.08 s using (B11) to compute the profile over 4800 points. Using (25) is slower due to the need to compute two integrals. In comparison, the reference solution was computed in 19 s (with adaptive quadrature). Both sets of errors approach machine precision over a large portion of the domain. Using (B11) appears to struggle near the singularity at $(\xi, \eta) = (0, 0)$ (this is to be expected from the blow-up of the integrand) and (25) appears to struggle more for small ξ , particularly at larger η due the need to evaluate a slowly (but still exponentially) decaying oscillatory integrand. We found that increasing *N* allows more accurate solutions to be computed near the boundary. We also remark that the concentration profile computed through the unified transform almost perfectly matches the solution computed through different approaches such as the approximate solution given by (26)[2]. The approximation (26) was found to agree to about 12 decimal places in the domain interior.

Finally, Fig. 6 shows the value of *N* needed in order to ensure five digits of accuracy, as a function of ξ . This shows that accurate solutions can be computed extremely efficiently, especially with the integral form (B11).

5.2. Results for Test Case 2

Now we present the results obtained from Eq. (30) for an oscillatory boundary condition and an exponentially distributed initial condition. For this analysis, we set $\omega = 1$ and $A_i = C_i/2$ in equations (27) and (29). Fig. 7 shows the concentration as a function of distance for different times ($\eta = 5$, 10, 20 and 30). In addition, Fig. 7 compares the results computed from the unified transform (30) with the numerical solution of the ADE subject to the conditions described in Section 4.2. The numerical solution was computed using MATLAB's built-in function pdepe which runs a variable time-stepping method, second order in the spatial resolution [36]. As shown in Fig. 7, both methods agree well.

Fig. 8 shows the convergence of the unified transform as we increase *N*. In the results depicted in Fig. 8 (left), we see exponential convergence as before. For completeness, we have also shown the convergence of pdepe as $\Delta x \downarrow 0$ (Fig. 8, right). As expected, the convergence of pdepe is algebraic with error of order $\mathcal{O}(\Delta x^2)$, however, this still serves to verify the computation of the solution using the unified transform. Finally, we have shown the solution over a range of values of ξ and η , as well as the absolute error of the unified transform over the domain for N = 250 in Fig. 9. As before, we see errors close to machine precision over the domain, with larger errors nearer the boundary $\xi = 0$. Again the method was very fast, taking 0.85s in total for evaluation at all 6000 points to produce Fig. 9.

5.3. An example where the integral transforms cannot be computed analytically

As a final example, we will consider the case where the functions \widehat{C}_o and \widetilde{g}_o themselves are computed numerically. Note that we only need to compute these at $\mathcal{O}(N)$ points. We will use



Fig. 4. Left: Exponential convergence of the unified transform as measured via a relative L^2 norm. We have shown convergence using both (25) and the simplified version (B11). Right: Algebraic convergence of Crank-Nicolson (CN, order 2) and backward Euler (BE, order 1) finite-difference methods.



Fig. 5. Top: Concentration profile. Bottom left: Absolute error when using (25). Bottom right: Absolute error when using (B11).



Fig. 6. Left: Value of N needed for five digits of accuracy when using (25). Right: Value of N needed for five digits of accuracy when using (B11).

Gaussian quadrature with M_1 nodes to compute \tilde{g}_o . Suitable quadrature rules for the computation of \hat{C}_o depend on the decay properties of $C_o(x)$ as $x \to \infty$. We will consider the case of

 $C_o(x) = 2\cos(x)\exp(-x).$

Note that \hat{C}_o can be computed analytically, but instead we shall use Gaussian-Laguerre quadrature with M_2 nodes. Altogether, this leads

to a method with computational cost $\mathcal{O}(N(M_1 + M_2))$. In general, the error of such a method will depend on the error of the computation of the integrals (a full discussion of which is beyond the scope of this paper, but there is a large literature on the convergence methods of different quadrature rules). For this example, we will take

$$g_o(t) = J_0(t) + 1,$$



Fig. 7. Dimensionless concentration $\chi(\xi, \eta)$ versus dimensionless distance ξ for $\eta = 5, 10, 20$ and 30. Comparison between the proposed analytical solution with the fully numerical solution (computed using $\Delta x = 0.5$).

where J_0 denotes the Bessel function of the first kind of order 0. For these choices of functions, the absolute error in the computation of the spectral functions decreases exponentially in M_1 , M_2 .

Fig. 10 (left) shows the exponential convergence of the method for $M_1 = N$ and $M_2 = \text{round}(2N/5)$. For larger times, a larger value of *N* is needed. This is because a larger value of M_2 is needed for the quadrature to compute \tilde{g}_o via integrating g_o over a larger interval. We have also shown the solution computed over a grid of 1200 points in Fig. 10 (right), with N = 600, $M_1 = 600$ and $M_2 = 240$, taking 11s with a maximum absolute error bounded by $\approx 10^{-14}$. We verified this against the solution using pdepe where the solutions agreed to about five decimal places. However, the unified transform is able to gain much more accurate solutions. Finally, to examine how the error depends on varying M_1 and M_2 , Table 1 shows the average absolute error (over the same computational grid) with N = 500 and various M_1 and M_2 . We see that large M_1 is more critical to computing very accurate solutions as expected due to the exponential decay of C_o .

6. Summary

We have implemented the unified transform to obtain the solution of the advection-dispersion equation on a half-line with a Dirichlet boundary condition (constant and time varying). The unified transform method contains classical approaches as special cases. In particular, we highlight the following key advantages of this new method: (1) it allows one to construct a representation of the solution which is uniformly convergent at the boundaries; (2) it can be implemented numerically in a straightforward manner; and (3) the unified transform is computationally efficient for large times, since it does not require discretization of the time domain. The solution obtained by the new method is in excellent agreement with the computational results obtained through numerical methods and other analytical solutions. In particular, it gives rise to a numerical method converging exponentially as the number of evaluations N increases and only has $\mathcal{O}(N)$ complexity.

Although the solution was obtained for advective-dispersive problems relevant for environmental engineering and hydrological applications, it can be employed to study a series of other applications regarding heat and mass transfer in mechanical, chemical,



Fig. 8. Left: Exponential convergence of the unified transform as measured via a relative L² norm for Case 2. Right: Algebraic convergence of MATLAB's pdepe (order 2).



Fig. 9. Left: Concentration profile. Right: Absolute error for N = 250.



Fig. 10. Left: Exponential convergence of the unified transform with increasing *N* for the case of using quadrature to evaluate the spectral functions. Right: Computed solution over various ξ and η . The maximum absolute error was bounded by $\approx 10^{-14}$.

Table 1 Average absolute error over the computational grid for various M_1 and M_2 .

<i>M</i> ₂	M_1				
	100	200	300	400	500
40	$3.15\times\mathbf{10^{-7}}$	8.26×10^{-10}	1.62×10^{-10}	1.60×10^{-10}	1.60×10^{-10}
80	3.14×10^{-7}	6.66×10^{-10}	2.55×10^{-12}	4.39×10^{-14}	1.35×10^{-14}
120	3.14×10^{-7}	$6.66 imes 10^{-10}$	2.54×10^{-12}	3.27×10^{-14}	2.41×10^{-15}
160	3.14×10^{-7}	6.66×10^{-10}	2.54×10^{-12}	3.27×10^{-14}	2.37×10^{-15}

petroleum and biomedical engineering [e.g., 37]. The solution obtained through the unified transform is well-suited to compute the concentration field from laboratory tracer experiments in soil columns where, in general, the velocity and dispersion coefficient are homogeneous. It is important to note that it is straightforward to implement the unified transform to problems posed on a finite interval [38,39]. Also, a major advantage of the unified transform is its applicability to a wide range of boundary conditions including Neumann, Robin and non-local boundary conditions. It can also be applied to some problems where classical methods apparently fail [40,41] and PDEs involving third order spatial derivatives (see point 6 in the introduction). Multidimensional applications and extensions to reactive transport will be presented in future work.

Finally, it should be emphasised that the unified transform is the only existing method which expresses the solution of linear evolution PDEs in terms of an integral involving a contour in the λ -complex plane. This opens the way for an efficient numerical evaluation of the solution. In this paper, for the first time, the advantage of the numerical implementation of the unified transform is elucidated.

Declaration of Competing Interest

None.

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Appendix A

Multiplying Eq. (3) with $e^{-i\lambda x}$ and integrating with respect to *x*, we obtain

$$\frac{\partial}{\partial t} \int_0^\infty e^{-\imath \lambda x} C(x,t) dx + \int_0^\infty e^{-\imath \lambda x} u \frac{\partial C(x,t)}{\partial x} dx = \int_0^\infty e^{-\imath \lambda x} D \frac{\partial^2 C(x,t)}{\partial x^2} dx.$$
(A1)

This equation can be further simplified to

$$\frac{\partial \widehat{C}(\lambda,t)}{\partial t} + u \int_0^\infty e^{-\iota\lambda x} \frac{\partial C(x,t)}{\partial x} dx = D \int_0^\infty e^{-\iota\lambda x} \frac{\partial^2 C(x,t)}{\partial x^2} dx.$$
(A2)

Integration by parts of the second term in the left-hand side of (A2) yields

$$u \int_0^\infty e^{-\imath\lambda x} \frac{\partial C(x,t)}{\partial x} dx = u \Big[-C(0,t) + \imath\lambda \widehat{C}(\lambda,t) \Big].$$
(A3)

For the right-hand side of (A2), we obtain the following expression after integration by parts:

$$D\int_{0}^{\infty} e^{-\imath\lambda x} \frac{\partial^{2} C(x,t)}{\partial x^{2}} dx = D \left[-\frac{\partial C(x,t)}{\partial x} \right|_{x=0} + \imath\lambda \int_{0}^{\infty} e^{-\imath\lambda x} \frac{\partial C(x,t)}{\partial x} dx \right].$$
(A4)

Integrating by parts again, we obtain

$$D\int_{0}^{\infty} e^{-\iota\lambda x} \frac{\partial^{2} C(x,t)}{\partial x^{2}} dx = D \left[-\frac{\partial C(x,t)}{\partial x} \Big|_{x=0} - \iota\lambda C(0,t) - \lambda^{2} \widehat{C}(\lambda,t) \right].$$
(A5)

Substituting Eqs. (A3) and (A5) into (A2) yields

$$\frac{\partial \widehat{C}(\lambda, t)}{\partial t} + (\lambda^2 D + \iota u \lambda) \widehat{C}(\lambda, t) = -\left[D \frac{\partial C(x, t)}{\partial x} \Big|_{x=0} - u C(0, t) \right]$$
(A6)
$$- \iota D \lambda C(0, t).$$

The above expression can be re-written as

$$\frac{\partial}{\partial t} \left[e^{\left(\lambda^2 D + i\lambda u\right)t} \widehat{C}(\lambda, t) \right] = - \left[D \frac{\partial C(\mathbf{x}, t)}{\partial \mathbf{x}} \Big|_{\mathbf{x} = 0} - u C(0, t) \right] e^{\left(\lambda^2 D + i\lambda u\right)t}$$
(A7)
$$- i D \lambda C(0, t) e^{\left(\lambda^2 D + i\lambda u\right)t}.$$

Integrating in time, Eq. (A7) becomes

$$e^{\left(\lambda^{2}D+\iota\lambda u\right)t}\widehat{C}(\lambda,t)-\widehat{C}(\lambda,0) = -\int_{0}^{t} \left[D\frac{\partial C(x,\tau)}{\partial x} \Big|_{x=0} - uC(0,\tau) \right] e^{\left(\lambda^{2}D+\iota\lambda u\right)\tau} d\tau$$

$$-\iota D\lambda \int_{0}^{t} C(0,\tau) e^{\left(\lambda^{2}D+\iota\lambda u\right)\tau} d\tau.$$
(A8)

Recalling that $C(0,t) = g_o(t)$, Eq. (A8) can be written in the more compact form

$$e^{\left(\lambda^{2}\mathsf{D}+\imath\lambda u\right)t}\widehat{C}(\lambda,t)=\widehat{C}_{\mathfrak{o}}(\lambda)-\widehat{g}_{1}(\lambda,t)-\imath\lambda\widehat{g}_{\mathfrak{o}}(\lambda,t),\tag{A9}$$

with $\hat{C}_o(\lambda)$, $\hat{g}_1(\lambda, t)$ and $\hat{g}_o(\lambda, t)$ defined in Eqs. (10)–(12). Eq. (A9) is identical to the global relation (9).

Appendix **B**

Analyzing Eq. (25), it is clear that $\lambda = 0$ is a removable singularity:

$$\lim_{\lambda \to 0} \left[\frac{1 - e^{-(\lambda^2 D + i\lambda u)t}}{\lambda} \right] = iut$$

Thus, before we split (25) with the goal of obtaining a simpler expression, we first deform ∂D^+ to go above $\lambda = 0$ (i.e. bypassing the singularity). We call this new contour *L*.

In order to simplify the above solution (25), we divide C(x, t) into two terms

$$C(x,t) = C_1(x) + C_2(x,t),$$
 (B1)

with

$$C_1(\mathbf{x}) = -\frac{\mathbf{i}C_i}{2\pi}I_1(\mathbf{x}); \tag{B2}$$

$$C_{2}(x,t) = \frac{iC_{i}}{2\pi} I_{2}(x,t),$$
(B3)

where $I_1(x)$ and $I_2(x, t)$ are given by

$$I_1(\mathbf{x}) = \int_L \left(\frac{2D\lambda + u}{u + D\lambda}\right) \frac{e^{i\lambda \mathbf{x}}}{\lambda} d\lambda, \tag{B4}$$

and
$$I_2(x,t) = \int_L \left(\frac{2D\lambda + u}{u + D\lambda}\right) \frac{e^{i\lambda x}}{\lambda} e^{-(\lambda^2 D + i\lambda u)t} d\lambda.$$
 (B5)

We can use Cauchy's theorem in the domain above *L*, where neither $\lambda = 0$ nor $\lambda = -ut/D$ are in this domain. As a consequence, we find $I_1(x) = 0$. In other words, $C_1(x) = 0$.

In order to simplify I_2 and compute $C_2(x, t)$, we observe that the argument of the exponential term in (B5) can be re-written as follows:

$$-(\lambda^{2}D + \iota\lambda u)t + \iota\lambda x = -tD[\lambda^{2} + \frac{\iota\lambda}{D}(u - \frac{x}{t})]$$

$$= -tD[\lambda^{2} + \frac{\iota\lambda}{D}v]$$

$$= -tD[\lambda^{2} + \frac{\iota\lambda}{D}v - (\frac{v}{2D})^{2} + (\frac{v}{2D})^{2}]$$

$$= -tD(\lambda + \iota\frac{v}{2D})^{2} - t\frac{v^{2}}{4D},$$

(B6)

with v := u - x/t. Therefore, $C_2(x, t)$, see Eq. (B3), becomes

$$I_{2}(x,t) = e^{-t\frac{y^{2}}{4D}} \int_{L} \left(\frac{2D\lambda + uu}{D\lambda + uu} \right) \frac{e^{-tD\left[\lambda + \frac{y^{2}}{2D}\right]^{2}}}{\lambda} d\lambda.$$
(B7)

If we employ the change of variables $\Lambda = \lambda + \iota v / (2D)$, the expression for I_2 can be re-written in the form

$$I_{2}(\mathbf{x},t) = e^{-t_{\overline{4D}}^{\nu^{2}}} \int_{\widetilde{L}} \left[\frac{2D\Lambda + \iota(u-\nu)}{D\Lambda + \iota(u-\frac{\nu}{2})} \right] \frac{e^{-tD\Lambda^{2}}}{\Lambda - \iota\frac{\nu}{2D}} d\Lambda,$$
(B8)

where the contour \tilde{L} is depicted in Fig. 11.



Fig. 11. Mapping the singularity $\lambda = 0$ to the point $\Lambda = vv/(2D)$, with v := u - x/t, in the Λ complex plane with the new contour \tilde{L} .

Comparing Eqs. (B7) and (B8), it follows that the point $\lambda = 0$ is mapped to the point $\Lambda = i\nu/(2D)$. Thus, we need to distinguish the following two cases: (1) u > x/t, and (2) u < x/t. These two cases are depicted in Fig. 11 where the new contour \tilde{L} is either above the real axis of the Λ plane if u > x/t, or below if u < x/t. In the second case, we can deform the contour \tilde{L} to the real axis; however, in the first case, we also obtain a pole contribution from $\Lambda = i\nu/(2D)$.

$$I_{2}(x,t) = \begin{cases} e^{-t\frac{u^{2}}{4D}} \widetilde{I}_{2}(x,t) &, \text{ if } u < \frac{x}{t} \\ e^{-t\frac{u^{2}}{4D}} \widetilde{I}_{2}(x,t) - 2i\pi &, \text{ if } u > \frac{x}{t} \end{cases}$$
(B9)

with

$$\widetilde{I}_{2}(x,t) \doteq \int_{-\infty}^{+\infty} \left[\frac{2D\Lambda + \iota(u-\nu)}{D\Lambda + \iota(u-\frac{\nu}{2})} \right] \frac{e^{-tD\Lambda^{2}}}{\Lambda - \iota\frac{\nu}{2D}} d\Lambda.$$
(B10)

Note that $\Lambda = -i(u - v/2)/D = -i(u + x/t)/(2D)$ never contributes to I_2 since it is below the real axis.

Making use of Eqs. (B3), (B9), (B10) and substituting into (B1) yields

$$C(x,t) = \begin{cases} \frac{iC_{i}}{2\pi} e^{-t\frac{y^{2}}{4D}} \int_{-\infty}^{+\infty} \left[\frac{2DA+i(u-v)}{DA+i(u-\frac{v}{2})} \right] \frac{e^{-tDA^{2}}}{A-\frac{v}{2D}} dA &, \text{ if } u < \frac{x}{t}; \\ \frac{iC_{i}}{2\pi} \left[-2i\pi + e^{-t\frac{y^{2}}{4D}} \int_{-\infty}^{+\infty} \left[\frac{2DA+i(u-v)}{DA+i(u-\frac{v}{2})} \right] \frac{e^{-tDA^{2}}}{A-\frac{v}{2D}} dA \right], & \text{ if } u > \frac{x}{t}. \end{cases}$$
(B11)

In summary, Eq. (25) can be re-written in the form (B11). To make this suitable for quadrature, we let $\tau = \sqrt{tD}A$ and rewrite the integral in the dimensionless form

$$\widetilde{I}_{2}(x,t) = \int_{-\infty}^{+\infty} \left[\frac{2\tau + \iota \frac{\zeta}{\sqrt{\eta}}}{\tau + \iota \frac{\sqrt{\eta} + \frac{\zeta}{\sqrt{\eta}}}{2}} \right] \frac{e^{-\tau^{2}}}{\tau - \iota \frac{\sqrt{\eta} - \frac{\zeta}{\sqrt{\eta}}}{2}} d\tau.$$
(B12)

For η close to ξ , we deform this contour into the complex plane to avoid the pole.

Appendix C. Supplementary material

Supplementary data associated with this article can be found, in the online version, at https://doi.org/10.1016/j.ijheatmasstransfer. 2019.05.018.

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