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One-dimensional turbulence (ODT): computationally efficient modeling and simulation of turbulent flows --Manuscript Draft--

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One-dimensional turbulence (ODT): computationally efficient modeling and simulation of turbulent flows

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Abstract

One-dimensional turbulence (ODT) is an accurate and computationally efficient technique for modeling turbulent flows. ODT has been applied to a wide range of flow problems including reaction, multiphase, differential diffusion, heat release, buoyancy, and wall flows. Applications include use as a standalone model and as a closure for large-eddy simulation (LES). Its strength lies in the ability to capture a full range of turbulent length and time scales. The ODT model is strongly coupled with its implementation, complicating its formulations. We present a modern, open-source, object-oriented C++ implementation of ODT. This code can be used as a starting point to understand, apply, and extend the ODT model, enabling its further application to turbulent flow research.

Key words: turbulence, reacting flows, one-dimensional turbulence

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Code Metadata

Nr.	Code metadata description	Please fill in this column
C1	Current code version	1.0
C2	Permanent link to code/repository	https://github.com/BYUignite/ODT
	used for this code version	
С3	Code Ocean compute capsule	N/A
C4	Legal Code License	MIT
C5	Code versioning system used	Git
C6	Software code languages, tools, and	C++, Python 3.x, YAML,
	services used	
C7	Compilation requirements, operat-	CMake 3.12+, Cantera, Git, Doxy-
	ing environments & dependencies	gen (optional)
C8	If available Link to developer docu-	ODT style guide
	mentation/manual	
С9	Support email for questions	davidlignell@byu.edu

Table 1: Code metadata (mandatory)

1. Motivation and significance

Turbulent flows characterize the vast majority of fluid flows in practical engineering applications, and simulations of turbulent flows provide researchers with valuable insights into complex systems, particularly reacting turbulent flows such as combustion processes. Turbulence is a complex phenomenon that affects the full range of a flow's length and time scales. As a result, resolving the entire flow field by numerically solving the Navier-Stokes equations of fluid flow, as is done in direct numerical simulations (DNS), requires substantial computational resources. DNS is a powerful research tool, but its high computational cost makes it intractable for simulating most practical engineering flows. In order to achieve numerical solutions to practical flow problems, researchers use alternative frameworks that model turbulence rather than resolving it directly.

Large-eddy simulations (LES) address the problem of wide-ranging length and time scales by combining direct resolution of grid-scale quantities, as in DNS, with subgrid modeling of smaller turbulence structures. The more complex the flow, the more modeling is required; for example, a jet flame simulation might require subgrid modeling for the combustion chemistry, radiative heat transfer, and soot chemistry in addition to turbulence structures, all of which form a tightly coupled system in which each model interacts heavily with the others. While subgrid modeling makes LES more computationally

affordable than DNS, it can introduce empiricism into simulations, which can lead to inaccurate results. Additionally, unresolved quantities are often parameterized in state space with empirical relationships or assumed distributions that lack universal applicability. LES is a valuable simulation tool, but its approach to turbulence modeling can introduce unwanted empiricism and make errors difficult to isolate and quantify.

The one-dimensional turbulence model (ODT) functionally reverses the LES approach, modeling large-scale turbulent advection and directly resolving small-scale flow structures, simulating the full range of length and time scales in a single dimension. Depending on the configuration, large-scale structures can be easier to model than small-scale structures, especially in reacting flows. ODT's direct resolution of fine scales in one dimension bypasses subgrid modeling issues that complicate LES. Previous studies show that ODT can attain accuracy comparable to DNS at a fraction of the computational cost [1, 2], making it an attractive tool for simulating turbulent flows. Because the ODT model is one-dimensional, it is limited to homogeneous or boundary-layer flows, such as jets, wakes, and mixing layers, but these types of flows are common in nature and central to turbulence research. ODT's computational efficiency and resolution of a full range of scales make it a valuable tool that complements experimental studies and other simulation tools like DNS and LES.

Early applications of ODT focused on homogeneous turbulence, wakes, and mixing layers [3–5]. Later extension to variable-density flows and a spatial downstream coordinate system facilitated its growth and application to more complex flows, including combustion in jet flames [6–13], counterflow flames [14], wall fires [15], and sooting flames [1, 16–19], as well as other particle flows [20–23]. ODT has also served to complement LES through subgrid modeling studies [24–26] and has been applied to various other flow configurations such as double-diffusive interfaces [27], Rayleigh-Taylor mixing [28], and stratified turbulence [29]. Most recently, the ODT code was extended to include cylindrical and spherical coordinate systems [30–32].

During the recent implementation of the cylindrical and spherical model formulations, the ODT code was drastically overhauled and reorganized, resulting in its current configuration. The ODT code presented here is a pared down version of the development code, representing the fundamental aspects of the ODT model and its most reliable functions. The example case in Section 3 is a representative sample of the ODT code's capabilities as it is presented here. Future releases will expand this code's functionality with additional features currently in development.

2. Software description

2.1. Model description

The ODT model is described in detail in the literature [3, 5, 30, 33, 34]; only a brief explanation will be given here. In ODT, turbulent advection is modeled with stochastic processes called eddy events, which punctuate the solution of unsteady, one-dimensional transport equations for mass, momentum, and enthalpy. The ODT code uses a Lagrangian finite-volume formulation for diffusive advancement in which mass stays constant within each grid cell while cell volumes increase or decrease according to cell dilation via an adaptive mesh refinement [34].

Transport equations for mass, momentum, and enthalpy in the temporal formulation of ODT take the following generic form, derived from the Reynolds Transport Theorem [35] for a given scalar quantity per unit mass β :

$$\frac{\mathrm{d}\beta}{\mathrm{d}t} = -\frac{j_{\beta,e}A_{x,e} - j_{\beta,w}A_{x,w}}{\rho V} + \frac{S_{\beta}}{\rho V}.$$
 (1)

Here, j_{β} is the diffusion flux of scalar β across the cell face area A_x where the subscripts e and w refer to the "east" and "west" faces of a given grid cell, respectively. S_{β} is the Lagrangian source term derived from the conservation law for β , ρ represents mass density, and V represents cell volume. In practice, we refer to the left hand term on the right side of Equation 1 as the "mixing term" and the right hand term on the right side of Equation 1 as the "source term". The generic transport equation differs slightly in the spatial formulation of ODT, but its form is similar, so we omit it here for brevity. The system of ordinary differential equations (ODEs) that results is well behaved at all grid points and in all geometries in their finite-volume forms. For details on transport equation derivation and use in both the temporal and spatial formulations of ODT, see [30].

Eddy events occur as a Poisson process in accordance with their eddy rates, where a given eddy event of size l and location x_0 has an eddy timescale t and an associated eddy rate 1/t. Three user-defined ODT parameters control the eddy event process: the eddy rate parameter C scales the rate of occurrence of the eddies; the viscous penalty parameter Z suppresses small eddies; and the large eddy suppression parameter β_{LES} constrains eddies to those with timescales proportional to the elapsed simulation time. Sampled eddies that do not fit the defined parameters are rejected and not applied to the domain.

Eddy events modify domain variables using "triplet maps", as illustrated in Figure 1. In an eddy region of size l, transported property profiles are modified (mapped) by making three copies, compressing each spatially by a

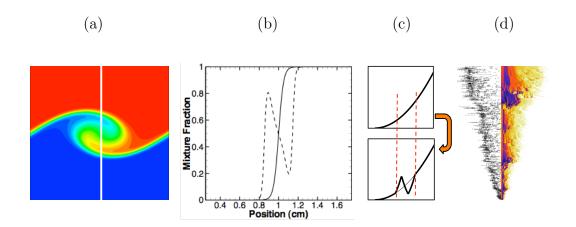


Figure 1: Schematic diagram of a triplet map. From left to right: (a) simulation of a single "eddy", a two-dimensional Kelvin-Helmholtz instability with a vertical white line-of-sight; (b) mixture fraction profile along the indicated line-of-sight before and after the eddy showing increased gradients and surface area; (c) a triplet map operates on a profile to produce a mapped profile with three copies, compressed and lined up, with the middle copy spatially inverted; (d) illustrative eddy event sizes and locations with corresponding temperature profile for a single ODT realization of a wall fire, adapted from [15].

factor of three, and realigning them with the middle copy spatially reversed. This map is continuous, conserves all quantities and statistical moments, increases scalar gradients, and decreases length scales consistent with turbulent eddies in real flows. Subsequent eddies in the same region will result in a cascade of scales. The eddy rates, used to sample and select eddy events, depend on eddy size and the local kinetic energy consistent with turbulent scaling laws.

Eddy events occur concurrently with diffusive advancement via solution of the system of unsteady one-dimensional transport equations. In this way, the ODT code marches in time or space until it reaches the simulation end point. Due to the stochastic nature of eddy events, each ODT simulation, or realization, is different, even when it is provided with the same input parameters. In order to obtain statistically stable data for a given set of parameters, we run many realizations with the same input parameters and then gather statistics over these realizations. This is done via post-processing tools, which are provided in the ODT package.

2.2. Software Architecture

The ODT package consists primarily of an object-oriented C++ code responsible for running flow simulation cases and generating data. The package also contains auxiliary data processing and visualization tools, written mostly

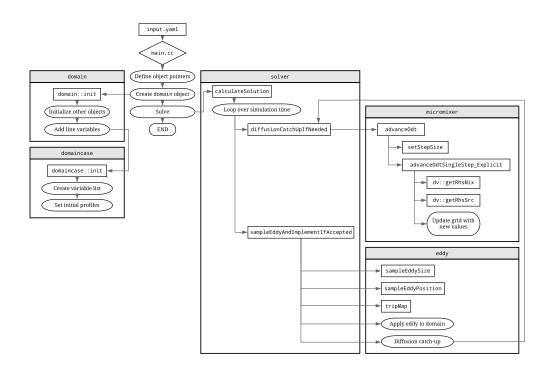


Figure 2: Structural outline of the ODT code, including its major class objects. For brevity, this diagram makes two simplifications. First, it assumes that diffusive advancement in the micromixer uses an explicit solution method; in practice, users specify either an explicit, semi-implicit, or Strang split method. Second, it implies that all of the eddy events sampled by the eddy object are applied to the domain; in reality, eddy events are filtered through several rejection tests (omitted from this figure for clarity) before acceptance and application to the domain.

in Python. The post-processing tools are case-specific but will be addressed generally in Section 2.3.

Figure 2 illustrates the ODT code's most important objects and structural features. User inputs are provided to the executable in YAML [36] format via input.yaml; the location of the specific input.yaml file to be used is determined by the case name and case type specified in the run script. The main function defines storage for the main objects, but, once created, the domain object is responsible for object initialization as well as variable initialization and storage via a case-specific domaincase object.

Three primary objects handle the code's main functions: the solver, micromixer, and eddy objects. The solver coordinates the ODT solution process, marching along the simulation time and invoking diffusive advancement and eddy events. The micromixer handles diffusive advancement by

setting step sizes, interacting with the transported domain variables, and solving the system of ODEs defined by Equation 1 (or its equivalent in the spatial formulation). The micromixer includes three solution methods that can be specified in input.yaml: a first-order explicit Euler method (pic-tured in Figure 2), a first-order semi-implicit method, and a second order Strang splitting method [37]. The latter two methods are useful for treating stiff chemistry, with source terms integrated using CVODE [38]. In reacting flow cases, enthalpy and chemical species mass fractions are transported, and physical properties and kinetic rates are computed with Cantera [39]. The micromixer is also the code's primary point of interaction with the mesher object (not pictured in Figure 2), which manages the adaptive grid functions. Finally, the eddy object manages eddy event selection and implementation, which proceeds as described in Section 2.1.

2.3. Workflow

This section outlines the process a user goes through in order to successfully build the ODT code and run a simulation. ODT is a standalone, self-contained package, and users interact with its files primarily via the command line rather than a graphical interface. For more details, please see the package documentation.

Within the main download package, several directories organize the ODT code. The source, build, and run directories contain the ODT source code, compilation tools, and run scripts, respectively. The input directory contains subdirectories corresponding to several possible case types, each populated with the appropriate input files. The data directory, initially empty, holds the raw data files and runtime information generated by ODT, as well as post-processed data files generated from within the post directory. Finally, the docs directory contains documentation optionally generated by Doxygen [40] during the build process.

2.3.1. Building ODT

The ODT build process is automated with CMake. The user navigates to the build directory and edits the CMake configuration file. The CMake configuration file specifies Cantera's location and must be changed to reflect the local installation location. For reacting flows, the chemical reaction rates can be computed by Cantera (default), or a from a user-specified function, specified at compile time. Once the configuration settings are updated, the user runs CMake to apply the changes.

YAML is required in order to process input files. If it has not yet been installed, it must be built and installed at this point. For convenience, the ODT package automates this process. YAML need only be installed once

for a given instance of the ODT package. Rebuilding the ODT code with CMake does not affect the YAML installation.

Once CMake and YAML have been prepared, the user can build the ODT code with the make command, which places the code executable in the run directory. The most common errors that occur during the build process involve incorrect file paths or incomplete installation of required packages. See the build documentation for troubleshooting help.

Once the code is built, the user may optionally build a local copy of the documentation via Doxygen, which must be previously installed. This step is not required in order to run the ODT code. As an alternative, documentation can be accessed at the code repository wiki.

2.3.2. Input files

User-modified input files are located within the input directory, which contains subdirectories that correspond to various case types that can be run with ODT. At minimum, a case's subdirectory must contain an input.yaml file, but may contain other files, or subfolders with supplementary information

The input directory also contains the gas_mechanisms subdirectory, which contains chemical mechanism files that can be used in reacting flow cases. For cases in which a chemical mechanism is not required, the not_used.xml mechanism file is specified in input.yaml. Note that the chemical mechanism chosen for the case and specified in the input file must match the mechanism specified in the CMake configuration file during the ODT build process.

Input files contain simulation parameters in a human-readable format parsed by YAML. Within input.yaml, parameters are organized into sections, several of which are common to all input files. Parameter values present in the provided input files may be considered defaults that can be used to run a successful simulation of that case type. General code parameters are stored in the params class. Variables needed in a simulation but not provided in an input file (not recommended), are either given defaults specified in the params class or result in an error message. Details about individual parameters, including usage and typical values, are covered in the documentation.

2.3.3. Running ODT

To run a simulation, users must then navigate to the run directory, which contains the odt.x executable and several possible run scripts. The simplest option is runOneRlz.sh, which runs one realization of ODT in the specified configuration. In this run script, the user must alter two variables near the top of the file: inputDir, which specifies which input directory and files to

use; and caseName, which provides a name for the simulation and the data files it outputs. The runManyRlz.sh script runs many realizations in serial, one after the other; it differs from runOneRlz.sh only in that the user must also alter the nRlz variable, which specifies the number of realizations to run. Each realization has a different seed value, which can be randomized or directly specified in the input file. To run the simulation with either runOneRlz.sh or runManyRlz.sh, save the run script and execute it at the command line. Users will see some output on the command line, but no data, which is instead output to the data directory.

ODT simulations can also be run in parallel using MPI. Two run scripts, slrmJob.sh and slrmJob_array.sh, are configured using Slurm [41], a common workload manager used for massively parallel computing resources. This allows many ODT realizations to run in parallel rather than in serial, reducing overall simulation time. Individual realizations are independent (embarrassingly parallel), but users must take care with case names and input file changes to ensure that individual realizations or entire cases are not overridden accidentally.

2.3.4. Data files and post-processing

For a given simulation, data is output to the data directory, which contains subdirectories for each simulation, specified by the caseName variable in the run script. Figure 3 illustrates the structure of the data directory and the locations of files within it. Each case folder is subdivided into input, runtime, data, and post. The input folder contains a copy of the input files used for the simulation; runtime contains runtime output, including eddy event information; data contains the raw data files; and post contains post-processed data files once they are generated.

To use post-processing tools, users navigate to the post directory. Within the post directory, data processing tools are organized by case type, which is specified in input.yaml and determines case-specific variables and simulation setup parameters (refer to the domaincase object in Figure 2). Each set of post-processing tools is different, but may contain some combination of Python scripts (often coordinated by a driver.py file), experimental data files for comparison and plot generation, or other supplementary files. Post-processed data and generated plots are deposited in the data directory, within the appropriate caseName/post folder.

3. Example Case: Canonical Jet Flame

ODT is uniquely suited to reacting flow simulations. Here, we present illustrative ODT simulation results of a round, turbulent jet flame based on

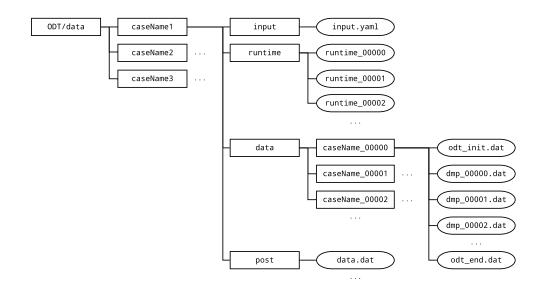


Figure 3: Data folder structure and file locations. Boxes indicate folders and ovals indicate files. For a given simulation, caseName will be replaced with the case name specified in the run script. The number attached to caseName folder names and to runtime data files indicates the realization number. The number attached to dmp files indicates the output time step, or "dump" time, of the data file. These correspond to the order of the time steps listed in input.yaml.

and compared to the experimental DLR-A flame of Meier et al. [42]. This canonical flame configuration has been used extensively to study and validate turbulent combustion models [43–48].

The DLR-A fuel stream is mixture of 22.1% CH_4 , 33.2% H_2 , and 44.7% N_2 (by volume) that issues into dry air via a nozzle with an inner diameter of 8 mm at a mean exit velocity of 42.2 m·s⁻¹. The coflow air stream issues from a concentric nozzle 140 mm in diameter at a velocity of 0.3 m·s⁻¹. The reported jet Reynolds number is 15,200.

Previous ODT studies of turbulent jet flames have used the temporal planar formulation, but the spatial cylindrical formulation developed recently [30] more closely matches the experimental configuration. This simulation uses the experimentally reported velocity profiles and jet dimensions. The fuel was diluted with N₂ in the experimental flame to minimize radiative heat losses, and radiation is ignored in the simulation. This flame has a low Reynolds number, and the combustion chemistry proceeds quickly. The ODT simulation transports the chemical species O₂, N₂, CH₄, H₂, H₂O, and CO₂. We assume that reactions proceed to the products of complete combustion and apply simple, fast reaction rates according to the following chemical equations:

$$CH_4 + 2O_2 \to CO_2 + 2H_2O,$$
 (2)

$$H_2 + \frac{1}{2}O_2 \to H_2O.$$
 (3)

These assumptions are reasonable for illustrating ODT in the DLR-A reacting jet configuration with variable properties and heat release.

This simulation uses ODT parameters C=20, $\beta_{LES}=17$, and Z=400. The values of C and β_{LES} were adjusted to give good agreement with the experimental data, and the value of Z is the same as the spatial simulations in [15]. 1024 independent flow realizations were performed in parallel and the results ensemble averaged. Downstream distance y and radial position r are normalized by the jet diameter D.

Figure 4 displays the simulation results: axial mean and centerline values for (a) axial velocity v, (b) mixture fraction ξ , and (c) temperature T. The ODT results track the experimental data well for all three variables. The centerline temperature peaks about 100 K above the experimental data, but this small difference can be attributed to thermal radiation (which was neglected in this simulation) and the assumption that reactions proceed to the products of complete combustion rather than their equilibrium state. The centerline velocity shows a fast initial decrease that can be attributed to diffusion. Similarly, the increase in centerline velocity RMS values is delayed; this occurs due to the elapsed time model for large eddy suppression, which

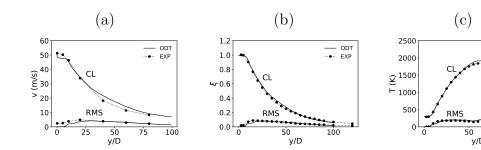


Figure 4: DLR jet flame example case results: (a) mean axial velocity and RMS velocity along the centerline versus downstream location; (b) mean axial mixture fraction and RMS mixture fraction along the centerline versus downstream location; (c) mean axial temperature (K) and RMS temperature (K) along the centerline versus downstream location.

limits disturbances in the early stages of the flow to small eddies that occur on the jet edges away from the centerline.

To replicate this example case, build the code with the CHEMISTRY = SIMPLEDLR flag in the user_config file and edit the desired run script with inputDir = "../input/jetFlame/DLR_A" and a new case name such as caseName = "jetFlame_example". The input file for this case, located at ODT/input/jetFlame/DLR_A/input.yaml, contains the appropriate parameters and does not need to be modified to match this example case. See the code documentation for data post-processing instructions.

4. Impact

The ODT model is complex, challenging to implement, and difficult to cleanly separate from its implementation, all of which presents barriers to practitioners who wish to use the model. For instance, triplet mapa can be represented mathematically but are usually implemented as previously described, assuming an adaptive grid with variable cell sizes. Alternative formulations are used for fixed grids. The code presented here reduces these barriers by providing an open-source, well-documented, modularly styled code that can be used directly or modified to suit particular research goals. Its structure facilitates addressing new research questions, such as soot-flame-radiation-chemistry interactions, multi-phase reacting flows using Lagrangian particle models (such as diesel sprays), and coupled near-wall heat and momentum transfer. The code presented here has been extended to and demonstrated in such systems already, facilitated by the object-oriented structure. For example, this code has been extended to solve steady laminar flamelets [?], requiring inheritance from the solver, micromixer, and dv (domain

variable) classes.

The structure of the ODT code, using an adaptive mesh and a Lagrangian formulation for diffusive advancement of transport equations, has several advantages, including simplicity of formulation and computational efficiency. The adaptive mesh minimizes the number of computational cells required and can be adapted to any variable of interest. This is especially important for flows with complex and expensive chemistry and flows with widely varying resolution demands such as near-wall flows. The Lagrangian formulation implies that cell faces move with fluid velocity (such as is present during flame expansion, for example). This treatment is inherently stable, removing the need for complex and diffusive hyperbolic conservation law treatments and their associated CFL constraints. The conservation equations, as in Equation 1, are similarly simplified.

The code has been run on Linux, Mac, and Windows machines, including parallel clusters of up to 512 processor cores. Several research groups have used the code and contributed to its development. These include groups at Sandia National Laboratories, Brandenburg University of Technology Cottbus-Senftenberg, Chalmers University, and Combustion Science and Engineering, Inc. The latter company is the only organization known to use the code in a commercial setting, though on a combustion research project.

5. Conclusion

A clean and extensible implementation of the one-dimensional turbulence (ODT) model has been presented in an open-source, object-oriented C++ framework, motivated by advanced turbulent reacting flow research. The build process, directory structure, input file format, run process, and post-processing procedure were described in detail, including a representative reacting jet example case. Access to a proven, documented code and a modern build procedure will aid researchers in applying and extending the model to new research questions.

329 6. Conflict of Interest

There are no known conflicts of interest associated with this publication and there has been no significant financial support for this work that could have influenced its outcome.

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