

Resource Summary Report

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CVODE

RRID:SCR_014696

Type: Tool

Proper Citation

CVODE (RRID:SCR_014696)

Resource Information

URL: <http://computation.llnl.gov/projects/sundials/cvode>

Proper Citation: CVODE (RRID:SCR_014696)

Description: Solver for stiff and nonstiff ordinary differential equation (ODE) systems (initial value problem), given in explicit form $y' = f(t,y)$, based on Adams and BDF methods. It is one of the programs that makes up the Suite of Nonlinear and Differential/Algebraic Equation Solvers (SUNDIALS)., THIS RESOURCE IS NO LONGER IN SERVICE. Documented on September 16,2025.

Synonyms: SUNDIALS CVODE

Resource Type: data processing software, software application, data analysis software, software resource

Keywords: data analysis software, solver, equation solver, ordinary differential equation, differential equation, stiff, nonstiff

Availability: THIS RESOURCE IS NO LONGER IN SERVICE

Resource Name: CVODE

Resource ID: SCR_014696

Record Creation Time: 20220129T080321+0000

Record Last Update: 20260815T182510+0000

Ratings and Alerts

No rating or validation information has been found for CVODE.

No alerts have been found for CVODE.

Data and Source Information

Source: [SciCrunch Registry](#)

Usage and Citation Metrics

We found 56 mentions in open access literature.

Listed below are recent publications. The full list is available at [NIF](#) .

Owen MJ, et al. (2025) IonBench: A benchmark of optimisation strategies for mathematical models of ion channel currents. PLoS computational biology, 21(8), e1013319.

Babaev D, et al. (2025) Mathematical modeling of pharmacokinetics and pharmacodynamics of losartan in relation to CYP2C9 allele variants. Frontiers in systems biology, 5, 1504077.

Kutumova EO, et al. (2024) Physiologically based pharmacokinetic model for predicting the biodistribution of albumin nanoparticles after induction and recovery from acute lung injury. Heliyon, 10(10), e30962.

Mate-Kole EM, et al. (2024) Mathematical complexities in radionuclide metabolic modelling: a review of ordinary differential equation kinetics solvers in biokinetic modelling. Journal of radiological protection : official journal of the Society for Radiological Protection, 44(2).

Yip HYK, et al. (2024) Integrative modeling uncovers p21-driven drug resistance and prioritizes therapies for PIK3CA-mutant breast cancer. NPJ precision oncology, 8(1), 20.

Mahardika T NQ, et al. (2024) Evaluation of cardiac pro-arrhythmic risks using the artificial neural network with ToR-ORD in silico model output. Frontiers in physiology, 15, 1374355.

Birgiolas J, et al. (2023) NeuroML-DB: Sharing and characterizing data-driven neuroscience models described in NeuroML. PLoS computational biology, 19(3), e1010941.

Miller C, et al. (2022) Multiscale modelling of desquamation in the interfollicular epidermis. PLoS computational biology, 18(8), e1010368.

Ataullakhanov FI, et al. (2022) Significance of two transmembrane ion gradients for human erythrocyte volume stabilization. PloS one, 17(12), e0272675.

Whittaker DG, et al. (2022) Ion channel model reduction using manifold boundaries. Journal of the Royal Society, Interface, 19(193), 20220193.

Britton OJ, et al. (2022) A population of in silico models identifies the interplay between Nav 1.8 conductance and potassium currents as key in regulating human dorsal root ganglion neuron excitability. F1000Research, 11, 104.

Marques CST, et al. (2021) Reduced Reaction Mechanisms for Ethanol under Ultra-lean Conditions in Internal Combustion Engines. ACS omega, 6(1), 206.

Yang B, et al. (2021) Reverse engineering gene regulatory network based on complex-valued ordinary differential equation model. BMC bioinformatics, 22(Suppl 3), 448.

Hendrix M, et al. (2021) cellmlmanip and chaste_codegen: automatic CellML to C++ code generation with fixes for singularities and automatically generated Jacobians. Wellcome open research, 6, 261.

Morii Y, et al. (2020) Optimization of one-parameter family of integration formulae for solving stiff chemical-kinetic ODEs. Scientific reports, 10(1), 21330.

Kopa? D, et al. (2020) First-Principles-Based Multiscale Modelling of Nonoxidative Butane Dehydrogenation on Cr₂O₃(0001). ACS catalysis, 10(24), 14732.

Tötsch N, et al. (2020) Bayesian Data Integration Questions Classic Study on Protease Self-Digest Kinetics. ACS omega, 5(25), 15162.

Gupta S, et al. (2020) Parallel Tempering with Lasso for model reduction in systems biology. PLoS computational biology, 16(3), e1007669.

Salazar-Cavazos E, et al. (2020) Multisite EGFR phosphorylation is regulated by adaptor protein abundances and dimer lifetimes. Molecular biology of the cell, 31(7), 695.

Heldt FS, et al. (2020) A Single Light-Responsive Sizer Can Control Multiple-Fission Cycles in Chlamydomonas. Current biology : CB, 30(4), 634.