

Virtual Reactor

VR[™] – III-V Edition

Software for Modeling of Epitaxial Growth of Group-III Arsenides and Phosphides by Metal-Organic Chemical Vapor Deposition

Version 7.6

Code Description



STR, IP Holding, LLC., Richmond, VA, USA Copyright © 2000-2017 by STR, IP Holding, LLC. All rights reserved. Published 2017.

This manual is the confidential and proprietary product of STR, IP Holding, LLC. Any unauthorized use, reproduction, or disclosure of this manual is strictly prohibited. (Subject to limited use within the STR End-User License Agreement only.)

CGSim[™], VR[™], PolySim[™], CVDSim[™], HEpiGaNS[™], SimuLED[™], SiLENSe[™], RATRO[™], SpeCLED[™], SimuLAMP[™], SELES[™], FETIS[™] are registered trademarks, brands, and proprietary products of STR, IP Holding, LLC.

User Support: <u>vr-support@str-soft.com</u> Software Sales: <u>STR-sales@str-soft.com</u> Phone: +7 812 320 4390 Fax: +7 812 326 6194 STR Group, Inc. <u>www.str-soft.com</u> Engels av. 27, P.O. Box 89, 194156, St. Petersburg, Russia



Software Overview

Virtual Reactor (VR[™]) is a family of two-dimensional software tools designed for the simulation of growth of bulk crystals and thin films from the vapor phase. VR[™] III-V Edition</sup> is designed for modeling of epitaxy of thin films of Group-III arsenides and phosphides.

Virtual Reactor provides comprehensive information about numerous physical processes responsible for crystal growth and film epitaxy. This includes information on the distribution of temperature and gas species concentration, heat fluxes, deposition rate and other parameters in the overall reactor and along all walls of the reactor units. This, in turn, provides wide possibilities for profound investigations of the phenomena underlying the growth, which allows optimization of the reactor geometry and growth conditions.

The problem is considered in axisymmetric or plane 2D approximation. The growth simulation includes modeling of the heat transfer, gas mixture flow and reactive species mass transport, in particular, multi-component diffusion and chemical reactions in the gas domain. The software employs a homogeneous chemistry model involving precursor decomposition and original heterogeneous chemistry models.

The Virtual Reactor is designed with a friendly user interface (GUI) which is aimed at minimization of the user efforts needed for the problem specification. The user needs to prescribe reactor geometry and specify the growth conditions.

Unstructured triangular and quadrilateral non-matched computational grids are used in numerical simulation. The grid generation is carried out for each geometry block providing the required grid density.

Series computations of the heat- and mass transport at a set of computation stages with user-specified values of the precursor flow rates, reference temperature, coil position and total gas pressure allow parametric study of the epitaxial growth dependence on the operating conditions.

Run-time and post-processing visualization is available within the GUI, presenting the two-dimensional and onedimensional distributions of temperature and other variables. In addition, Virtual Reactor is supplied with external visualization tool View 2D. The computational results are also stored in plain text files allowing a post-processing analysis using Tecplot and other commercial graphical packages.



Key Features

VR™ III-V Edition software tool includes the following models:

- > AIAs epitaxy from TMAI and AsH₃ with H₂ as carrier
- > GaAs epitaxy from TMGa and AsH₃ with H₂ as carrier
- > GaAs epitaxy from TMGa and TBAs with H₂ as carrier
- > InAs epitaxy from TMIn and AsH₃ with H₂ as carrier
- > InAs epitaxy from TMIn and TBAs with H_2 as carrier
- > $AI_xGa_{1-x}As$ epitaxy from TMAI, TMGa, and AsH₃ with H₂ as carrier
- In_xGa_{1-x}As epitaxy from TMIn, TMGa, and AsH₃ with H₂ as carrier
- > In_xGa_{1-x}As epitaxy from TMIn, TMGa, and TBAs with H₂ as carrier
- > AIP epitaxy from TMAI and PH_3 with H_2 as carrier
- > GaP epitaxy from TMGa and PH₃ with H₂ as carrier
- > GaP epitaxy from TMGa and TBP with H₂ as carrier
- > InP epitaxy from TMIn and PH₃ with H₂ as carrier
- > InP epitaxy from TMIn and TBP with H₂ as carrier
- > $In_xGa_{1-x}P$ epitaxy from TMIn, TMGa, and PH₃ with H₂ as carrier
- In_xGa_{1-x}P epitaxy from TMIn, TMGa, and TBP with H₂ as carrier
- > Al_xIn_yGa_{1-x-y}P epitaxy from TMAI, TMIn, TMGa, and PH₃ with H₂ as carrier

The mass transport and chemistry models incorporated into VR[™] III-V Edition can be applied to the following types of MOVPE reactors:

- > Planetary reactors
- High-speed rotation Turbodisc reactors
- Horizontal reactors
- Close-coupled showerhead reactors
- > Home-made reactors that can be simulated within a 2D plane or axisymmetric model

The following options are incorporated in the VR[™] III-V Edition software tool:

- > Modeling of the global heat transfer problem in a MOCVD reactor, including
 - Inductive heating (for axisymmetric problems only). The computation of the Joule heat sources due to inductive heating is carried out by solving the Maxwell equations.



- Resistive heating. The resistive heating is modeled by assigning a uniform heat release distribution in the heater. The user can split the heater block into several units and specify either the global heat release in the heater or the power density in each unit.
- Temperature fitting. The program predicts power needed to maintain the desired temperature(s) in a reference point(s).
- Conductive heat transfer in solid materials. The thermal conductivity of the materials used in the growth system can be prescribed by the user as a function of temperature. Anisotropic thermal conductivity can be assigned.
- Convective and radiative heat transfer in transparent gas blocks. The view-factor technique is used to model the radiation heat exchange. The surface emissivities are assigned as constant values.
- Radiative heat transfer in semitransparent solid blocks. The two-waveband radiation model is used to account for radiative heat transfer in semitransparent solid materials, which are characterized by threshold wavelength.
- > Modeling of the mass transport in the reactor includes
 - Non-isothermal flow of gas mixture. Heat conductivity and viscosity of the mixture are calculated in terms of molecular kinetic theory.
 - ✓ Multi-component diffusion of reactive species.
 - ✓ Homogeneous chemistry.
 - Instantaneous values of the growth rate and compound composition and the averaged values calculated accounting for the substrate rotation.
- > Turbulent flow considered using an algebraic turbulence model
- > Database of Material Properties, User Manual, Context Help, and files with examples of computations.
- Modeling of surface chemistry on chemically reactive surfaces of the substrate and parasitic deposition on the reactor side walls accounting for the kinetic limitations, which are calculated by the program as a function of temperature and local gas mixture composition.
- > Visualization of the obtained results.
 - ✓ Substrate rotation is accounted for by averaging the instantaneous distributions of the growth rate.
 - Calculation of the growth rate non-uniformity accounting for the edge exclusion.

More Info

More detailed information about the VR[™] III-V Edition can be requested by the e-mail address <u>VR-support@str-soft.com</u>.



Support

Hot-line support is provided on request. The support includes supply of updated versions released during the license period and technical consulting on the VR[™] III-V Edition</sup> operation.