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## **Preface**

Open Command Environment for Analysis (OCEAN) lets you set up, simulate, and analyze circuit data without starting Virtuoso Analog Design Environment L, XL or GXL.

This manual describes OCEAN and the commands required to set up, simulate, and analyze circuit data using OCEAN. This manual assumes that you are familiar with analog design and simulation using the Virtuoso Analog Design Environment. You should also be proficient in Cadence<sup>®</sup> SKILL language programming.

The preface contains the following:

- Scope
- <u>Licensing Requirements</u>
- Related Documentation
- Customer Support
- Feedback about Documentation
- Typographic and Syntax Conventions
- Identifiers Used to Denote Data Types

## Scope

Unless otherwise noted, the functionality described in this guide can be used in both mature node (for example, IC6.1.8) and advanced node (for example, ICADVM20.1) releases.

Label	Meaning
(ICADVM20.1 Only)	Features supported only in ICADVM20.1 advanced nodes and advanced methodologies releases.
(IC6.1.8 Only)	Features supported only in mature node releases.

## **Licensing Requirements**

You must have the Analog\_Design\_Environment\_L licence to use OCEAN. For information on licensing, see <u>Virtuoso Software Licensing and Configuration Guide</u>.

## **Related Documentation**

#### What's New and KPNS

- <u>Virtuoso Analog Design Environment XL What's New</u>
- Virtuoso Analog Design Environment XL Known Problems and Solutions

### Installation, Environment, and Infrastructure

- Cadence Installation Guide.
- <u>Virtuoso Design Environment User Guide</u>.
- Cadence SKILL Language User Guide
- Cadence SKILL Language Reference
- Cadence SKILL++ Object System Reference
- Virtuoso Design Environment SKILL Reference
- Virtuoso Design Environment SKILL Reference

#### Preface

- <u>Virtuoso Analog Design Environment L SKILL Language Reference</u>
- <u>Virtuoso Analog Design Environment XL SKILL Language Reference</u>

#### Virtuoso Tools

- <u>Virtuoso Analog Design Environment L User Guide</u>
- <u>Virtuoso Analog Design Environment XL User Guide</u>
- <u>Virtuoso Analog Design Environment GXL User Guide</u>
- <u>Virtuoso Analog Distributed Processing Option User Guide</u>

## **Additional Learning Resources**

### **Video Library**

The <u>Video Library</u> on the Cadence Online Support website provides a comprehensive list of videos on various Cadence products.

To view a list of videos related to a specific product, you can use the *Filter Results* feature available in the pane on the left. For example, click the *Virtuoso Layout Suite* product link to view a list of videos available for the product.

You can also save your product preferences in the Product Selection form, which opens when you click the *Edit* icon located next to *My Products*.

### **Virtuoso Videos Book**

You can access certain videos directly from Cadence Help. To learn more about this feature and to access the list of available videos, see Virtuoso Videos.

## **Rapid Adoption Kits**

Cadence provides a number of <u>Rapid Adoption Kits</u> that demonstrate how to use Virtuoso applications in your design flows. These kits contain design databases and instructions on how to run the design flow.

In addition, Cadence offers the following training courses the SKILL programming language, which you can use to customize, extend, and automate your design environment:

#### Preface

- SKILL Language Programming Introduction
- SKILL Language Programming
- Advanced SKILL Language Programming.

To explore the full range of training courses provided by Cadence in your region, visit Cadence Training or write to training\_enroll@cadence.com.

**Note:** The links in this section open in a separate web browser window when clicked in Cadence Help.

## **Help and Support Facilities**

Virtuoso offers several built-in features to let you access help and support directly from the software.

- The Virtuoso *Help* menu provides consistent help system access across Virtuoso tools and applications. The standard Virtuoso *Help* menu lets you access the most useful help and support resources from the Cadence support and corporate websites directly from the CIW or any Virtuoso application.
- The Virtuoso Welcome Page is a self-help launch pad offering access to a host of useful knowledge resources, including quick links to content available within the Virtuoso installation as well as to other popular online content.

The Welcome Page is displayed by default when you open Cadence Help in standalone mode from a Virtuoso installation. You can also access it at any time by selecting *Help – Virtuoso Documentation Library* from any application window, or by clicking the *Home* button on the Cadence Help toolbar (provided you have not set a custom home page).

For more information, see <u>Getting Help</u> in *Virtuoso Design Environment User Guide*.

## **Customer Support**

For assistance with Cadence products:

- Contact Cadence Customer Support
  - Cadence is committed to keeping your design teams productive by providing answers to technical questions and to any queries about the latest software updates and training needs. For more information, visit <a href="https://www.cadence.com/support">https://www.cadence.com/support</a>.
- Log on to Cadence Online Support

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Customers with a maintenance contract with Cadence can obtain the latest information about various tools at <a href="https://support.cadence.com">https://support.cadence.com</a>.

## **Feedback about Documentation**

You can contact Cadence Customer Support to open a service request if you:

- Find erroneous information in a product manual
- Cannot find in a product manual the information you are looking for
- Face an issue while accessing documentation by using Cadence Help

You can also submit feedback by using the following methods:

- In the Cadence Help window, click the *Feedback* button and follow instructions.
- On the Cadence Online Support <u>Product Manuals</u> page, select the required product and submit your feedback by using the <u>Provide Feedback</u> box.

## **Understanding Cadence SKILL**

Cadence SKILL is a high-level, interactive programming language based on the popular artificial intelligence language, Lisp. It lets you customize and extend your design environment. Using SKILL you can validate the steps of your algorithm incrementally before incorporating them into a larger program.

For more information about the SKILL language, see <u>Getting Started</u> in the SKILL Language User Guide.

### **Using SKILL Code Examples**

The SKILL APIs in this user manual are explained with illustrative code examples.

You can copy these examples from the manual and paste them directly into the Command Interpreter Window (CIW) or use the code in non-graphical SKILL mode.

## Sample SKILL Code

The following code sample shows the syntax of a SKILL API that accepts three arguments.

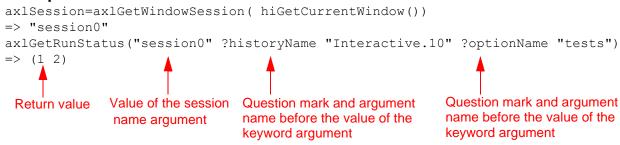
#### axIGetRunStatus

The first argument  $t\_sessionName$  is a required argument, where t signifies the data type of the argument. The second and third arguments ?optionName  $t\_optionName$  and ?historyName  $t\_historyName$  are optional keyword arguments (identified by a question mark), which are specified in name-value pairs and can be placed in any order during the function call.

#### Preface

The return value is the value that the SKILL API returns after evaluating the expression. In this case, it is a list of status values, 1 statusValues.

#### **Example**



### **Accessing API Help**

Quick reference information for SKILL APIs is available from the CIW and the SKILL API Finder. To access the reference information for a particular SKILL API, do one of the following:

- Type help <function name> in the CIW.
- Type startFinder ([?funcName t\_functionName]) in the CIW.
- Start the <u>SKILL API Finder</u> from the CIW by choosing *Tools Finder* or type cdsFinder on the UNIX command line.

In the Search in field of the displayed Cadence SKILL API Finder window, type the SKILL API name for which you want to display the help information and click Go.

The matches for the searched SKILL API appear in the Results area.

To view the complete documentation of the searched SKILL API, select the API name in the Results area and click the More Info button. The complete documentation of the selected SKILL API appears in a new Cadence Help window.

## **Typographic and Syntax Conventions**

The following typographic and syntax conventions are used in this manual.

text	Indicates names of manuals, menu commands, buttons, and fields.
text	Indicates text that you must type exactly as presented. Typically used to denote command, function, routine, or argument names that must be typed literally.
$z\_argument$	Indicates text that you must replace with an appropriate argument value. The prefix (in this example, $z_{-}$ ) indicates the data type the argument can accept and must not be typed.
	Separates a choice of options.
{ }	Encloses a list of choices, separated by vertical bars, from which you <b>must</b> choose one.
[ ]	Encloses an optional argument or a list of choices separated by vertical bars, from which you <b>may</b> choose one.
[ ?argName t_arg ]	
	Denotes a <i>key argument</i> . The question mark and argument name must be typed as they appear in the syntax and must be followed by the required value for that argument.
• • •	Indicates that you can repeat the previous argument.
	Used with brackets to indicate that you can specify zero or more arguments.
	Used without brackets to indicate that you must specify at least one argument.
,	Indicates that multiple arguments must be separated by commas.
=>	Indicates the values returned by a Cadence <sup>®</sup> SKILL <sup>®</sup> language function.
/	Separates the values that can be returned by a Cadence SKILL language function.

If a command-line or SKILL expression is too long to fit within the paragraph margins of this document, the remainder of the expression is moved to the next line and indented. In code excerpts, a backslash (\) indicates that the current line continues on to the next line.

## **Identifiers Used to Denote Data Types**

Data type identifiers are used to indicate the type of value required by an API argument. These data types are denoted by a single letter that is prefixed to the argument label and is separated from the argument by an underscore; for example, t is the data type in  $t\_viewName$ . Data types and underscores are used only as identifiers; they must not be typed when specifying the argument in a function.

Prefix	Internal Name	Data Type
а	array	array
А	amsobject	AMS object
b	ddUserType	DDPI object
В	ddCatUserType	DDPI category object
C	opfcontext	OPF context
d	dbobject	Cadence database object (CDBA)
е	envobj	environment
f	flonum	floating-point number
F	opffile	OPF file ID
g	general	any data type
G	gdmSpecIIUserType	generic design management (GDM) spec object
h	hdbobject	hierarchical database configuration object
I	dbgenobject	CDB generator object
K	mapiobject	MAPI object
1	list	linked list
L	tc	Technology file time stamp
m	nmpIIUserType	nmpll user type
Μ	cdsEvalObject	cdsEvalObject
n	number	integer or floating-point number
0	userType	user-defined type (other)
p	port	I/O port
q	gdmspecListIIUserType	gdm spec list

### Preface

Prefix	Internal Name	Data Type
r	defstruct	defstruct
R	rodObj	relative object design (ROD) object
S	symbol	symbol
${\mathcal S}$	stringSymbol	symbol or character string
t	string	character string (text)
T	txobject	transient object
и	function	function object, either the name of a function (symbol) or a lambda function body (list)
U	funobj	function object
V	hdbpath	hdbpath
W	wtype	window type
SW	swtype	subtype session window
dw	dwtype	subtype dockable window
X	integer	integer number
Y	binary	binary function
&	pointer	pointer type

For more information, see <u>Cadence SKILL Language User Guide</u>.

## Introduction to OCEAN

This chapter provides an introduction to Open Command Environment for Analysis (OCEAN). In this chapter, you can find information about

- Types of OCEAN Commands
- OCEAN Online Help
- OCEAN Syntax Overview
- Parametric Analysis
- Distributed Processing
- Plotting Simulation Results

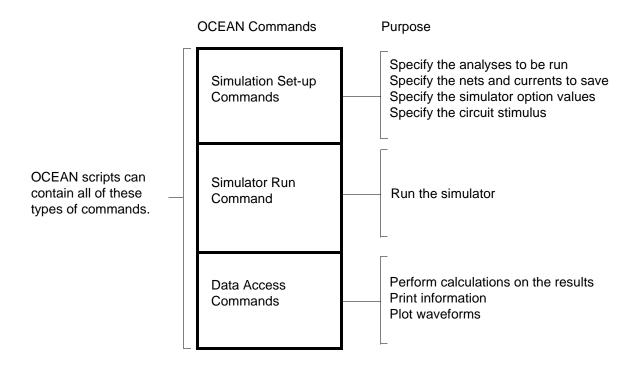
OCEAN lets you set up, simulate, and analyze circuit data. OCEAN is a text-based process that you can run from a UNIX shell or from the Command Interpreter Window (CIW). You can type OCEAN commands in an interactive session, or you can create scripts containing your commands, then load those scripts into OCEAN. OCEAN can be used with any simulator integrated into the Virtuoso® Analog Design Environment.

Typically, you use the Virtuoso® Analog Design Environment when creating your circuit (in Composer) and when interactively debugging the circuit. After the circuit has the performance you want, you can use OCEAN to run your scripts and test the circuit under a variety of conditions. After making changes to your circuit, you can rerun your scripts. OCEAN lets you

- Create scripts that you can run repeatedly to verify circuit performance
- Run longer analyses such as parametric analyses and statistical analyses more effectively
- Run long simulations in OCEAN without starting the Virtuoso® Analog Design Environment graphical user interface
- Run simulations from a nongraphic, remote terminal

## **Types of OCEAN Commands**

You can create OCEAN scripts to accomplish the full suite of simulation and data access tasks that you can perform in the Virtuoso® Analog Design Environment. An OCEAN script can contain three types of commands, as shown in the following figure:



All the parameter storage format (PSF) information created by the simulator is accessible through the OCEAN data access commands. (The data access commands include all of the Virtuoso® Analog Design Environment calculator functions.)

You can use the <u>history</u> command to view the command history from the current session and the most recently terminated session.

## **OCEAN Online Help**

Online help is available for all the OCEAN commands when you are in an OCEAN session. To get help for a specific OCEAN command, type the following:

```
ocnHelp( 'commandName )
```

This command returns an explanation of the command and examples of how the command can be used.

#### Introduction to OCEAN

To get a listing of all the different types of commands in OCEAN, type the following: ocnHelp()

For more information, see "ocnHelp" on page 186.

## **OCEAN Syntax Overview**

OCEAN is based on the Virtuoso<sup>®</sup> SKILL programming language and uses SKILL syntax. All the SKILL language commands can be used in OCEAN. This includes if statements, case statements, for loops, while loops, read commands, print commands, and so on.

The most commonly used SKILL commands are documented in this manual. However, you are not limited to these commands. You can use any SKILL routine from any SKILL manual.

## **Common SKILL Syntax Characters Used in OCEAN**

This section provides an overview of some basic SKILL syntax concepts that you need to understand to use OCEAN. For more information about SKILL syntax, see <u>Chapter 3</u>, <u>"Introduction to SKILL."</u>

#### **Parentheses**

Parentheses surround the arguments to the command. The command name is followed immediately by the left parenthesis, with no intervening space.

#### **Examples**

The following example shows parentheses correctly enclosing two arguments to the path command.

```
path( "~/simulation1/schematic/psf" "~/simulation2/schematic/psf" )
```

In the next example, the space after the command name causes a syntax error.

```
Syntax error.

path ( "~/simulation1/schematic/psf" "~/simulation2/schematic/psf" )
```

### Introduction to OCEAN

#### **Quotation Marks**

Quotation Marks are used to surround string values. A string value is a sequence of characters, such as "abc".

In the following example, the directory names provided to the path command are strings, which must be surrounded by quotation marks.

```
path( "~/simulation1/schematic/psf" "~/simulation2/schematic/psf" )
```

#### Convention

In this manual, a SKILL convention is used to let you know when an argument must be a string. When you see the prefix  $t_{-}$ , you must substitute a string value (surrounded by quotation marks) for the argument. Consider the following syntax statement:

```
desVar( t desVar1 g value1 t desVar2 g value2)
```

In this case, there are two string values that must be supplied:  $t_{desVar1}$  and  $t_{desVar2}$ . (The  $g_{prefix}$  indicates a different type of argument. For more information about prefixes, see Chapter 4, "Working with SKILL.")

#### **Recovering from an Omitted Quotation Mark**

Accidentally omitting a closing quotation mark from an OCEAN command can cause great confusion. For example, typing the incorrect command

```
strcat( "rain" "bow )
```

appears to hang OCEAN. In an attempt to recover, you type a Control-c. That gives you a prompt but it does not fix the problem, as you discover when you then type the correct command.

```
strcat( "rain" "bow" )
```

Again, you have to type a Control-c and OCEAN responds with another message.

```
^C*Error* parser: interrupted while reading input
```

If you find yourself in this situation, do not press a Control-c. Instead, recover by entering a quotation mark followed by a right square bracket ( ] ). This procedure reestablishes a normal OCEAN environment and you can then reenter the correct command.

```
ocean> strcat( "rain" "bow )
"]
"rainbow ) "
ocean> strcat( "rain" "bow" )
"rainbow"
ocean>
```

#### Introduction to OCEAN

### Single Quotation Marks

The single quotation mark indicates that an item is a symbol. Symbols in SKILL correspond to constant enums in C. In the context of OCEAN, there are predefined symbols. The simulator that you use also has predefined symbols. When using symbols in OCEAN, you must use these predefined symbols.

### **Examples**

In the following example, tran is a symbol and must be preceded by a single quotation mark. The symbol tran is predefined. You can determine what the valid symbols for a command are by checking the valid values for the command's arguments. For example, if you refer to "analysis" on page 91, you see that the valid values for the first argument include 'tran.

```
analysis ( 'tran ... )
```

The list of items you can save with the save command is also predefined. You must choose from this predefined list. See "save" on page 155 and refer to the valid values for the  $s\_saveType$  argument. The 'v symbol indicates that the item to be saved is the voltage on a net.

```
save( 'v "net1" )
```

#### Convention

In this manual, a SKILL convention is used to let you know when an argument must be a symbol. When you see the prefix  $s_{-}$ , you must substitute a symbol (preceded by a single quotation mark) for the argument. Consider the following syntax statement:

```
selectResults( s resultsName ) => t / nil
```

In this case, there is one symbol that must be supplied:  $s\_resultsName$ . For the selectResults command, there is a different mechanism that lets you know the list of predefined symbols. If you type the following command, with no arguments, the list of predefined symbols is returned: results() => ( dc tran ac )

**Note:** Depending on which results are selected, the values returned by the results command vary.

#### **Question Mark**

The question mark indicates an optional keyword argument, which is the first part of a keyword parameter. A keyword parameter has two components:

■ The first component is the keyword, which has a question mark in front of it.

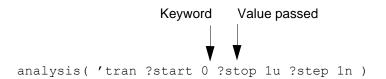
#### Introduction to OCEAN

■ The second component is the value being passed, which immediately follows the keyword.

Keyword parameters, composed of these keyword/value pairs, are always optional.

### **Examples**

In the following example, all the arguments to the analysis command except 'tran are keyword/value pairs and are optional.



For example, you can use ?center and ?span instead of ?start and ?stop. You also can omit ?start altogether because it is an optional argument.

#### Convention

In this manual, a SKILL convention is used to let you know when arguments are optional. Optional arguments are surrounded by square brackets []. In the following example, all of the keyword/value pairs are surrounded by square brackets, indicating that they are optional.

## **Data Types Used in OCEAN**

The following table shows the internal names and prefixes for the SKILL data types that are used in OCEAN commands.

Data Type	Internal Name	Prefix	
floating-point number	flonum	f	
any data type	general	g	
linked list	list	1	
integer, floating-point number, or complex number		n	
user-defined type		0	

### **OCEAN Reference** Introduction to OCEAN

Data Type	Internal Name	Prefix
I/O port	port	р
symbol	symbol	S
symbol or character string		S
character string (text)	string	t
window type		W
integer number	fixnum	X

For more information about SKILL datatypes, see Chapter 4, "Working with SKILL."

#### **OCEAN Return Values**

You get return values from most OCEAN commands and can use these values in other OCEAN commands.

The following table shows some examples in which the return value from a command is assigned to a variable.

Assigning a Return Value to a Variable	Resulting Value for the Variable
a=desVar("r1" 1k)	a=1k
a=desVar("r1" 1k "r2" 2k)	a=2k
a=desVar("r1")	a=1k, assuming r1 was set in a desVar command
a=desVar("r2")	a=2k, assuming r2 was set in a desVar command

## **Design Variables in OCEAN**

Design variables in OCEAN function as they do in the Virtuoso® Analog Design Environment. Design variables are not assigned in the order specified. Rather, they are reordered and then assigned. Consider the following example:

```
desVar( "a" "b+1" )
desVar( "b" 1 )
```

#### Introduction to OCEAN

You might expect an error because a is assigned the value b+1 before b is assigned a value. However, OCEAN reorders the statements and sends them as follows:

```
desVar( "b" 1 )
desVar( "a" "b+1" )
```

After the reordering, there is no error. (b is equal to 1 and a is equal to 2.)

Suppose you run a simulation, then specify the following:

```
desVar("b" 2)
```

You might expect a to be equal to 2, which was the last value specified. Instead, a is reevaluated to b+1 or 3.

This approach is similar to how the design variables are used in simulation. For example, consider a circuit that has the following resistor:

```
R1 1 0 resistor r=b
```

If you change the value of b, you expect the value of R1 to change. You do not expect to have to netlist again or retype the R1 instantiation.

This approach is used in the Virtuoso® Analog Design Environment. Users are not expected to enter design variables in a particular order. Rather, the design variables are gathered during the design variable search then reordered before they are used.

You can also use a conditional expression containing the ternary operator (?:) to specify the value of a design variable, as shown below:

```
desVar( "b" 1 )
desVar( "a" "((b>1)?1:0)" )
```

**Note:** Do not use simulator reserved words as design variable names. For more information, see the <u>Reserved Words</u> section in the *Virtuoso Analog Design Environment User Guide*.

## outputs() in OCEAN

Throughout this manual are examples of nets and instances preceded by a "/" as well as examples without the "/". There is a significant difference between the two.

If you create a design in the Virtuoso® Analog Design Environment and save the OCEAN file, all net and instance names will be preceded with a "/", indicating they are schematic names. The netlist/amap directory must be available to map these schematic names to names the simulator understands. (If your design command points to the raw netlist in the netlist directory, the amap directory is there.)

#### Introduction to OCEAN

If you create a design or an OCEAN script by hand, or move the raw netlist from the netlist directory, the net and instance names might not be preceded with "/". This indicates that simulator names are used, and mapping is not necessary.

If you are unsure whether schematic names or simulator names are used, after  $selectResult(S\_resultsName)$ , type outputs() to see the list of net and instance names.

**Note:** Although you can move the raw netlist file from the netlist directory, it is not advised. There are other files in the netlist directory that are now required to run OCEAN.

# **Parametric Analysis**

There are two ways you can run parametric analyses in OCEAN:

- You can use the paramAnalysis command (recommended approach).
- You can use a SKILL for loop.

Using the paramAnalysis command is an easier approach. With this command, you can set up any number of nested parametric analyses in an OCEAN script. The paramRun command runs all the parametric analyses. When the analysis is complete, the data can be plotted as a family of curves. The following example shows how you might use nested parametric analyses:

```
paramAnalysis( "rl" ?start 200 ?stop 600 ?step 200
    paramAnalysis( 'rs ?start 300 ?stop 700 ?step 200
    )
)
paramRun ()
```

In this example, the outer loop cycles through r1, and the inner loop cycles through rs as follows:

Loop through r1 from 200 to 600 by 200.

Loop through rs from 300 to 700 by 200.

Run.

End the first loop.

End the second loop.

So, for r1=200, rs equals 300, 500 and then 700. Then, for r1=400, rs equals 300, 500, and then 700. Finally, for r1=600, rs equals 300, 500, and then 700

#### Introduction to OCEAN

Use a SKILL for loop only if the paramAnalysis command is not adequate. You can use the SKILL for loop to set up any number of variable-switching runs. After all the simulations are complete, you have to work with the results directories individually. The following example shows how you might use SKILL loops for parametric analyses.

```
Cload = list( 2p 4p 6p 8p )
foreach( val Cload
   desVar( "Cload" val )
   a=resultsDir( sprintf( nil "./demo/Cload=%g" val ) )
   printf( "%L", a )
   run( )
)

foreach( val Cload
   openResults( sprintf( nil "./<dir>/Cload=%g" val ) )
   selectResults( 'ac )
   plot( vdb( "vout" ) )
}
```

### **Data Access Without Running a Simulation**

You can retrieve and use data from previous simulations at any time by opening the data with the openResults command. After opening the data, you can use any data access commands on this data. For more information, see <a href="#">Chapter 7</a>, "Data Access Commands."

You can use query commands such as results, outputs, and dataTypes to see what data is available to be opened.

# **Distributed Processing**

You can use OCEAN distributed processing commands to run simulations across a collection of computer systems. The distributed processing commands allow you to specify where and when jobs are run and allow you to monitor and control jobs in a variety of ways. Using distributed commands, you can

- Submit one or more jobs to a distributed processing queue
- Specify a host or group of hosts on which to distribute jobs
- View the status of jobs
- Specify when a job will run or in what sequence a group of jobs will run
- Suspend and resume jobs
- Cancel jobs

#### Introduction to OCEAN

For you to be able to use the distributed processing commands, your site administrator needs to set up the lists of machines to which jobs are submitted. Each list of machines is a group of hosts and is called a queue. Consult the <u>Virtuoso Analog Distributed Processing Option User Guide</u> for more information on how to configure systems for distributed processing. For information on the distributed processing commands for OCEAN, see <u>Chapter 12</u>, "OCEAN Distributed Processing Commands."

### **Blocking and Nonblocking Modes**

You can configure jobs to run in blocking or nonblocking mode. In blocking mode, execution of subsequent OCEAN commands is halted until the job completes. In nonblocking mode, the system does not wait for the first job to finish before starting subsequent jobs.

#### **Blocking Mode**

You must run jobs in blocking mode to be able to use the data resulting from a job in a subsequent command in an OCEAN script or batch run.

For example, if you want to run a simulation, select the tran results from that simulation, and then plot them, you

- 1. Configure the simulation with setup commands
- 2. Run the simulation with the run () command
- 3. Select the desired results with the selectResults ( 'tran) command
- 4. Plot the results with the plot () command

A job like the one in the example above must run in blocking mode so that the commands are processed sequentially. If the jobs in the example above are run in nonblocking mode, the selectResult command starts before the run command can return any data, and the selectResult command and the plot command cannot complete successfully.

#### **Nonblocking Mode**

If you are submitting several jobs that have no interdependencies, you can run them concurrently when hostmode is set to distributed.

For example, if you want to run two separate simulations as jobs, but do not want to wait until the first is complete before starting the second, you

1. Configure the first simulation with setup commands

#### OCEAN Reference Introduction to OCEAN

2. Configure a second simulation with setup commands

In the example above, the script starts the first job and then starts the second job without waiting for the first job to finish.

If you are running several commands, and some of them are data access commands, you can use the <u>wait</u> command to block a single job. The <u>wait</u> command is needed between the simulation and the data access commands to ensure the desired simulation is complete before the data is accessed.

For example, if you want to run two separate simulations as jobs (sim1 and sim2), and want to select and plot the results of the second simulation run, you

- 1. Configure the first simulation with setup commands
- 2. Run the simulation with a run(?jobPrefix "sim1") command
- **3.** Configure a second simulation with setup commands
- **4.** Run the second simulation with the run ( ?jobPrefix "sim2) command
- **5.** Cause the script to wait until the second simulation finishes before starting the selectResults command with the wait (sim2) command
- **6.** Select the desired results with the selectResults ('tran) command
- **7.** Plot the results with the plot ( ) command

In the example above, the script starts the first job and then starts the second job without waiting for the first job to finish. When the script reaches the wait command, it pauses until the second simulation runs and then selects the results to plot.

# **Plotting Simulation Results**

The simulation results can be plotted in Virtuoso Visualization and Analysis XL, which is supported in the OCEAN environment.

# **Using OCEAN**

This chapter explains the different ways you can use OCEAN to perform simulation tasks. In this chapter, you can find information about

- OCEAN Use Models
- Using OCEAN Interactively
- License Requirements
- Creating OCEAN Scripts
- Selecting Results
- Running Multiple Simulators
- OCEAN Tips

## **OCEAN Use Models**

There are two ways you can use OCEAN:

- You can use OCEAN interactively to perform simple tasks.
- You can use OCEAN in batch mode and provide the name of an existing (or parameterized) script as a command line argument. OCEAN scripts can be created:
  - □ From the Virtuoso® Analog Design Environment window with the command Session Save Script
  - Using a text editor

For information about creating scripts, see "Creating OCEAN Scripts" on page 46.

All the OCEAN commands are described in this manual, and online help is available for all these commands. For information about using the OCEAN online help, see <u>"OCEAN Online Help"</u> on page 30.

Using OCEAN

**Note:** The current version of OCEAN has some specific issues that are addressed in <u>Appendix 15, "OCEAN 4.4.6 Issues."</u> Refer this appendix before using OCEAN.

# **Using OCEAN Interactively**

You can run OCEAN from a UNIX prompt or from the Virtuoso<sup>®</sup> design framework II (DFII) Command Interpreter Window (CIW).

**Note:** The primary use model is to use OCEAN in a UNIX shell. Unless otherwise indicated, the rest of this chapter assumes that you are working from OCEAN in a UNIX shell.

### **Using OCEAN from a UNIX Shell**

To start OCEAN from a UNIX prompt, type the following command:

ocean

This command loads and reads the .oceanrc file. You can place OCEAN commands in your .oceanrc file, which is similar to the .cdsinit file. This file can contain any valid OCEAN command, function or SKILL initialization routine (excluding graphical dflI references, such as bindkeys and so on, which are not applicable to OCEAN). If you do not want to specify any startup initialization options for OCEAN, you do not need to create or add an .oceanrc file.

The OCEAN prompt appears indicating that you have started OCEAN:

ocean>

If you do not see this prompt after starting OCEAN, press Return. If you still do not see this prompt, you may have redefined the prompt with the setPrompt command. (This does not affect OCEAN; the prompt just will not indicate OCEAN is running.)

Now you can start typing OCEAN commands interactively. For an example of interactive use, see <u>"Interactive Session Demonstrating the OCEAN Use Model"</u> on page 45.

To guit the OCEAN executable from UNIX, type the following command:

exit

#### **OCEAN in Non-Graphical Mode**

OCEAN is an executable shell script that calls the AWD workbench and passes all its command-line options to it using the following shell command:

#### **Using OCEAN**

```
#! /bin/sh -
exec awd -ocean "$@"
```

This makes OCEAN highly dependent on the UNIX shell environment.

You can run OCEAN in a non-graphical mode by using the <code>-nograph</code> option with the <code>ocean</code> command. This disables the graphical options of the software. This option is useful if OCEAN is started on a machine that does not have X-Windows running.

**Note:** You can use the -nograph option to run the OCEAN job through a cron. Ensure that DISPLAY is set to ": 0". If the screen will be locked when the OCEAN cron job runs, use the allowaccess option with the xlock command on the UNIX prompt. For more information on the usage of xlock, type man xlock in a terminal window.

The <code>-nograph</code> option must only be used to replay logfiles that have been created interactively. For example, while using OCEAN with the <code>-nograph</code> option, your <code>oceanScript.ocn</code> file must have an <code>exit()</code> statement at the end followed by a newline. Otherwise, OCEAN hangs. The reason for this is that when the workbench is started in the non-graphical mode, it does not redirect standard I/O as it normally does; instead, it lets the SKILL human interface (HI) handle the standard I/O. HI expects an <code>explicitexit()</code> statement at the end of the OCEAN script and OCEAN exits only when it detects an <code>exit()</code> at EOF. The command is used as follows:

```
ocean -nograph < oceanScript.ocn > oceanScript.log
```

Alternatively, you can execute the OCEAN script using the -replay option. The command is used as follows:

```
ocean -nograph -replay oceanScript.ocn -log oceanScript.log
```

.oceanrc is automatically loaded while using ocean -nograph -replay command. If you use the virtuoso command, .oceanrc is not loaded automatically.

While using the <code>-nograph</code> option with <code>ocean</code>, if you find that simulation run messages are not being stored in the log file, check for the following environment variable in the <code>.cdsenv</code> file:

```
(envGetVal "spectre.envOpts" "firstRun" )
```

It must be set to nil as shown below for simulation run messages to be stored in it:

```
(envSetVal "spectre.envOpts" "firstRun" 'boolean nil)
```

For more information about this variable, see Appendix B of the *Virtuoso Analog Design Environment L User Guide*.

**Using OCEAN** 

### **Using OCEAN from the CIW**

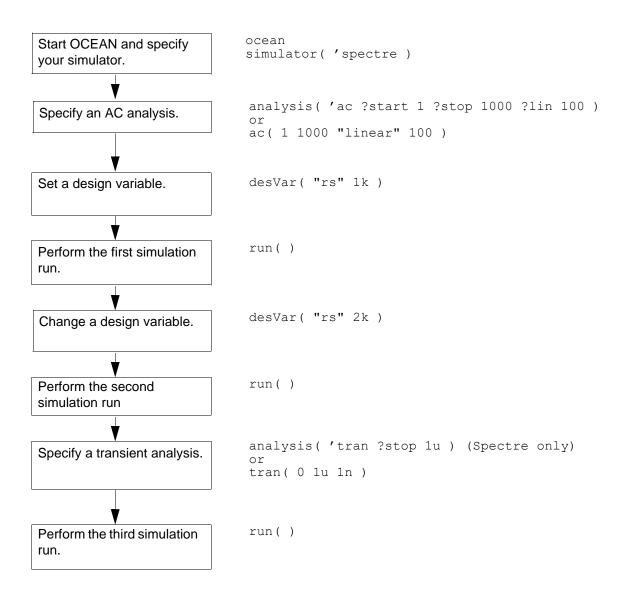
You can type OCEAN commands in the CIW after you bring up the Virtuoso® Analog Design Environment. (Starting the design environment loads the required OCEAN files.)

Your .oceanrc file is *not* automatically read when you start the DFII software (using the virtuoso command). Therefore, you might want to load your .oceanrc file manually in the CIW if you need information that it contains.

You can also use the <u>history</u> command from the CIW to list and reuse the most recently used commands.

### Interactive Session Demonstrating the OCEAN Use Model

The following figure shows a typical set of simulation tasks that you might perform interactively in OCEAN with the corresponding commands.



On the second and third run, the AC analysis runs because it is still active. If you do not want it to run, you must disable it with the following command:

```
delete( 'analysis 'ac )
```

The simulator is not called and run until the run () command is entered.

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The commands can be given in any order, as long as they are all defined before the run() command.

# **License Requirements**

You need licenses to run the simulator() and ocnSetXLMode() OCEAN commands. For more information on these commands, see <u>simulator</u>.

- To run the simulator() OCEAN command, you must have one of the following licenses. If one of these licenses are not already checked out, the first available license will be checked out in the following order:
  - 95200 Virtuoso(R) Analog Design Environment L
  - 95210 Virtuoso(R) Analog Design Environment XL
  - 95220 Virtuoso(R) Analog Design Environment GXL
- To run the ocnSetXLMode() OCEAN command, you must have one of the following licenses. If one of these licenses are not already checked out, the first available license will be checked out in the following order:
  - 95210 Virtuoso(R) Analog Design Environment XL
  - 95220 Virtuoso(R) Analog Design Environment GXL

**Note:** If you have run the <code>ocnSetXLMode()</code> command, running the <code>simulator()</code> command subsequently will not checkout an additional license.

If you do not want OCEAN to automatically checkout a higher tiered license—for example, if you do not want OCEAN to automatically checkout the 95210 license if the 95200 license is not available—set the following environment variable in your .cdsenv file:

```
asimenv.misc alwaysTryHigherTieredLicenseInOcean 'boolean nil
```

**Note:** If the alwaysTryHigherTieredLicenseInOcean environment variable is set to nil, errors are displayed if OCEAN is unable to checkout a license.

**Note:** The 95200 Virtuoso(R) Analog Design Environment L license is checked out when you load the ocean script. Exit Virtuoso, or run the <a href="https://ocentrology.org/ncmman4">ocnCloseSession()</a> command to release the license.

# **Creating OCEAN Scripts**

You can modify the included sample script files or create script files interactively from the Virtuoso® Analog Design Environment.

Using OCEAN

### **Creating Scripts Using Sample Script Files**

You can create your own script files with a text editor using the sample scripts as examples, or you can make copies of the sample scripts and modify them as needed using a text editor. The scripts can be found in the following directory:

your install dir/tools/dfII/samples/artist/OCEAN

Refer to the README file in this directory for information about the scripts.

### **Creating Scripts from the Analog Design Environment**

When you perform tasks in the design environment, the associated OCEAN commands are automatically stored in the <code>simulatorx.ocn</code> file in your <code>netlist</code> directory. For example, if you start the Virtuoso software, open the Virtuoso® Analog Design Environment window, and run a simulation using the Cadence SPICE simulator, a <code>cdsSpice0.ocn</code> file is created in your <code>netlist</code> directory. You can load this <code>cdsSpice0.ocn</code> script as described in "Loading OCEAN Scripts" on page 50.

### **Selectively Creating Scripts**

You can be selective about the information that is created in your .ocn script. The Virtuoso® Analog Design Environment has a feature that lets you create an OCEAN script based on the state of your current session. The following example illustrates how using this feature is different than using the automatic script generation feature.

Consider the following task flow:

- 1. Start the Virtuoso® Analog Design Environment.
- 2. Specify a DC analysis.
- 3. Select nets on the schematic to save.
- **4.** Run the simulation.
- **5.** Turn off the DC analysis.
- **6.** Select a transient analysis.
- 7. Run the simulation.
- **8.** Save the script from the Virtuoso® Analog Design Environment.

#### Using OCEAN

The script that is created, called oceanScript.ocn by default, contains only the selected nets, the transient analysis, and the run command. The script does not contain the DC analysis because it was turned off.

In contrast, the <code>simulator0.ocn</code> script, which is automatically created in the <code>netlist</code> directory, contains all of the commands, including the DC analysis and the current state of the analysis (on or off).

#### **Creating a Script**

To selectively create a script from Virtuoso Analog Design Environment L or,

**1.** Start the Virtuoso software.

virtuoso

The CIW appears.

**2.** From the CIW, choose *Tools – ADE L – Simulation*.

The Virtuoso Analog Design Environment window appears.

- **3.** Perform all of the design environment tasks that you want to capture in the script.
- **4.** Choose Session Save Script.

The Save Ocean Script to File form appears.

**5.** Click OK to accept the default file name (./oceanScript.ocn), or change the name for the file and click OK.

A script containing the OCEAN commands for the tasks you performed is created. For information about how to load this script, see "Loading OCEAN Scripts" on page 50.

#### **Controlling What Is Included in Scripts**

You can use .cdsenv variables to alter the OCEAN script that is created when you choose Session – Create Script in the Virtuoso Analog Design Environment. One variable allows you to include default environment settings in a script, two other variables allow you to run procedures before and after a script is created.

#### Including Default Control Statements

To save every control statement, including default statements, in your OCEAN script, add the following line to your .cdsenv file.

asimenv.misc saveDefaultsToOCEAN boolean t

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Setting saveDefaultsToOCEAN to t results in a complete dump of the current circuit design environment, defaults and all. Because the created OCEAN script contains all the settings, you might use this variable when you plan to archive a script, for example.

If saveDefaultsToOCEAN is not set to t, the created OCEAN script contains only those items that you explicitly set to some value other than their default.

#### Running Functions Before or After Creating a Script

The information in this section describes how you can specify functions to be run before or after a script is created. You can use these functions, for example, to add information at the beginning or end of a script. To use this capability follow these steps.

- 1. Decide when you want the functions to run.
  - Add the following line to your .cdsenv file to run the function preoceanFunc before the OCEAN script is created.

```
asimenv.misc preSaveOceanScript string "preOceanFunc"
```

Add the following line to your .cdsenv file to run the function postOceanFunc after the OCEAN script is created.

```
asimenv.misc postSaveOceanScript string "postOceanFunc"
```

**2.** Use the following syntax to specify the functions.

```
preOceanFunc( session fp )
postOceanFunc( session fp )
```

In this syntax, session is the OASIS session and fp is the file pointer to the OCEAN script file. For guidance on determining the session to use, see the <u>VirtuosoAnalog</u> <u>Design Environment L SKILL Language Reference</u>.

3. Load the functions in your .cdsinit file.

For example, you might add the following lines to your .cdsenv file.

```
asimenv.misc preSaveOceanScript string "MYfirstProc"
asimenv.misc postSaveOceanScript string "MYlastProc"
```

The functions MYfirstProc and MYlastProc might be defined like this.

```
procedure( MYfirstProc( session fp)
  fprintf(fp "; This will be the first line in the ocean script.\n")
)
procedure( MYlastProc( session fp)
  fprintf(fp "; This will be the last line in the ocean script.\n")
)
```

If these procedures are defined in a file called myOceanProcs.il, you can load them by adding to your .cdsinit file a command like the following.

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```
load "myOceanProcs.il"
```

When you choose Session – Create Script, first the preSaveOceanScript procedure is called, then the OCEAN script is created, then the postSaveOceanScript procedure is called.

### **Loading OCEAN Scripts**

You can load OCEAN scripts from OCEAN (in UNIX) or from the CIW.

#### From a UNIX Shell

To load an OCEAN script,

**1.** Type the following command to start OCEAN:

ocean

The OCEAN prompt appears.

2. Use the SKILL load command to load your script:

```
load( "script name.ocn" )
```

Messages about the progress of your script appear.

#### From the CIW

To load an OCEAN script,

1. Start the Virtuoso software

```
virtuoso &
```

The CIW appears.

2. In the text entry field, use the SKILL load command to load your script:

```
load( "script name.ocn" )
```

Messages about the progress of your script appear in the CIW.

# **Selecting Results**

You may use OCEAN to run several simulations on the same design and save the results in different result directories. You can then use Analog Design Environment XL to select the results and work with features like annotation etc.

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# Selecting Results Run from Worst Case Scripts for Cross-Probing or Back Annotating Operating Points

Assume that you have been using Ocean to create separate data directories for worst case parameter sweeps. Also, assume that the new directories you make are accessed with the resultsDir() ocean command in your Ocean script and that these directories are in the standard location where psf data is stored in the Analog Design Environment.

In the Analog Design Environment, psf data is stored in:

<runDir>/simulation/<testSchemName>/spectre/schematic/psf
where.

runDir is the directory where you invoke virtuoso&

testSchemName is your test schematic

This implies that your script should store the new directories under the schematic directory. Therefore, if c1, c2 and c3 are the worst case directories, they are located at:

<runDir>/simulation/<testSchemName>/spectre/schematic/c1
<runDir>/simulation/<testSchemName>/spectre/schematic/c2
<runDir>/simulation/<testSchemName>/spectre/schematic/c3

- 1. Choose Results Select
- 2. The Select Results form opens. Click Browse. A Unix Browser form appears.
- **3.** Navigate to the directory that contains your Ocean generated directories c1, c2, and c3.
- **4.** Click *OK* on the Unix Browser form. Now the *Select Results* Form should show c1, c2 and c3.
- **5.** Double-click c1, c2 or c3. Alternatively, you can also single-click c1, c2 or c3 and then choose *Update Results* and click *OK*. At this point the data is selected though there is no confirmation in the CIW. Now, you should be able to use *Results Direct Plot*, *Results Annotate* etc to see the results of that particular directory.

## **Selecting Results Run from Spectre Standalone**

After running spectre standalone, you can select results using the *Results Browser* and use calculator to plot the results. However, this does not allow you to use ADE features like *Results – Direct Plot* or *Results – Annotate*.

#### Using OCEAN

#### Consider that your data is in

<runDir>/simulation/<testSchemName>/spectre/schematic/psf.

where,

runDir is the directory where you invoke virtuoso&

testSchemName is your test schematic

- **1.** Choose *Tools Results Browser*. A pop up box appears. Enter your design path up to the spectre directory.
- **2.** Click *OK*, and the browser comes up.
- **3.** Click the schematic directory. The psf directory should appear.
- **4.** Click the directory with the data in it, psf. When you click the 'psf' directory you should see the tree expand with different results from your spectre stand alone simulation, e.g. tran.tran etc.
- **5.** Place the mouse pointer over the 'psf' node in the tree and press down the middle mouse key and scroll down to "create ROF". You should now see the psf directory change, and an intermediate node comes up --Run1-- betweenpsf/ and the results.
- **6.** Place the middle mouse pointer over the Run1 node, scroll down and select "Select Results".

**Note:** Even though there is a confirmation message in the CIW that the select was success, Analog Design Environment is not synced up to allow cross-probing and backannotation of operating points.

**7.** You may now use *Tools – Calculator* to select objects from the schematic. You can then choose 'plot' from the calculator, or different calculator operations.

**Note:** You CAN use *Tools – Calculator* but you CAN NOT use *Results – Direct Plot* or *Results – Annotate* etc.

# **Running Multiple Simulators**

There are times when you might want to run more than one simulator. You might be benchmarking simulators or comparing results. In OCEAN, you can only use one simulator per OCEAN session. If you change simulators, you must start a new OCEAN session. This is because some OCEAN command arguments are simulator specific, and therefore change when the simulator changes. For example, the arguments to the option command are

#### Using OCEAN

simulator specific. (No two simulators have the exact same options.) The analyses are typically simulator specific also.

# **OCEAN Tips**

The information in this section can help you solve problems that you encounter while using OCEAN.

■ While working in OCEAN, you might get the following SKILL error message:

```
*Error* eval: unbound variable - nameOfVariable
```

In this case, you need to see if you have an undeclared variable or if you are missing a single quotation mark (') or a quotation mark (") for one of your arguments. For example, the following command returns an error message stating that fromVal is an unbound variable because the variable has not been declared:

```
analysis('tran ?from fromVal)
```

However, the following pair of statements work correctly because from Val has a value (is bound).

```
fromVal=0
analysis('tran ?from fromVal)
```

If you get an error in an OCEAN session, you are automatically put into the SKILL debugger. In this case, you see a prompt similar to this:

```
ocean-Debug 2>
```

You can continue working. However, if you would like to get out of the debugger, you can type

```
debugQuit()
```

Now you are back to the normal prompt:

```
ocean>
```

- If it appears that OCEAN does not accept your input, or OCEAN appears to hang, then you may have forgotten to enter a closing quotation mark. Type "] to close all strings. For more information, and some examples, see <u>"Recovering from an Omitted Quotation Mark"</u> on page 32.
- In SKILL, the following formats are equivalent: (one two) and one (two). Results might be returned in either format. For example, OCEAN might return ac(tran) or (ac tran), but the two forms are equivalent.
- You can check your script for simple syntax errors by running SKILL lint. For example, you might use a command like

```
sklint -file myScript.ocn
```

**Using OCEAN** 

From within OCEAN, you can run SKILL lint by typing the following at the OCEAN prompt:

sklint(?file "yourOceanScript.ocn")

Running SKILL lint helps catch basic errors, such as unmatched parentheses and strings that are not closed.

# Introduction to SKILL

This chapter introduces you to the basic concepts that can help you get started with the Virtuoso<sup>®</sup> SKILL programming language. In this chapter, you can find information about

- The Advantages of SKILL
- Naming Conventions
- Arithmetic Operators
- Scaling Factors
- Relational and Logical Operators
- SKILL Syntax
- Arithmetic and Logical Expressions

# The Advantages of SKILL

The SKILL programming language lets you customize and extend your design environment. SKILL provides a safe, high-level programming environment that automatically handles many traditional system programming operations, such as memory management. SKILL programs can be immediately run in the Virtuoso environment.

SKILL is ideal for rapid prototyping. You can incrementally validate the steps of your algorithm before incorporating them in a larger program.

SKILL leverages your investment in Cadence technology because you can combine existing functionality and add new capabilities.

SKILL lets you access and control all the components of your tool environment: the User Interface Management System, the Design Database, and the commands of any integrated design tool. You can even loosely couple proprietary design tools as separate processes with SKILL's interprocess communication facilities.

# **Naming Conventions**

The recommended naming scheme is to use uppercase and lowercase characters to separate your code from code developed by Cadence.

All code developed by Cadence Design Systems typically names global variables and functions with up to three lowercase characters, that signify the code package, and the name starting with an uppercase character. For example, <code>dmiPurgeVersions()</code> or <code>hnlCellOutputs</code>. All code developed outside Cadence should name global variables by starting them with an uppercase character, such as <code>AcmeGlobalForm</code>.

# **Arithmetic Operators**

SKILL provides many arithmetic operators. Each operator corresponds to a SKILL function, as shown in the following table.

Sample SKILL Operators

Operators in Descending Precedence	Underlying Function
**	exponentiation
* /	multiply divide
+ -	plus minus
== !=	equal nequal
=	assignment

# **Scaling Factors**

SKILL provides a set of scaling factors that you can add to the end of a decimal number (integer or floating point) to achieve the scaling you want.

- Scaling factors must appear immediately after the numbers they affect. Spaces are not allowed between a number and its scaling factor.
- Only the first nonnumeric character that appears after a number is significant. Other characters following the scaling factor are ignored. For example, for the value 2.3mvolt, the m is significant, and the volt is discarded. In this case, volt is only for your reference.

### OCEAN Reference Introduction to SKILL

If the number being scaled is an integer, SKILL tries to keep it an integer; the scaling factor must be representable as an integer (that is, the scaling factor is an integral multiplier and the result does not exceed the maximum value that can be represented as an integer). Otherwise, a floating-point number is returned.

The scaling factors are listed in the following table. **Scaling Factors** 

Character	Name	Multiplier	Examples
Υ	Yotta	10 <sup>24</sup>	10Y [ 10e+25 ]
Z	Zetta	10 <sup>21</sup>	10Z [ 10e+22 ]
E	Exa	10 <sup>18</sup>	10E [ 10e+19 ]
Р	Peta	10 <sup>15</sup>	10P [ 10e+16 ]
Т	Tera	10 <sup>12</sup>	10T [ 1.0e13 ]
G	Giga	10 <sup>9</sup>	10G [ 10,000,000,000 ]
М	Mega	10 <sup>6</sup>	10M [ 10,000,000 ]
K	Kilo	10 <sup>3</sup>	10K [ 10,000 ]
%	percent	10 <sup>-2</sup>	5% [ 0.05 ]
m	milli	10 <sup>-3</sup>	5m [ 5.0e-3 ]
u	micro	10 <sup>-6</sup>	1.2u [ 1.2e-6 ]
n	nano	10 <sup>-9</sup>	1.2n [ 1.2e-9 ]
p	pico	10 <sup>-12</sup>	1.2p [ 1.2e-12 ]
f	femto	10 <sup>-15</sup>	1.2f [ 1.2e-15 ]
а	atto	10 <sup>-18</sup>	1.2a [ 1.2e-18 ]
Z	zepto	10 <sup>-21</sup>	1.2z [ 1.2e-21 ]
у	yocto	10 <sup>-24</sup>	1.2y [ 1.2e-24 ]

**Note:** The characters used for scaling factors depend on your target simulator. For example, if you are using cdsSpice, the scaling factor for M is different than shown in the previous table, because cdsSpice is not case sensitive. In cdsSpice, M and M are both interpreted as  $10^{-3}$ , so ME or ME is used to signify  $10^{6}$ .

# **Relational and Logical Operators**

This section introduces you to

■ Relational Operators: <, <=, >, >=, ==, !=

■ Logical Operators: !, &&, ||

### **Relational Operators**

Use the following operators to compare data values. SKILL generates an error if the data types are inappropriate. These operators all return t or nil. Sample Relational Operators

Operator	Arguments	Function	Example	Return Value
<	numeric	lessp	3 < 5 3 < 2	t nil
<=	numeric	leqp	3 <= 4	t
>	numeric	greaterp	5 > 3	t
>=	numeric	geqp	4 >=3	t
==	numeric string list	equal	3.0 == 3 "abc" == "ABc"	t nil
!=	numeric string list	nequal	"abc" != "ABc"	t

Knowing the function name is helpful because error messages mention the function (greaterp below) instead of the operator ( > ).

```
1 > "abc"
Message: *Error* greaterp: can't handle (1 > "abc")
```

### **Logical Operators**

SKILL considers nil as FALSE and any other value as TRUE. The and (&&) and or (||) operators only evaluate their second argument if it is required for determining the return result.

#### **Sample Logical Operators**

Operator	Arguments	Function	Example	Return Value
&&	general	and	3 && 5 5 && 3 t && nil nil && t	5 3 nil nil
	general	or	3    5 5    3 t    nil nil    t	3 5 t

The && and | | operators return the value last computed. Consequently, both && and | | operators can be used to avoid cumbersome if or when expressions.

The following example illustrates the difference between using && and | | operators and using if or when expressions.

#### You do not need to use

```
If (usingcolor then
currentcolor=getcolor()
else
currentcolor=nil
)
Instead use
```

currentcolor=usingcolor && getcolor()

#### **Using &&**

When SKILL creates a variable, it gives the variable a value of unbound to indicate that the variable has not been initialized yet. Use the boundp function to determine whether a variable is bound. The boundp function

- Returns t if the variable is bound to a value
- Returns nil if the variable is not bound to a value

#### Introduction to SKILL

Suppose you want to return the value of a variable trMessages. If trMessages is unbound, retrieving the value causes an error. Instead, use the expression

boundp( 'trMessages ) && trMessages

#### Using ||

Suppose you have a default name, such as noName, and a variable, such as userName. To use the default name if userName is nil, use the following expression:

userName || "noName"

# **SKILL Syntax**

This section describes SKILL syntax, which includes the use of special characters, comments, spaces, parentheses, and other notation.

### **Special Characters**

Certain characters are special in SKILL. These include the *infix* operators such as less than (<), colon (:), and assignment (=). The following table lists these special characters and their meaning in SKILL.

**Note:** All nonalphanumeric characters (other than \_ and ?) must be preceded (*escaped*) by a backslash (\) when you use them in the name of a symbol. **Special Characters in SKILL** 

Character	Name	Meaning
\	backslash	Escape for special characters
( )	parentheses	Grouping of list elements, function calls
[ ]	brackets	Array index, super right bracket
,	single quotation mark	Specifies a symbol (quoting the expression prevents its evaluation)
"	quotation mark	String delimiter
,	comma	Optional delimiter between list elements
;	semicolon	Line-style comment character

### Special Characters in SKILL

Character	Name	Meaning
+, -, *, /	arithmetic	For arithmetic operators; the /* and */ combinations are also used as comment delimiters
!,^,&,	logical	For logical operators
<,>,=	relational	For relational and assignment operators; < and > are also used in the specification of bit fields
?	question mark	If first character, implies keyword parameter
%	percent sign	Used as a scaling character for numbers

### **White Space**

White space sometimes takes on semantic significance and a few syntactic restrictions must therefore be observed.

Write function calls so the name of a function is immediately followed by a left parenthesis; no white space is allowed between the function name and the parenthesis. For example

```
f(a b c) and g() are legal function calls, but f (a b c) and g () are illegal.
```

The unary minus operator must immediately precede the expression it applies to. No white space is allowed between the operator and its operand. For example

```
-1, -a, and - (a*b) are legal constructs, but - 1, - a, and - (a*b) are illegal.
```

The binary minus (subtract) operator should either be surrounded by white space on both sides or be adjacent to non-white space on both sides. To avoid ambiguity, one or the other method should be used consistently. For example:

a - b and a-b are legal constructs for binary minus, but a -b is illegal.

#### Comments

SKILL permits two different styles of comments. One style is block oriented, where comments are delimited by /\* and \*/. For example:

```
/* This is a block of (C style) comments comment line 2 comment line 3 etc. \begin{tabular}{ll} \star/ \end{array}
```

#### Introduction to SKILL

The other style is line- oriented where the semicolon (;) indicates that the rest of the input line is a comment. For example:

```
\mathbf{x} = \mathbf{1} ; comment following a statement; comment line 1; comment line 2 and so forth
```

For simplicity, line-oriented comments are recommended. Block-oriented comments cannot be nested because the first \*/ encountered terminates the whole comment.

#### **Role of Parentheses**

Parentheses () delimit the names of functions from their argument lists and delimit nested expressions. In general, the innermost expression of a nested expression is evaluated first, returning a value used in turn to evaluate the expression enclosing it, and so on until the expression at the top level is evaluated. There is a subtle point about SKILL syntax that C programmers, in particular, must notice.

#### Parentheses in C

In C, the relational expression given to a conditional statement such as if, while, and switch must be enclosed by an outer set of parentheses for purely syntactical reasons, even if that expression consists of only a single Boolean variable. In C, an if statement might look like

```
if (done) i=0; else i=1;
```

#### Parentheses in SKILL

In SKILL, parentheses are used for specifying lists, calling functions, delimiting multiple expressions, and controlling the order of evaluation. You can write function calls in prefix notation

```
(fn2 arg1 arg2) or (fn0)
```

as well as in the more conventional algebraic form

```
fn2(arq1 arq2) or fn0()
```

The use of syntactically redundant parentheses causes variables, constants, or expressions to be interpreted as the names of functions that need to be further evaluated. Therefore,

Never enclose a constant or a variable in parentheses by itself; for example, (1), (x).

#### Introduction to SKILL

■ For arithmetic expressions involving *infix* operators, you can use as many parentheses as necessary to force a particular order of evaluation, but never put a pair of parentheses immediately outside another pair of parentheses; for example, ((a + b)): the expression delimited by the inner pair of parentheses would be interpreted as the name of a function.

For example, because if evaluates its first argument as a logical expression, a variable containing the logical condition to be tested should be written without any surrounding parentheses; the variable by itself is the logical expression. This is written in SKILL as

```
if ( done then i = 0 else i = 1)
```

#### **Line Continuation**

SKILL places no restrictions on how many characters can be placed on an input line, even though SKILL does impose an 8,191 character limit on the strings being entered. The parser reads as many lines as needed from the input until it has read in a complete form (that is, expression). If there are parentheses that have not yet been closed or binary *infix* operators whose right sides have not yet been given, the parser treats carriage returns (that is, newlines) just like spaces.

Because SKILL reads its input on a form-by-form basis, it is rarely necessary to "continue" an input line. There might be times, however, when you want to break up a long line for aesthetic reasons. In that case, you can tell the parser to ignore a carriage return in the input line, by preceding it immediately with a backslash (\).

```
string = "This is \
a test."
=> "This is a test."
```

# **Arithmetic and Logical Expressions**

Expressions are SKILL objects that also evaluate to SKILL objects. SKILL performs a computation as a sequence of function evaluations. A SKILL *program* is a sequence of expressions that perform a specified action when evaluated by the SKILL interpreter.

There are two types of primitive expressions in SKILL that pertain to OCEAN: constants and variables.

#### **Constants**

A *constant* is an expression that evaluates to itself. That is, evaluating a constant returns the constant itself. Examples of constants are 123, 10.5, and "abc".

#### Introduction to SKILL

#### **Variables**

A *variable* stores values used during the computation. The variable returns its value when evaluated. Examples of variables are a, x, and init var.

When the interpreter evaluates a variable whose value has not been initialized, it displays an error message telling you that you have an unbound variable. For example, you get an error message when you misspell a variable because the misspelling creates a new variable.

```
myVariable
```

causes an error message because it has been referenced before being assigned, whereas myVariable = 5

works.

When SKILL creates a variable, it gives the variable an initial value of unbound. It is an error to evaluate a variable with this value because the meaning of unbound is that-value-which-represents-no-value. unbound is not the same as nil.

### **Using Variables**

You do not need to declare variables in SKILL as you do in C. SKILL creates a variable the first time it encounters the variable in a session. Variable names can contain

- Alphanumeric characters
- Underscores ( \_ )
- Question marks
- Digits

The first character of a variable must be an alphanumeric character or an underscore. Use the assignment operator to store a value in a variable. You enter the variable name to retrieve its value.

### **Creating Arithmetic and Logical Expressions**

Constants, variables, and function calls can be combined with the *infix* operators, such as less than (<), colon (:), and greater than (>) to form arithmetic and logical expressions. For example: 1+2, a\*b+c, x>y.

### **OCEAN Reference** Introduction to SKILL

Υo	u can f	form	arbitrarily	complicated	expressions	by	combining	any	number	of the	primitive
ex	oressic	ns d	lescribed	above.							

Introduction to SKILL

# **Working with SKILL**

This chapter provides information on using SKILL functions. It includes information on the types of SKILL functions, the types of data accepted as arguments, how data types are used, and how to declare and define functions. In this chapter, you can find information about

- SKILL Functions
- Data Types
- Arrays
- Concatenating Strings (Lists)
- Declaring a SKILL Function
- SKILL Function Return Values
- Syntax Functions for Defining Functions

# **SKILL Functions**

There are two basic types of SKILL functions:

- Functions carry out statements and return data that can be redirected to other commands or functions.
- Commands are functions that carry out statements defined by the command and return t or nil. Some commands return the last argument entered, but the output cannot be redirected.

# **Data Types**

SKILL supports several data types, including integer and floating-point numbers, character strings, arrays, and a highly flexible linked list structure for representing aggregates of data. The simplest SKILL expression is a single piece of data, such as an integer, a floating-point

### Working with SKILL

number, or a string. SKILL data is case sensitive. You can enter data in many familiar ways, including the following:

### Sample SKILL Data Items

Data Type	Syntax Example
integer	5
floating point number	5.3
text string	"Mary had a little lamb"

For symbolic computation, SKILL has data types for dealing with symbols and functions.

For input/output, SKILL has a data type for representing I/O ports. The table below lists the data types supported by SKILL with their internal names and prefixes.

Data Types Supported by SKILL

Data Type	Internal Name	Prefix
array	array	а
boolean		b
floating-point number	flonum	f
any data type	general	g
linked list	list	I
floating-point number or integer n		
user-defined type		0
I/O port	port	p
symbol	symbol	S
symbol or character string		S
character string (text)	string	t
window type		W
integer number	fixnum	х

#### **Numbers**

SKILL supports the following numeric data types:

#### Working with SKILL

- Integers
- Floating-point

Both integers and floating-point numbers may use scaling factors to scale their values. For information on scaling factors, see <u>"Scaling Factors"</u> on page 56.

#### **Atoms**

An *atom* is any data object that is not a grouping or collection of other data objects. Built into SKILL are several special atoms that are fundamental to the language.

nil The nil atom represents both a false logical condition and an

empty list.

t The symbol t represents a true logical condition.

Both nil and t always evaluate to themselves and must never be used as the name of a variable.

unbound

To make sure you do not use the value of an uninitialized variable, SKILL sets the value of all symbols and array elements initially to unbound so that such an error can be detected.

#### **Constants and Variables**

Supported constants and variables are discussed in <u>"Arithmetic and Logical Expressions"</u> on page 3-14.

## Strings

Strings are sequences of characters; for example, "abc" or "123". A string is marked off by quotation marks, just as in the C language; the empty string is represented as "". The SKILL parser limits the length of input strings to a maximum of 8,191 characters. There is, however, no limit to the length of strings created during program execution. Strings of more than 8,191 characters can be created by applications and used in SKILL if they are not given as arguments to SKILL string manipulation functions.

#### Working with SKILL

When typing strings, you specify

- Printable characters (except a quotation mark) as part of a string without preceding them with the backslash (\) escape character
- Unprintable characters and the quotation mark itself by preceding them with the backslash (\) escape character, as in the C language

# **Arrays**

An *array* represents aggregate data objects in SKILL. Unlike simple data types, you must explicitly create arrays before using them so the necessary storage can be allocated. SKILL arrays allow efficient random indexing into a data structure using familiar syntax.

- Arrays are not typed. Elements of the same array can be different data types.
- SKILL provides run-time array bounds checking. The array bounds are checked with each array access during runtime. An error occurs if the index is outside the array bounds.
- Arrays are one dimensional. You can implement higher dimensional arrays using single dimensional arrays. You can create an array of arrays.

# Allocating an Array of a Given Size

Use the declare function to allocate an array of a given size.

- The declare function returns the reference to the array storage and stores it as the value of week.
- The type function returns the symbol array.

# **Concatenating Strings (Lists)**

#### Concatenating a List of Strings with Separation Characters (buildString)

buildString makes a single string from the list of strings. You specify the separation character in the third argument. A null string is permitted. If this argument is omitted, buildString provides a separating space as the default.

#### **Concatenating Two or More Input Strings (strcat)**

strcat creates a new string by concatenating two or more input strings. The input strings are left unchanged.

```
strcat( "l" "ab" "ef" ) => "labef"
```

You are responsible for any separating space.

```
strcat( "a" "b" "c" "d" ) => "abcd" strcat( "a" "b" "c" "d" ) => "a b c d"
```

### Appending a Maximum Number of Characters from Two Input Strings (strncat)

strncat is similar to strcat except that the third argument indicates the maximum number of characters from string2 to append to string1 to create a new string. string1 and string2 are left unchanged.

### **Comparing Strings**

### Comparing Two Strings or Symbol Names Alphabetically (alphalessp)

alphalessp compares two objects, which must be either a string or a symbol, and returns t if arg1 is alphabetically less than arg2. alphalessp can be used with the sort function to sort a list of strings alphabetically. For example:

```
stringList = '( "xyz" "abc" "ghi" )
sort( stringList 'alphalessp ) => ("abc" "ghi" "xyz")
```

The next example returns a sorted list of all the files in the login directory:

#### Working with SKILL

```
sort( getDirFiles( "~" ) 'alphalessp )
```

#### **Comparing Two Strings Alphabetically (strcmp)**

strcmp compares two strings. (To test if two strings are equal or not, you can use the equal command.) The return values for strcmp are explained in the following table.

Return Value	Meaning
1	string1 is alphabetically greater than string2.
0	string1 is alphabetically equal to string2.
-1	string1 is alphabetically less than string2.

```
strcmp( "abc" "abb" )=> 1
strcmp( "abc" "abc")=> 0
strcmp( "abc" "abd")=> -1
```

# Comparing Two String or Symbol Names Alphanumerically or Numerically (alphaNumCmp)

alphaNumCmp compares two string or symbol names. If the third optional argument is not nil and the first two arguments are strings holding purely numeric values, a numeric comparison is performed on the numeric representation of the strings. The return values are explained in the following table.

Return Value	Meaning
1	arg1 is alphanumerically greater than arg2.
0	arg1 is alphanumerically identical to arg2.
-1	arg2 is alphanumerically greater than arg1.

# **Declaring a SKILL Function**

To refer to a group of statements by name, use the procedure declaration to associate a name with the group. The group of statements and the name make up a SKILL function.

- The name is known as the function name.
- The group of statements is the function body.

## Working with SKILL

To run the group of statements, mention the function name followed immediately by ().

The clearplot command below erases the Waveform window and then plots a net.

```
procedure( clearplot( netname )
    clearAll( )
    plot( v (netName))
)
```

# **Defining Function Parameters**

To make your function more versatile, you can identify certain variables in the function body as formal parameters.

When you start your function, you supply a parameter value for each formal parameter.

# **Defining Local Variables (let)**

Local variables can be used to establish temporary values in a function. This is done using the let statement. When local variables are defined, they are known only within the let statement and are not available outside the let statement.

When the function is defined, the let statement includes the local variables you want to define followed by one or more SKILL expressions. The variables are initialized to nil. When the function runs, it returns the last expression computed within its body. For example:

```
procedure( test ( x )
        let(( a b )
        a=1
        b=2
        x * a+b
        )
)
```

- The function name is test.
- The local variables are a and b.
- The local variables are initialized to nil.
- The return value is the value of x \* a + b.

## OCEAN Reference Working with SKILL

# **SKILL Function Return Values**

All SKILL functions compute a data value known as the return value of the function. Throughout this document, the right arrow (=>) denotes the return value of a function call. You can

- Assign the return value to a SKILL variable
- Pass the return value to another SKILL function

Any type of data can be a return value.

# **Syntax Functions for Defining Functions**

SKILL supports the following syntax functions for defining functions. You should use the procedure function in most cases.

# procedure

The procedure function is the most general and is easiest to use and understand.

The procedure function provides the standard method of defining functions. Its return value is the symbol with the name of the function. For example:

```
procedure( trAdd( x y )
    "Display a message and return the sum of x and y"
    printf( "Adding %d and %d ... %d \n" x y x+y )
    x+y
    ) => trAdd
trAdd( 6 7 ) => 13
```

## Working with SKILL

### **Terms and Definitions**

function, procedure

In SKILL, the terms *procedure* and *function* are used interchangeably to refer to a parameterized body of code that can be executed with parameters in the function call bound to the parameters in the function definition. SKILL can represent a function as both a hierarchical list and as a function object.

argument, parameter The terms argument and parameter are used interchangeably.

The arguments in a function call correspond to the formal

arguments in the declaration of the function.

expression A use of a SKILL function, often by means of an operator

supplying required parameters.

function body The collection of SKILL expressions that define the function's

algorithm.

Working with SKILL

# **OCEAN Environment Commands**

The following OCEAN environment commands let you start, control, and quit the OCEAN environment.

<u>appendPath</u>

<u>path</u>

prependPath

<u>setup</u>

**history** 

<u>ocnSetSilentMode</u>

### **OCEAN Environment Commands**

# appendPath

```
appendPath(
    t_dirName1 ... [ t_dirNameN ]
   )
   => t dirNameN / nil
```

# **Description**

Appends a new path to the end of the search path list. You can append as many paths as you want with this command.

## **Arguments**

t_dirName1	Directory path.
t_dirNameN	Additional directory paths.

#### Value Returned

t_dirNameN	Returns the last path specified.
nil	Returns nil and prints an error message if the paths cannot be appended.

## **Example**

```
appendPath( "/usr/mnt/user/processA/models" )
=> "/usr/mnt/user/processA/models"
```

Adds /usr/mnt/user/processA/models to the end of the current search path.

```
appendPath( "/usr/mnt/user/processA/models" "/usr/mnt/user/processA/models1")
=> "/usr/mnt/user/processA/models"
```

Adds /usr/mnt/user/processA/models and /usr/mnt/user/processA/models1 to the end of the current search path.

### **OCEAN Environment Commands**

# path

```
path(
    t_dirName1 ... [ t_dirNameN ]
)
=> 1 pathList / nil
```

## **Description**

Sets the search path for included files.

This command overrides the path set earlier using any of these commands: <u>path</u>, <u>appendPath</u>, or <u>prependPath</u>.

Using this command is comparable to setting the Include Path for the direct simulator, or the modelPath for socket simulators in the Virtuoso® Analog Design Environment user interface. You can add as many paths as you want with this command.

## **Arguments**

t_dirName1	Directory path.
t_dirNameN	Additional directory path.

#### Value Returned

l_pathList	Returns the entire list of search paths specified.
nil	Returns nil and prints an error message if the paths
	cannot be set.

### **Example**

```
path( "~/models" "/tmp/models" )
=> "~/models" "/tmp/models"
```

Specifies that the search path includes /models followed by /tmp/models.

```
path()
=> "~/models" "/tmp/models"
```

Returns the search path last set.

### **OCEAN Environment Commands**

# prependPath

```
prependPath(
    t_dirName1 ... [ t_dirNameN ]
    )
    => undefined / nil
```

# **Description**

Adds a new path to the beginning of the search path list. You can add as many paths as you want with this command.

## **Arguments**

t_dirName1	Directory path.
t_dirNameN	Additional directory path.

#### Value Returned

undefined	The return value for this command/function is undefined.
nil	Returns nil and prints an error message if the paths cannot be added.

# **Example**

```
prependPath( "/usr/mnt/user/processB/models" )
=> "/usr/mnt/user/processB/models"
```

Adds /usr/mnt/user/processB/models to the beginning of the search path list.

```
prependPath( "/usr/mnt/user/processB/models" "/usr/mnt/user/processB/models2")
=> "/usr/mnt/user/processB/models"
```

Adds /usr/mnt/user/processB/models and /usr/mnt/user/processB/models2 to the beginning of the search path list.

```
prependPath()
=> "/usr/mnt/user/processB/models" "~/models" "/tmp/models"
```

Returns the search path last set.

# **OCEAN Environment Commands**

# setup

```
setup(
    [?numberNotation s_numberNotation ]
    [?precision x_precision ]
    [?reportStyle s_reportStyle ]
    [?charsPerLine x_charsPerLine ]
    [?messageOn g_messageOn ]
)
    => t / nil
```

# **Description**

Specifies default values for parameters.

# **OCEAN Environment Commands**

# Arguments

?numberNotation	Specifies the notation for printed information.
$s\_numberNotation$	Valid values: 'suffix, 'engineering, 'scientific, 'none
	Default value: 'suffix
	The format for each value is 'suffix: 1m, 1u, 1n, etc.; 'engineering: 1e-3, 1e-6, 1e-9, etc.; 'scientific: 1.0e-2, 1.768e-5, etc.; 'none.
	The value 'none is provided so that you can turn off formatting and therefore greatly speed up printing for large data files.
?precision x_precision	Specifies the number of significant digits that are printed. Valid values: 1 through 16
	Default value: 6
?reportStyle s_reportStyle	Specifies the format of the output of the <u>report</u> command. Valid values: spice, paramValPair
	Default value: paramValPair
	The spice format is:
	Param1 Param2 Param3
	Name1valuevalue
	Name2valuevalue
	Name3valuevalue
	The paramValPair format is:
	Name1
	Param1=value Param2=value Param3=value
	Name2
	Param1=value Param2=value Param3=value
	Name3
	Param1=value Param2=value Param3=value

### **OCEAN Environment Commands**

?charsPerLine Specifies the number of characters per line output to the

 $x\_charsPerLine$  display.

Default value: 80

?messageOn  $g_messageOn$  Specifies whether error messages are turned on.

Valid values: t, nil

Default value: t, which specifies that messages are

turned on.

#### Value Returned

t Returns t if the value is assigned to the name.

nil Returns nil if there is a problem.

### **Example**

```
setup( ?numberNotation 'engineering )
=> t
```

Specifies that any printed information is to be in engineering mode by default.

```
setup( ?precision 5 )
=> t
```

Specifies that 5 significant digits are to be printed.

```
setup(?numberNotation 'suffix ?charsPerLine 40 ?reportStyle 'spice ?messageOn t)
```

Sets up number notation to suffix format, characters per line to 40, reporting style to Spice, and error message to ON.

#### **OCEAN Environment Commands**

# history

## **Description**

Displays the command history. By default, it prints the last 20 commands from the current session and the most recently terminated session. More commands can be printed by giving a number as an argument.

### **Arguments**

 $x_number$ 

The number of previously entered commands to be listed.

Default value: 20

### Value Returned

t

Returns t to indicate that the commands from history have been listed.

# **Example**

```
history
1 simulator('spectre)
2 design( "tests" "simple" "schematic")
3 analysis( 'tran ?start 0 ?stop 1u ?step 10n )
4 run()
=> t
```

Displays the most recently used commands. To reuse any of these commands, use the following methods at the ocean prompt:

■ ocean> !1

This executes the command numbered 1, which in this example is simulator ('spectre).

ocean> !des

This executes the last command whose prefix starts with des in the history. In this example, it is the second command listed, that is, design( "tests" "simple" "schematic").

# **OCEAN Environment Commands**

Note: To run history in CIW, the syntax is:

<space>!<commandNumber>

For example:

<space>!1

This executes the command numbered 1 from the CIW.

### **OCEAN Environment Commands**

### ocnSetSilentMode

```
ocnSetSilentMode(
    g_silentMode
)
=> t
```

# **Description**

Filters out OCEAN warning and information messages and allows only error messages to be written. This functionality is useful while running the OCEAN scripts when you might want to skip all OCEAN messages except errors.

### **Arguments**

g\_silentMode Accepts boolean values t or nil.

Set to t to suppress the OCEAN warning and information

messages.

Set to nil to allow all OCEAN messages to be displayed.

## **Value Returned**

t

Returns  $\ensuremath{\mathtt{t}}$  to indicate the successful assignment of the passed argument.

## **Example**

```
ocnSetSilentMode(t) => t
```

Suppresses the ocean warning messages

```
ocnSetSilentMode(nil) => t
```

Displays the ocean warning messages

# **Simulation Commands**

The following OCEAN simulation commands let you set up and run your simulation.

<u>ac</u>

analysis

<u>converge</u>

connectRules

<u>createFinalNetlist</u>

createNetlist

<u>dc</u>

definitionFile

<u>delete</u>

deleteOpPoint

<u>design</u>

<u>desVar</u>

discipline

displayNetlist

**envOption** 

<u>evcdFile</u>

<u>evcdInfoFile</u>

forcenode

globalSigAlias

globalSignal

ic

# **Simulation Commands**

<u>includeFile</u>
<u>modelFile</u>
nodeset
<u>noise</u>
<u>ocnCloseSession</u>
<u>ocnDisplay</u>
<u>ocnDspfFile</u>
ocnSpefFile
<u>ocnPspiceFile</u>
ocnGetAdjustedPath
<u>ocnGetInstancesModelName</u>
<u>off</u>
<u>option</u>
<u>restore</u>
<u>resultsDir</u>
<u>run</u>
<u>save</u>
<u>saveOpPoint</u>
<u>saveOption</u>
<u>simulator</u>
<u>solver</u>
<u>stimulusFile</u>
<u>store</u>
<u>temp</u>
<u>hlcheck</u>
<u>ocnAmsSetOSSNetlister</u>
<u>ocnAmsSetUnlNetlister</u>

### Simulation Commands

### ac

```
ac(
    g_fromValue
    g_toValue
    g_ptsPerDec
)
    => undefined / nil
ac(
    g_fromValue
    g_toValue
    t_incType
    g_points
)
    => undefined / nil
```

# **Description**

Specifies an AC analysis.

To know more about this analysis, see the simulator-specific user guide.

#### Simulation Commands

### **Arguments**

g\_fromValue Starting value for the AC analysis.

g\_toValue Ending value.

g ptsPerDec Points per decade.

*t\_incType* Increment type.

Valid values:

For the Spectre<sup>®</sup> circuit simulator, "Linear",

"Logarithmic", or "Automatic". For other simulators,

"Linear" or "Logarithmic".

g\_points Either the linear or the logarithmic value, which depends on

t\_incType.

#### Value Returned

undefined The return value for this command/function is undefined.

nil Returns nil and prints an error message if the analysis is not

specified.

### **Example**

ac(1 10000 2)

Specifies an AC analysis from 1 to 10,000 with 2 points per decade.

ac(1 10000 "Linear" 100)

Specifies an AC analysis from 1 to 10,000 by 100.

ac(1 5000 "Logarithmic" 10)

Specifies an AC analysis from 1 to 5000 with 10 logarithmic points per decade.

# **OCEAN Reference** Simulation Commands

# analysis

```
analysis(
    s_analysisType
    [ ?analysisOption1 g_analysisOptionValue1 ]... [ ?analysisOptionN
    g_analysisOptionValueN ]
    )
    => undefined / nil
```

# Description

Specifies the analysis to be simulated.

You can include as many analysis options as you want. Analysis options vary, depending on the simulator you are using. To include an analysis option, replace <code>analysisOption1</code> with the name of the desired analysis option and include another argument to specify the value for the option. If you have an AC analysis, the first option/value pair might be <code>[?from 0]</code>.

**Note:** Some simplified commands are available for basic SPICE analyses. See the ac, dc, tran, and noise commands. Use the ocnHelp( 'analysis') command for more information on the analysis types for the simulator you choose.

#### Simulation Commands

### **Arguments**

s analysisType

Type of the analysis. The valid values for this argument depend on the analyses that the simulator contains. The basic SPICE2G-like choices: 'tran, 'dc, 'ac, and 'noise.

?analysisOption1 g analysisOptionValue1

Analysis option. The analysis options available depend on which simulator you use. (See the documentation for your simulator.) If you are using the Spectre® circuit simulator, see the information about analysis statements in the *Virtuoso Spectre Circuit Simulator Reference* for analysis options you can use.

?analysisOptionN g analysisOptionValueN

Any subsequent analysis option. The analysis options that are available depend on which simulator you use. (See the documentation for your simulator.)

#### Value Returned

undefined

The return value for this command/function is undefined.

nil

Returns nil and prints an error message if there is a problem specifying the analysis.

### **Examples**

## Example1

```
analysis( 'ac ?start 1 ?stop 10000 ?lin 100 )
```

The above example runs an AC analysis from 1Hz to 10000Hz with a linear step of 100Hz for the Spectre® circuit simulator.

#### Example 2

```
analysis('tran ?start 0 ?stop 1u ?step 10n )
```

The above example specifies that a <u>transient</u> analysis is performed from 0 to 1u with a sweep value of 10n for the Spectre® circuit simulator.

#### Simulation Commands

## Example 3

The above example specifies that the temperature sweep starts at -50 and stops at 100 for the Spectre® circuit simulator when DC analysis is performed.

### Example 4

```
analysis('dc ?saveOppoint t )
```

The above example saves the operating point information when performing Spectre DC analysis.

### Example 5

```
analysis('xf ?start 0 ?stop 100 ?lin 2 ?dev "v3" ?param "dc" ?freq 1 ?probe "v4")
```

The above example specifies that a dc <u>transfer function</u> analysis is performed from 0 to 100 with a linear sweep of 2 and frequency of 1 on the probe v4.

# Example 6

```
analysis('sens ?analyses_list list("dcOp" "dc" "ac") ?output_list list("I7:3"
"OUT")
```

The above example specifies that a Spectre <u>sensitivity</u> analysis is performed with the dcOp, dc and ac analyses and the output is saved to the instance 17:3.

### Example 7

```
analysis ('noise ?start 1 ?stop 10e6 ?oprobe "V4")
```

The above example specifies that a Spectre noise analysis is performed from 1 to 10e6 and probes the operating point on instance V4.

# Example 8

```
analysis( 'dcmatch ?oprobe "/PR1" )
analysis( 'dcmatch ?param "temp" ?start "24" ?stop "26 ?lin "5" )
```

The example above specifies that a Spectre dcmatch analysis for the temperature of operating point PR1 is performed from 24 to 26°F with a linear sweep of 5.

#### Simulation Commands

## Example 9

```
analysis('pz ?freq "2" ?readns "./abc" ?oppoint "rawfile" ?fmax "4.5GHz" ?zeroonly
"no" ?prevoppoint "no" ? restart "no" ?annotate "no" ?stats "no"
)
```

The example above specifies that a Spectre pz analysis is performed on the operating point rawfile with a frequency of 2 and fmax of 4.5GHz.

### Example 10

```
analysis('stb ?start "10" ?stop "10G" ?dec "10" ?probe "/PR1" ?prevoppoint "yes" ?readns "./abc" ?save "lvl" ?nestlvl "1" ?oppoint "logfile" ?restart "yes" ?annotate "no" ?stats "yes" )
```

The above example specifies that a Spectre stability analysis is performed from 10 to 10G on the probe instance PR1.

### Example 11

```
analysis('pss ?fund "100M" ?harms "3" ?errpreset "moderate" )
```

The above example specifies that the Spectre <u>pss</u> RF analysis is performed with fundamental frequency of 100M, solution harmonics set to 3, and a moderate error level.

#### Example 12

```
analysis('pnoise ?start "1K" ?stop "30M" ?log "20" ?maxsideband "3"
?oprobe "/rif" ?iprobe "/rf" ?refsideband "0" )
```

The above example specifies that a Spectre <u>pnoise</u> RF analysis is performed from 1K to 30M with a maxsideband of 3, a logarithmic sweep of 20, and input and output probes set to rf and rif respectively.

### Example 13

```
analysis('pac ?sweeptype "relative" ?relharmnum "" ?start "700M" ?stop "800M"
?lin "5" ?maxsideband "3")
```

The above example specifies that a Spectre <u>pac</u> RF analysis is performed from 700M to 800M with a relative linear sweep of 5 and a maxsideband of 3.

# Example 14

```
analysis('pxf ?start "10M" ?stop "1.2G" ?lin "100" ?maxsideband "3" ?p "/Plo" ?n "/qnd!" )
```

#### Simulation Commands

The above example specifies that a Spectre <u>pxf</u> RF analysis is performed from 10M to 1.2G with a linear sweep of 100 and maxsideband of 3 between the open terminals Plo and qnd!.

### Example 15

```
analysis('qpss ?funds list("flo" "frf") ?maxharms list("0" "0") ?errpreset
"moderate" ?param "prf" ?start "-25" ?stop "-10" ?lin "5" )
```

The example above specifies that a Spectre <u>qpss</u> RF analysis is performed from -25 to -10 with a linear sweep of 5 for the fundamental frequencies of flo and frf. Here, the number of harmonics of each fundamental to be considered is 0, the number of parameters to be updated are saved in prf, and the error level is set to moderate.

### Example 16

```
analysis('qpac ?start "920M" ?stop "" ?clockmaxharm "0" )
```

The above example specifies that a Spectre <u>qpac</u> analysis is performed from 920M with clockmaxharm set to 0.

## Example 17

```
analysis('sp ?start "100M" ?stop "1.2G" ?step "100" ?donoise "yes" ?oprobe "/PORT0" ?iprobe "/RF" )
```

The above example specifies that the Spectre sp (S - parameter) analysis is performed from 100M to 1.2G with a sweep value of 100 between the open terminals /Port0 and /RF. Additionally, noise analysis is also performed.

### Simulation Commands

# converge

```
converge(
    s_convName
    t_netName1
    f_value1 ... [ t_netNameN f_valueN ]
)
    => undefined / nil
```

# **Description**

Sets convergence criteria on nets.

To know more about convergence, refer to the chapter <u>Helping a Simulation to Converge</u> of the *Virtuoso Analog Design Environment L User Guide*.

# **Arguments**

s_convName	Name of the convergence type. Valid values are one of nodeset ic and forcenode.
	<b>Note:</b> forcenode is not supported for the spectre simulator.
t_netName1	Name of the net to which you want to set convergence criteria.
f_value1	Voltage value for the net
t_netNameN	Name of the additional net
f_value	Voltage value for the additional net

### Value Returned

undefined	The return value for this command/function is undefined.
nil	Returns nil and prints an error message if the function fails.

### **Example**

```
converge( 'ic "/I0/net1" 5 )
```

Sets the convergence name for the initial condition net1 to 5 volts.

# Simulation Commands

converge( 'nodeset "/I0/net1" 5 )

Sets the convergence name for nodeset of net1 to 5 volts.

## connectRules

```
connectRules(
     t ruleName
     [ ?lib t libName ]
     [ ?view t viewName ]
     [ ?baseRule t baseRule ]
     [ ?moduleInfo 1 moduleInfo ]
     [ ?resolutionInfo 1 resolutionInfo ]
     [ ?commonParam 1_commonParam ]
     [ ?userDefined s userDefined ]
    => t / nil
connectRules (
     t ruleName
     => t / nil
connectRules (
     ?none s tag
    => t / nil
```

The following arguments are composed of other arguments as described below:

# Description

Sets connect rules for a given AMS OCEAN session required by the elaborator. To specify multiple connect rules, use this command multiple times. To add a connect rule to an OCEAN session, you can either choose a built-in rule from the <code>connectLib</code> library (by specifying <code>t\_ruleName</code>, <code>t\_libName</code> and <code>t\_viewName</code>) or one of your own compiled built-in connect rules (by specifying <code>t\_ruleName</code>, <code>t\_libName</code> and <code>t\_viewName</code>). To add a user defined connect rule to an OCEAN session specify <code>s\_userDefined</code>. To modify an existing built-in rule, you need to specify <code>t\_baseRule</code> (the name of the built-in rule that needs be modified), specify a new name (by specifying <code>t\_ruleName</code>, <code>t\_libName</code> and <code>t\_viewName</code>) and also specify one or more of the optional arguments.

### Simulation Commands

You can use the <code>delete('connectRules)</code> command to delete one or more specified connect rules. See the examples provided with the <u>delete</u> command.

You can use <code>ocnDisplay('connectRules)</code> to view the currently active connect rules in an OCEAN session. You may use <code>ocnDisplay('connectRules' all)</code> to display all information about all active connect rules in an OCEAN session.

Note: This command is applicable only when ams is the selected simulator.

#### Simulation Commands

## **Arguments**

t_ruleName	Name of the connect rule that you want to use in the

current session.

t\_libName Name of the library that contains a list of user-compiled

connect rules. If you do not specify this, the connect rules

are assumed to be in the default location.

t viewName Name of the view of the selected cell.

t baseRule Name of the connect rule that you want to modify.

1\_moduleInfo Arguments that need to be updated for a specified connect rule. The arguments may include s mode,

s direction1, s direction2, s discipline1 and

s discipline2.

Valid values for s mode are: null, split, merged.

s\_direction1 and s\_direction2 work as a pair. Valid combinations are: both null, input/output,

output/input, inout/inout.

s\_discipline1 and s\_discipline2 also work as a pair. Either they should both be null or they should both

have values.

t\_resolutionInfo Names of disciplines that need to be resolved to another

discipline. The value specified overwrites the

1 resolutionInfo in the base rule or in the existing

connect rule.

t commonParam One or more parameters that you want to modify for all

modules or a set of modules. Although the same result can be achieved by using the <code>l\_moduleInfo</code> argument, <code>l commonParam</code> facilitates updating parameters for all

modules in one go.

#### Simulation Commands

$s\_userDefined$	Name of the user defined connect rule that you want to
	use in the current session. Specify 3step as the value of
	<pre>s_userDefined and specify t_ruleName, t_libName</pre>
	and t_viewName to add a user defined connect rule for
	the Cellview-based netlister flow. Specify irun as the
	value of s userDefined and specify t ruleName,

s fileName or both to add a user defined connect rule for the OSS-based netlister with irun flow. Any other argument specified when adding a user defined connect

Name of the user defined connect rule that you want to

rule will be ignored.

The option used to indicate that no connect rules are to s tag

be used for the current session.

#### Value Returned

Returns t if the specifed connect rules are set. t

nil Returns nil and prints an error message otherwise.

# **Example**

connectRules("ConnRules 5V full")

Sets ConnRules 5V full as the current connect rule from the default connectLib located in your hierarchy.

```
connectRules ("CustomRules 9V high" ?lib "myConnectLib" ?view "myViewName")
```

Sets CustomRules 9V high from myConnectLib, where myConnectLib contains a list of user-compiled connect rules and myViewName is the specified view name.

```
connectRules("connRule3" ?lib "lib2" ?view "view2" ?baseRule "ConnRules 18V full"
?description "updated directions" ?moduleInfo ((?name "E2L" ?direction1 "input" ?direction2 "output")))
```

Checks if connRule3 exists in the session. If it does, it updates direction1 to input and direction2 to output for E2L and description for this rule. If this rule does not exist, then it takes the base values as values from ConnRules 18V full and updates direction1, direction2, and description and names the new rule as connRule3.

```
connectRules("connRule3" ?lib "lib2" ?view "view2" ?moduleInfo ((?name "E2L" ?mode
"split")))
```

Checks if connRule3 exists. If it does not exist, as no base rule is specified, a relevant error message appears. If the rule exists, it would update mode to split for the existing connect rule connRule3 for the module E2L.

#### Simulation Commands

```
connectRules("connRule3" ?lib "lib2" ?view "view2" ?description "desc123"
?moduleInfo ((?name "E2L" ?mode "split" ?direction1 "input" ?direction2 "output"))
?resolutionInfo nil)
```

If connRule3 does not exist and the base rule is not specified but description, moduleInfo and resolutionInfo are specified, the connect rule connRule3 is added with the values specified for moduleInfo, resolutionInfo and description.

**Note:** In this case no checks are done (that is, module names and parameter names are not checked against base information as no base rule information is available). This command is applicable while using the connectRules command as saved in ocean.

```
connectRules("connRule3" ?lib "lib2" ?view "view2" ?moduleInfo ((?name "L2E"
?paramInfo (("vsup" "1.7")("vtlo" "3.2")))
```

Updates the parameters vsup and vtlo for the existing rule connRule3 in the L2E module.

```
connectRules("connRule3" ?lib "lib2" ?view "view2" ?resolutionInfo (("r1" "e1
e2")("r2" "e4 e5")) ?commonParam (("vsup" "1.2") ("vtlo" "3.4" ("L2E" "Bidir"))
```

Updates resolutionInfo for the existing connect rule connRule3. The old resolutionInfo value for this rule is replaced with the new information. It also updates the vsup parameter to 1.2 for all connRule3 modules and updates vtlo to 3.4 for the modules L2E and Bidir.

```
connectRules("connRule3" ?lib "lib2" ?view "view2" ?userDefined 3step)
```

Sets connRule3 from view2 of lib2 as a user defined connect rule for the Cellview-based netlister flow.

```
connectRules("connRule3" ?userDefined irun)
```

Sets connRule3 from the connectLib library as a user defined connect rule for the OSS-based netlister with irun flow.

```
connectRules("connRule3" ?userDefined irun ?file "file1")
```

Sets connRule3 from file1 as a user defined connect rule for the OSS-based netlister with irun flow.

```
connectRules(?userDefined irun ?file "file1")
```

No user-defined connect rule name is specified for the OSS-based netlister with irun flow. Hence, the first rule found in file1 will be used for AMS simulation.

```
connectRules(?none t)
=> t
```

Sets the current connect rule to None so that no connect rule is provided to ncelab during elaboration.

```
delete('connectRules list("mylib" "myrule" "myview") list("mylib1" "myrule1"
"myview1"))
```

### Simulation Commands

Deletes the connect rule myrule in the library mylib with the view myview. It also deletes the connect rule myrule1 in the library mylib1 with the view myview1.

```
delete('connectRules list("" "rule1" ""))
```

Deletes the specified connect rule rule1 from the default connectLib library.

### Simulation Commands

# createFinalNetlist

```
createFinalNetlist(
    )
    => t / nil
```

# **Description**

Creates the final netlist for viewing purposes. The netlist also can be saved but is not required to run the simulator.

**Note:** This command works only for socket simulators. For direct simulators, such as spectre, use createNetlist instead.

# **Arguments**

None.

#### Value Returned

t Returns t if the final netlist is created.

nil Returns nil and prints an error message otherwise.

### **Example**

```
createFinalNetlist()
```

Creates the final netlist for the current simulation run.

#### Simulation Commands

### createNetlist

```
createNetlist(
    [ ?recreateAll g_recreateAll ]
    [ ?display g_display ]
    )
    => t filename / nil
```

# **Description**

Creates the simulator input file.

If the design is specified as cellview, this command netlists the design, if required, and creates the simulator input file. When the  $g_recreateAll$  argument is set to t and the design is specified as cellview, all the cells in the design hierarchy are renetlisted, before creating the simulator input file. If the design is specified as netlist file, that netlist is included in the simulator input file. Also see the <u>design</u> function.

When the  $g_{display}$  option is set to t (or nil) the netlist file is displayed (or undisplayed) to the user. By default,  $g_{display}$  it set to 't (true).

Specifies if the netlist is to be displayed or not.

Note: This command does not work with socket simulators.

### **Arguments**

```
?g_recreateAll g_recreateAll Specifies if the netlist needs to be recreated or not. ?display g_display
```

#### Value Returned

t_fileName	Returns the name of the simulator input file.
nil	Returns nil otherwise

# **Example**

```
createNetlist()
=> "/usr/foo/netlist/input.scs"
```

Creates simulator input file for the current simulation run.

#### Simulation Commands

```
design( "test" "mytest" "spectre")
createNetlist( ?recreateAll t )
=>"/usr/foo/netlist/input.scs"
```

Netlists and creates simulator input file for the current simulation run.

```
design( "test" "mytest1" "spectre")
createNetlist( ?recreateAll t ?display nil )
=>"/usr/foo/netlist/input.scs"
```

Netlists and creates simulator input file for the given simulation run but does not display the input.scs file in a new window which may be annoying to the user. By default ?display option is set to 't meaning netist file would be displayed. This can be turned ON/OFF via ?display set to t/nil

**Note:** If you regenerate the netlist after changing the design in a different Virtuoso session, the netlist is not updated with the design changes. To update the netlist with the current cellview, run the <a href="mailto:ddsRefresh">ddsRefresh</a> command before running the <a href="mailto:createNetlist">createNetlist</a> command as shown below:

```
ddsRefresh( ?cellview t )
=> t
createNetlist( ?recreateAll t )
=> "/usr/foo/netlist/input.scs"
```

### Simulation Commands

# dc

```
dc(
    t_compName
    [ t_compParam ]
    g_fromValue
    g_toValue
    g_byValue
)
    => undefined / nil
```

# **Description**

Specifies a DC sweep analysis with limited options. If other analysis options are needed, use the  ${\tt analysis}$  command.

To know more about this analysis, see the simulator-specific user guide.

**Note:**  $t\_compParam$  is valid only for the spectre simulator.

# **Arguments**

t_compName	Name of the source (or component, for the Spectre® circuit simulator) to sweep.
t_compParam	For the Spectre® circuit simulator, the component parameter to be swept.
g_fromValue	Starting value for the DC analysis.
g_toValue	Ending value.
g_byValue	The increment at which to step through the analysis.

#### Value Returned

undefined	The return value for this command/function is undefined.
nil	Returns nil and prints an error message if the analysis is not specified.

# **Example**

```
dc("v1" "dc" 0 5 1)
dc("r1" "r" 0 5 1)
```

# Simulation Commands

Specifies two DC sweep analyses for the Spectre® circuit simulator.

dc("v1" 0 5 1)

Specifies one DC sweep analysis for a simulator other than the Spectre® circuit simulator.

#### Simulation Commands

# definitionFile

```
definitionFile(
    t_fileName [ t_fileName2 ... t_fileNameN ]
   )
    => 1 fileNames / nil
```

### **Description**

Specifies definitions files to be included in the simulator input file.

Definitions files define functions and global variables that are not design variables. Examples of such variables are model parameters or internal simulator parameters. To know more about definitions files, see the section *Using a Definitions File* in *Chapter 3* of the *Virtuoso Analog Design Environment L User Guide*.

**Note:** This command does not work with socket simulators.

## **Arguments**

t_fileName	The name of the definition file that would typically contain
	functions or parameter statements.

#### Value Returned

$1\_file extit{Names}$	A list of the file names specified; returned on success.
nil	Otherwise nil is returned.

## **Example**

```
definitionFile( "functions.def" "constants.def" )
=> ("functions.def" "constants.def")
```

Includes functions.def and constants.def files in the simulator input file.

```
definitionFile( )
=> ("functions.def" "constants.def")
```

Returns the definition files set earlier.

### Simulation Commands

# delete

```
delete(
    s_command
    [ g_commandArg1 ] [ g_commandArg2 ] ...
)
    => t / nil
```

# **Description**

Deletes all the information specified.

The  $s\_command$  argument specifies the command whose information you want to delete. If you include only this argument, all the information for the command is deleted. If you supply subsequent arguments, only information specified by these arguments is deleted, and not all the information for the command.

# **Arguments**

s_command	Command that was initially used to add the items that are now being deleted.
	Valid values: analysis, connectRules, discipline, globalSignal, desVar, path, save, ic, forcenode, nodeset
$g\_commandArg1$	Argument corresponding to the specified command.
g_commandArg2	Additional argument corresponding to the specified command.

#### Value Returned

t Returns t if the information is deleted.

nil Returns nil if there is an error.

# **Example**

```
delete( 'save )
=> t
```

#### Deletes all the saves.

```
delete( 'save 'v )
=> t
```

### Simulation Commands

Deletes *only* the nets. The rest of the information can be saved in subsequent simulations.

```
delete( 'save "net23" )
=> t
```

Deletes only net23. The rest of the information can be saved in subsequent simulations.

```
delete( 'monteCarlo )
=> t
```

Turns off the monteCarlo command and sets everything back to the defaults.

# **OCEAN Reference** Simulation Commands

# deleteOpPoint

```
deleteOpPoint(
    t_instName
    [ @rest l_args ]
)
    => t / nil
```

# **Description**

Deletes the specified operating point instance.

# **Arguments**

t_instName	Name of the operating point instance to be deleted.
@rest <i>l_args</i>	List of optional arguments that can be passed to this function.

### **Values Returned**

t	Returns t when the function runs successfully.
nil	Otherwise, returns nil if there is an error.

# **Example**

```
deleteOpPoint( "/I8/Q3" )
```

This example deletes the operating point instance 18/Q3.

### Simulation Commands

# design

# **Description**

Specifies the directory path to the netlist of a design or the name of a design to be simulated.

For the *lib*, *cell*, *view* version of the design command, you can specify the mode (r, w, or a, representing read, write, or append) in which the design should be opened.

#### Simulation Commands

# **Arguments**

t_cktFile	Directory path to the netlist followed by the name of
-----------	---

the netlist file. Name of the netlist file must be

netlist. The netlistHeader and

netlistFooter files must be in the same directory.

Otherwise, cktFile is a pre-existing netlist file from

another source.

t 1ib Name of the library that contains the design.

t cell Name of the design.

t view View of the design (typically schematic).

t\_mode The mode in which the design should be opened. The

value can be r, w or a, representing read, write, and append, respectively. The default mode is append. Read-only designs can be netlisted only by direct netlisters, and not socket. The w mode should

not be used as it overwrites the design.

#### Value Returned

t	cktFile	Returns the name of the design if successful.

l (lib cell view)

Returns the name of the view for an Virtuoso® Analog

Design Environment design if successful.

nil Returns nil and prints an error message if there is a

problem using the specified design.

### **Examples**

#### Example 1

```
design( "./opampNetlist/netlist" )
=> netlist
```

specifies that netlist, a netlist file, be used in the simulation.

### Simulation Commands

# Example 2

```
design( "tests" "simple" "schematic" )
=> (tests simple schematic)
```

Specifies that the schematic view of the simple design from your tests library be used in the simulation.

# Example 3

```
design("mylib" "ampTest" "schematic" "a")
=> (mylib ampTest schematic)
```

Specifies that the schematic view of the ampTest design from your mylib library be appended to the simulation.

# Example 4

```
design()
=> (mylib ampTest schematic)
```

Returns the lib-cell-view being used in the current session. If a design has not been specified, it returns nil.

#### Simulation Commands

### desVar

```
desVar(
    t_desVar1 f_value1 ... [ t_desVarN f_valueN ]
   )
    => undefined / nil
```

# **Description**

Sets the values of design variables used in your design. You can set the values for as many design variables as you want.

To know more about design variables, refer to the Chapter 3, <u>Design Variables and Simulation Files for Direct Simulation</u> of the Virtuoso Analog Design Environment L User Guide.

# **Arguments**

t_desVar1	Name of the design variable.
f_value1	Value for the design variable.
t_desVarN	Name of an additional design variable.
f_valueN	Value for the additional design variable.

#### Value Returned

undefined	The return value for this command/function is undefined.
nil	Returns nil and prints an error message if the assignments fail.

## **Example**

```
desVar()
```

Returns the design variables set last, if any. Otherwise, it returns nil.

```
desVar( "rs" 1k )
```

Sets the rs design variable to 1k.

```
desVar( "r1" "rs" "r2" "rs*2" )
```

Sets the r1 design variable to rs, or 1k, and sets the r2 design variable to rs\*2, or 2k.

# Simulation Commands

a = evalstring( desVar( "rs")) / 2

Sets a to 1k/2 or 500.

**Note:** evalstring is necessary because desVar returns a string.

#### Simulation Commands

# discipline

```
discipline(
    g_discipline1 [ g_discipline2 ... ]
)
    => t / nil
```

# **Description**

Adds discrete disciplines to the existing set of disciplines for a given 'ams' OCEAN session. You can use delete('discipline) to delete one or more specified disciplines. You can use ocnDisplay('discipline) to view the currently active disciplines in an OCEAN session.

**Note:** This command is applicable only when ams is the simulator.

## **Arguments**

g_discipline1	Name of a discrete discipline to be added.
g_discipline2	Names of additional discrete disciplines to be added.

### Value Returned

t	Returns t if the discipline is added.
nil	Returns nil or prints an error message otherwise.

## **Example**

```
discipline( "logic1" "logic2" '("logic3") )
```

Disciplines to be added can be either strings or lists containing the discipline name. If no disciplines have been added so far, this sample command adds the three discrete disciplines logic1, logic2 and logic3 to the session; otherwise, it adds these three disciplines to the existing set of disciplines.

```
discipline ("LL")
```

Adds discipline LL to the existing set of disciplines. If logic1, logic2 and logic3 are already added, LL is added as the fourth discipline.

```
delete('discipline "logic2" "LL")
```

Deletes disciplines logic2 and LL from the session.

# **OCEAN Reference** Simulation Commands

delete('discipline)

Deletes all the specified disciplines in the session.

#### Simulation Commands

# displayNetlist

```
displayNetlist(
    )
    => t / nil
```

# **Description**

Displays the concatenated AMS complete design info file used in a given AMS OCEAN session. The concatenated file displays the cell-based netlisting of the cellviews used in the configuration along with the analog control file and the TCL file generated by AMS-ADE. This command is applicable for both solvers – Spectre and UltraSim.

**Note:** This command is applicable only when ams is the simulator.

# **Arguments**

None.

#### Value Returned

t

Returns t if the concatenated design information file.

nil

Returns nil or prints an error message otherwise.

# **Example**

```
displayNetlist()
=> t
```

Displays the concatenated design information file.

#### Simulation Commands

# envOption

```
envOption(
    s_envOption1 g_value1 ... [ s_envOptionN g_valueN ]
   )
    => undefined / nil
```

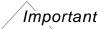
### **Description**

Sets environment options.

To get the list of environment options that can be set for a simulator, first set the simulator and then run the OCEAN online help command ocnHelp('envOption). For example,

```
simulator('spectre)
ocnHelp('envOption)
```

The above command displays a list of environment options that can be set for spectre.



To specify an include file, use the includeFile command, not the envOption command. To set a model path, use the path command, not the envOption command.

To know more about environment options, see the section *Environment Options* in *Chapter 2* of the *Virtuoso Analog Design Environment L User Guide*.

### Simulation Commands

# **Arguments**

s envOption1 Name of the first environment option to set.

g value1 Value for the option.

s envOptionN Name of an additional environment option to set.

g valueN Value for the option.

#### Value Returned

undefined The return value for this command/function is undefined.

nil Returns nil if there are problems setting the option.

### **Example**

```
envOption( 'paramRangeCheckFile "./myDir/range.check")
```

Sets the paramRangeCheckFile environment option.

```
envOption( 'initFile "./myDotSFiles/init" )
```

Sets the initFile environment option.

```
envOption( 'updateFile "./myDotSFiles/update")
```

Sets the updateFile environment option.

### Simulation Commands

### evcdFile

```
evcdFile(
    t_evcdFileName
)
    => t evcdFileName / nil
```

# **Description**

Sets an EVCD file for a given UltraSim OCEAN session. You also need to specify an EVCD info file while using this command. You can specify only one EVCD file for a session. You may use ocnDisplay ('evcdFile) to view the currently active EVCD file.

**Note:** This command is applicable for the UltraSim simulators.

### **Arguments**

t event i le value i i le le value i i session.	t	evcdFileName	The name of the EVCD file to be used for session.
---	---	--------------	---

#### Value Returned

t_evcdFileName	The EVCD file name is the output if the command is successful.
nil	Otherwise, nil is returned.

### **Example**

```
evcdFile("/tmp/evcdFile.dat")
=> "/tmp/evcdFile.dat"
```

Specifies /tmp/evcdFile.dat as the EVCD file to be used for current UltraSim OCEAN session.

#### Simulation Commands

### evcdInfoFile

```
evcdInfoFile(
    t_evcdInfoFileName
)
    => t evcdInfoFileName / nil
```

### **Description**

Sets a EVCD info file for a given UltraSim OCEAN session. You also need to specify an EVCD file while using this command. You can specify only one EVCD info file for a session. You may use ocnDisplay ('evcdInfoFile) to view the currently active EVCD info file.

**Note:** This command is applicable only for the UltraSim simulator.

### **Arguments**

```
t evcdInfoFileName
```

The name of the EVCD info file to be included.

### **Value Returned**

```
t evcdInfoFileName
```

The EVCD info file name is the output if the

command is successful.

nil Otherwise, nil is returned.

# **Example**

```
evcdInfoFile("/tmp/evcdInfoFile.dat")
=> "/tmp/vcdInfoFile.dat"
```

Specifies /tmp/evcdInfoFile.dat as the EVCD file to be used for current UltraSim OCEAN session.

### Simulation Commands

### forcenode

```
forcenode(
    t_netName1 f_value1 ... [ t_netNameN f_valueN ]
    )
    => undefined / nil
```

# **Description**

Holds a node at a specified value.

To know more about convergence, refer to the chapter <u>Helping a Simulation to Converge</u> of the *Virtuoso Analog Design Environment L User Guide*.

**Note:** This is not available for the spectre simulator. Refer to the documentation for your simulator to see if this feature is available for your simulator.

## **Arguments**

t_netName1	Name of the net.
f_value1	Voltage value for the net.
t_netNameN	Name of an additional net.
f_valueN	Voltage value for the net.

### Value Returned

undefined	The return value for this command/function is undefined.
nil	Returns nil and prints an error message.

# **Example**

```
forcenode( "net1" 5 "net34" 2 )
```

Sets the force nodes of "net1" to 5 and "net34" to 2.

#### Simulation Commands

# globalSigAlias

```
globalSigAlias(
    g_signalList1 [ g_signalList2 ... ]
)
=> t / nil
```

## **Description**

Removes all the previous signal aliases and creates the specified aliases. The signal names in each of the signal lists are marked as aliases of each other. Each of the signal lists is a set of signal names that are to be aliased. The signal names should match the names that were specified using the <u>globalSignal</u> command. To unalias all signal, pecify nil instead of signal lists.

**Note:** This command is applicable only when AMS is the simulator.

## **Arguments**

g_signalList(n)	A list of signals that are to be marked as aliases of each
	other.

#### Value Returned

t	Returns t when previous signal aliases have been removed successfully and new aliases are created according to the signal lists provided.
nil	Returns nil and prints an error message if the function was unsuccessful.

# Example

```
globalSigAlias('("sig1" "sig2") '("sig4" 'sig5" 'sig8"))
```

Removes the previous signal aliases and marks sig1 and sig2 as aliases of each other and sig4, sig5 and sig8 as aliases of each other. The signal names in each of the signal lists are marked as aliases of each other.

```
globalSigAlias("signal2" "signal6" "signal3")
```

If there is just one list of signals to be aliased, it can be given without the list. In this case, signal2, signal6 and signal3 are marked as aliases of each other.

### Simulation Commands

# globalSignal

```
globalSignal(
    [?name t_signalName]
    [?lang t_langName]
    [?wireType t_wireType]
    [?discipline t_discipline]
    [?ground t_ground]
    @Rest args
)
    => t / nil
```

# **Description**

Adds or modifies a global signal for a given AMS OCEAN session needed by the elaborator. If the global signal already exists in the session, the values are updated. If it does not exist, a global signal with the specified name is added. In case of a vector signal, the range information can be appended with the name of the signal.

**Note:** This command is applicable only when AMS is the simulator.

#### Simulation Commands

## **Arguments**

?name t\_signalName

The name of the global signal.

?lang t langName

The namespace within which the signal is entered. It is used to map the signal name to Verilog-AMS.

Valid Values: CDBA, Spectre, Spice, Verilog-

AMS

Default Value: CDBA

?wireType t wireType

Indicates the Verilog type of the signal

declaration.

Valid Values: wire, supply0, supply1, tri, tri0, tri1, triand, trior, trireq, wand, wor,

wreal

Default Value: wire

?discipline t discipline

A string value to indicate the discipline of

the signal.

?ground t ground

Valid Values: YES, NO

Default Value: NO

@Rest args List

List of optional arguments that can be passed to this function.

#### **Value Returned**

t Returns t when a global signal has been successfully added

or modified.

nil Returns nil and prints an error message if the function was

unsuccessful.

### Simulation Commands

### **Example**

```
globalSignal("signal1" ?wireType "tri")
```

Adds the global signal signal1 with wire type as tri, default language as CDBA, and ground as NO to the list of global signals if it has not already been added. If it already exists, then it updates the wire type for signal1.

```
globalSignal("signal2" ?lang "Spectre" ?discipline "electrical")
```

Adds signal2 with language as Spectre, discipline as electrical, and ground as NO to the list of global signals if it is not already added. If it already exists, then it updates language to Spectre and discipline to electrical.

```
delete('globalSignal "sig1" "sig2")
```

Deletes sig1 and sig2 after unaliasing them if they are in aliased sets.

```
delete('globalSignal)
```

Deletes all user-specified global signals.

### Simulation Commands

### ic

```
ic(
    t_netName1 f_value1 ... [ t_netNameN f_valueN ]
)
=> undefined / nil
```

# **Description**

Sets initial conditions on nets in a transient analysis.

To know more about convergence, refer to the chapter <u>Helping a Simulation to Converge</u> of the *Virtuoso Analog Design Environment L User Guide*.

# **Arguments**

t_netName1	Name of the net.
f_value1	Voltage value for the net.
t_netNameN	Name of an additional net.
f_valueN	Voltage value for the net.

### Value Returned

undefined	The return value for this command/function is undefined.
nil	Returns nil and prints an error message.

# Example

```
ic( "/net1" 5 "/net34" 2 )
```

Holds the nodes of "/net1" at 5 and "/net34" at 2.

#### Simulation Commands

### includeFile

```
includeFile(
    t_fileName
)
    => t fileName / nil
```

### **Description**

Includes the specified file in the final netlist of the simulator for the current session.

#### Notes:

- 1. This command is not available for the direct simulator. Use the modelFile or stimulusFile command instead.
- 2. Using this command is comparable to using the Environment Options form of the Virtuoso® Analog Design Environment to name an include file and specify that the syntax for the file be that of the target simulator. If you want the include file to be in Cadence-SPICE circuit simulator syntax, you must edit the raw netlist file (which has a .c or .C suffix), and manually add the include file.

# **Arguments**

t fileName Name of the file to include in the final netlist.

#### Value Returned

t_fileName	Returns the name of the file if successful.
nil	Returns nil and prints an error message otherwise.

### **Example**

```
includeFile( "~/projects/nmos" )
=> "~/projects/nmos"
```

Includes the nmos file in the final netlist of the simulator for the current session.

```
includeFile()
=>"~/projects/nmos"
```

Returns the includeFile, if one was set earlier. Otherwise, it returns nil.

#### Simulation Commands

### modelFile

```
modelFile(
        [ g_modelFile1 [ g_modelFile2 ... ] ]
)
=> 1 modelFile
```

### **Description**

Specifies model files to be included in the simulator input file.

This command returns the model files used. When model files are specified through the arguments, the model files are set accordingly. Use of full paths for the model file is recommended.

### **Arguments**

g_modelFile1	This argument can be a string to specify the name of the model file.
g_modelfile2	This argument can be a list of two strings to specify the name of the model file and the name of the section.

#### Value Returned

l_modelfile	A list of all the model file/section pairs.
nil	Returned when no file section pairs have been specified with the current call or a previous call of this command. The nil
	value is also returned when an error has been encountered.

### Example

```
modelFile( "bjt.scs" "nmos.scs" )
=>( ("bjt.scs" "") ("nmos.scs" "") )
modelFile( "bjt.scs" '("nmos.scs" "typ") 'my_models )
=> ( ("bjt.scs" "") ("nmos.scs" "typ") ("my_models" "") )
modelFile()
=> ( ("bjt.scs" "") ("nmos.scs" "") )
```

Returns the modelFile, if one was set earlier. Otherwise, it returns nil.

### Simulation Commands

### nodeset

```
nodeset(
    t_netName1 f_value1 ... [ t_netNameN f_valueN ]
)
=> undefined / nil
```

# **Description**

Sets the initial estimate for nets in a DC analysis, or sets the initial condition calculation for a transient analysis.

To know more about convergence, refer to the chapter <u>Helping a Simulation to Converge</u> of the *Virtuoso Analog Design Environment L User Guide*.

### **Arguments**

t_netName1	Name of the net.
f_value1	Voltage value for the net.
t_netNameN	Name of an additional net.
$f_valueN$	Voltage value for the net.

#### Value Returned

undefined	The return value for this command/function is undefined.
nil	Returns nil and prints an error message otherwise.

### **Example**

```
nodeset( "net1" 5 "net34" 2 )
```

Sets the initial estimates of "net1" to 5 and "net34" to 2.

### Simulation Commands

# noise

```
noise(
    t_output
    t_source
)
=> undefined / nil
```

# **Description**

Specifies a noise analysis.

**Note:** This command cannot be used with the spectre simulator.

# **Arguments**

t_	_output	Output node
t_	source	Input source

### Value Returned

undefined The return value for this command/function is undefined.

nil Returns nil and prints an error message If there is a problem

specifying the analysis.

### **Example**

```
noise( "n1" "v1" )
```

Specifies a noise analysis.

### Simulation Commands

## ocnCloseSession

```
ocnCloseSession(
    )
    => t / nil
```

# **Description**

Closes the current OCEAN session without saving any settings made during the session. The command has no effect if no session is currently active.

### **Value Returned**

t Returns t when the current session is successfully closed.

nil Returns nil if there is a problem closing the active session.

### **Example**

```
ocnCloseSession()
=> t
```

Closes the current OCEAN session.

#### Simulation Commands

# ocnDisplay

```
ocnDisplay(
    [?output t_filename | p_port ]
    s_command
    [g_commandArg1] [g_commandArg2] ...
)
=> t / nil
```

# Description

Displays all the information specified.

The  $s\_command$  argument specifies the command whose information you want to display. If you include only this argument, all the information for the command displays. If you supply subsequent arguments, only those particular pieces of information display as opposed to displaying all the information for that command. If you provide a filename as the <code>?output</code> argument, the <code>ocnDisplay</code> command opens the file and writes the information to it. If you provide a port (the return value of the SKILL <code>outfile</code> command), the <code>ocnDisplay</code> command appends the information to the file that is represented by the port.

### Simulation Commands

# **Arguments**

?output t_filename	File in which to write the information. The <code>ocnDisplay</code> command opens the file, writes to the file, then closes the file. If you specify the filename without a path, the <code>ocnDisplay</code> command creates the file in the directory pointed to by your SKILL Path. To find out what your SKILL path is, type <code>getSkillPath()</code> at the OCEAN prompt.
p_port	Port (previously opened with outfile) through which to append the information to a file. You are responsible for closing the port. See the <u>outfile</u> command for more information.
s_command	Command that was initially used to add the items that are now being displayed.
	Valid values: Most simulation setup commands. The commands that are supported include design, analysis, tran, ac, dc, noise, resultsDir, temp, option, desVar, path, includeFile, modelFile, stimulusFile, definitionFile, saveOption, envOption, save, converge, ic, forcenode, nodeset, simulator, setup, restore, saveSubckt

#### **Value Returned**

g commandArg1

g commandArg2

t Displays the information and returns t.

nil Returns nil and prints an error message if there are problems

Argument corresponding to the specified command.

Additional argument corresponding to the specified command.

displaying the information.

# Example

```
ocnDisplay( 'optimizeGoal )
=> t
```

Displays all the optimizeGoal information.

```
ocnDisplay( 'analysis 'tran )
=> t
```

Displays only transient analyses.

# **Simulation Commands**

```
ocnDisplay( 'save )
=> t
```

# Displays all the keeps.

```
ocnDisplay( ?output myPort 'analyis )
=> t.
```

Displays and writes all the analyses to the port named myPort.

#### Simulation Commands

# ocnDspfFile

```
ocnDspfFile(
    t_dspfFile [ t_dspfFile1 ... t_dspfFileN ]
   )
    => t dspfFile(s) / nil
```

# **Description**

Sets the parasitic (dspf, spf) files to be used in a Spectre OCEAN session. You can use this command to specify a list of parasitic files to be included in the control file. You can use ocnDisplay('dspfFile) to view the currently active parasitic (dspf, spf) files in an OCEAN session.

**Note:** This command is applicable for Spectre simulator. For AMS, it works only when Spectre is selected as the solver.

## **Arguments**

 $t\_dspfFile$  The name of the parasitic (dspf, spf) file to be included.  $t\_dspfFile1...t\_dspfFileN$ 

The name of the additional parasitic (dspf, spf) files to be included.

#### Value Returned

t\_dspfFile Lists the names of the parasitic (dspf, spf) files.

nil Returns nil if there are problems displaying the information.

## **Example**

```
ocnDspfFile("/tmp/file1.dspf" "/tmp/file2.dspf")
=> ("/tmp/file1.dspf" "/tmp/file2.dspf")
```

Displays the /tmp/file1.dspf and /tmp/file2.dspf parasitic files to be used for current Spectre OCEAN session.

#### Simulation Commands

# ocnSpefFile

```
ocnSpefFile(
    t_SpefFile [ t_SpefFile1 ... t_SpefFileN ]
   )
    => t SpecFile(s) / nil
```

# **Description**

Sets the parasitic (spef) files to be used in a Spectre OCEAN session. You can use this command to specify a list of parasitic files to be included in the control file. You can use ocnDisplay('SpefFile) to view the currently active parasitic (spef) files in an OCEAN session.

**Note:** This command is applicable for Spectre simulator. For AMS, it works only when Spectre is selected as the solver.

## **Arguments**

```
t\_SpefFile The name of the parasitic (spef) file to be included. t\_SpefFile1...t\_SpefFileN
```

The name of the additional parasitic (spef) files to be included.

#### Value Returned

```
t_SpefFile Lists the names of the parasitic (spef) files.

nil Returns nil if there are problems displaying the information.
```

# **Example**

```
ocnSpefFile("/tmp/file1.spef" "/tmp/file2.spef")
=> ("/tmp/file1.spef" "/tmp/file2.spef")
```

Displays the /tmp/file1.spef and /tmp/file2.spef parasitic files to be used for current Spectre OCEAN session.

#### Simulation Commands

# ocnPspiceFile

```
ocnPspiceFile(
    t_PSpiceFile
    [ t_PSpiceFile1 ... t_PSpiceFileN ]
    )
    => t PSpiceFile(s) / nil
```

# **Description**

Sets the PSpice files to be used in a Spectre OCEAN session. Use this command to specify a list of PSpice files to be included in the control file.

**Note:** This command is applicable for the Spectre simulator. For AMS, it works only when Spectre is selected as the solver.

# **Arguments**

```
t_PSpiceFile The name of the PSpice file to be included.

t PSpiceFile1...t PSpiceFileN
```

The name of the additional PSpice files to be included.

#### Value Returned

$t\_{PSpiceFile}$	Lists the names of the PSpice files.
nil	Returns nil if there are problems displaying the information.

### **Example**

```
ocnPspiceFile("/tmp/file1.sp" "/tmp/file2.sp")
=> ("/tmp/file1.sp" "/tmp/file2.sp")
```

Returns the /tmp/file1.sp and /tmp/file2.sp PSpice files to be used for the current Spectre OCEAN session.

### Simulation Commands

# ocnGetAdjustedPath

```
ocnGetAdjustedPath(
    t_libName
    t_cellName
    t_viewName
    t_netName)
=> t_adjustedPath / nil
```

# **Description**

Reduces the given hierarchical net path to the shortest hierarchical name that is equivalent to this net.

## **Arguments**

t_libName	Library name of the top cellview of the design.
t_cellName	Cell name of the top cellview of the design.
t_viewName	View name of the top cellview of the design.
t_netName	A single concatenated string for the instance hierarchy with "/" as the hierarchy separator in the string.

### Value Returned

t_adjustedPath	The reduced net name. If the net is local to this cell view only, the reduced net name is the same as the provided net name.
nil	Returns nil if there is a problem returning the adjusted path.

# **Example**

```
ocnGetAdjustedPath( "mylib" "test" "schematic" "I7/I3/gnd")
=> "/gnd"
```

The return value is "/gnd" because the gnd net is connected from this point up to the top level of hierarchy.

### Simulation Commands

### ocnGetInstancesModelName

# **Description**

This function returns the model name used by the instance in opened simulation results.

### **Arguments**

1\_instance Name of the instance in the simulation result or the schematic.

### Value Returned

1\_instancenilReturns nil if no result is open.

### **Examples**

```
ocnGetInstancesModelName()
=> (("/I8/Q4" "trpnp")
     ("/I8/Q3" "trpnp")
     ("/I8/Q2" "trpnp")
     ("/I8/Q1" "trnpn")
     ("/I8/Q0" "trnpn")
     ("/I8/C0" "capacitor")
     ("/I2" "isource")
     ("/I8/M1" "trpmos")
     ("/I8/M3" "trpmos")
     ("/I8/M2" "trnmos")
     ("/I8/M5" "trnmos")
     ("/R1" "resistor")
     ("/R0" "resistor")
     ("/I8/R0" "resistor")
     ("/V2" "vsource")
     ("/I1/V2" "vsource")
```

### Simulation Commands

```
("/I1/V0" "vsource")
)
ocnGetInstancesModelName("/R1")
=> ("/R1" "resistor")
ocnGetInstancesModelName(list("/R1" "/I8/Q1"))
=> (("/R1" "resistor") ("/I8/Q1" "trnpn") )
```

#### Simulation Commands

#### off

```
off(
    s_command
    [ g_commandArg1 ] [ g_commandArg2 ]
    )
    => t / nil
```

### **Description**

Turns off the specified information.

This command is currently available only for the analysis and restore commands. The first argument specifies the command whose information you want to turn off. If you include only this first argument, all the information for the command is turned off. If you supply subsequent arguments, only those particular pieces of information are turned off as opposed to turning off all the information for that command. The information is not deleted and can be used again.

### **Arguments**

s_command	Command that was initially used to add the items that are now being turned off.	
	Valid value: restore	
$g\_commandArg1$	Argument corresponding to the specified command.	
g_commandArg2	Additional argument corresponding to the specified command.	

#### Value Returned

t	Returns t if the information is turned off.
nil	Returns nil and prints an error message if there are
	problems turning off the information.

### **Example**

```
off( 'restore )
=> t
```

Turns off the restore command.

### Simulation Commands

```
off( restore 'tran )
=> t
```

Turns off the transient restore command.

#### Simulation Commands

# option

```
option(
    [?categ s_categ]
    s_option1 g_value1 [ s_option2 g_value2 ] ...)
    => undefined / nil
```

### **Description**

Specifies the values for built-in simulator options. You can specify values for as many options as you want.

### **Arguments**

s_categ	Type of simulator to be used.
	Valid values: analog if the options are for an analog simulator, digital for a digital simulator, or mixed for a mixed-signal simulator
	Default value: analog
s_option1	Name of the simulator option.
g_value1	Value for the option.
s_option2	Name of an additional simulator option.
g_value2	Value for the option.

#### **Value Returned**

undefined	The return value for this command/function is undefined.
nil	Returns nil and prints an error message if there are
	problems setting the option.

### **Example**

```
option( 'abstol 1e-10 )

Sets the abstol option to 1e-10.

option( 'delmax 50n )

Sets the delmax option to 50n.
```

#### Simulation Commands

option()

Returns the category list for simulation options, including analog, digital, and mixed.

option(?categ 'analog)

Returns all the simulator options for the analog simulator currently set. For example, if the set simulator is spectre, it returns the valid simulator options for spectre.

#### Simulation Commands

#### restore

```
restore(
    s_analysisType
    t_filename
)
    => undefined / nil
```

### **Description**

Tells the simulator to restore the state previously saved to a file with a store command.

This command is not available for the Spectre® circuit simulator, with which you can use the store/restore options: readns, readforce, write, or writefinal.

**Note:** Restore is available for the cdsSpice and hspiceS simulators.

### **Arguments**

s_analysisType	Type of the analysis.		
	Valid values: dc or tran		
t filename	Name of the file containing the saved state.		

#### Value Returned

undefined	The return value for this command/function is undefined.
nil	Returns nil and prints an error message if there are problems restoring the information.

### Example

```
restore( 'dc "./storeFile" )
=> ./storeFile
```

Initializes the simulator to the state saved in the storeFile file.

```
restore( 'tran "./tranStoreFile" )
=> ./tranStoreFile
```

Initializes the simulator to the state of a transient analysis saved in the transtoreFile file.

#### Simulation Commands

### resultsDir

```
resultsDir(
    t_dirName
)
    => undefined / nil
```

### **Description**

Specifies the directory where the PSF files (results) are stored.

If you do not specify a directory with this command, the PSF files are placed in .../psf to the netlist directory.

**Note:** The directory you specify with resultsDir is also where the *simulator*.out file is created.

**Note:** Some simulators are designed to *always* put their results in a specific location. For these simulators, resultsDir has no effect. You might use this command when you want to run several simulations using the same design and want to store each set of results in a different location. If this command is not used, the results of an analysis are overwritten with each simulation run.

### **Arguments**

t dirName Directory w	where the PSF files are to be stored.
-----------------------	---------------------------------------

#### Value Returned

undefined	The return value for this command/function is undefined.
nil	Returns nil and prints an error message if there is a problem with that directory.

### **Example**

```
resultsDir("~/simulation/ckt/spectre/schematic/psf") =>
    "~/simulation/ckt/spectre/schematic/psf"
```

Specifies the psf directory as the directory in which to store the PSF files.

```
resultsDir() => "~/simulation/ckt/spectre/schematic/psf"
```

Returns the results directory.

#### Simulation Commands

#### run

```
run (
     [ ?jobName t_jobName ]
     [ ?drmsCmd t_drmsCmd ])
    => s jobName / nil
run(
     [ analysisList ]
     [ ?jobName t jobName ]
     [ ?host t hostName ]
     [ ?queue t queueName ]
     [ ?startTime t startTime ]
     [ ?termTime t termTime ]
     [ ?dependentOn t dependentOn ]
     [ ?mail t mailingList ]
     [ ?block s block ]
     [ ?notify s notifyFlag ]
     [ ?lsfResourceStr s lsfResourceStr ]
    => s_jobName / nil
run (
    => t dirName / nil
run(
    s_analysisType1 - s_analysisTypeN
    => t dirName / nil
```

### Description

Starts the simulation or specifies a time after which an analysis should start. If distributed processing is not available on the system or is not enabled, the arguments specific to distributed processing (see *Arguments* section below for list of arguments specific to distributed processing) are ignored and the simulation runs locally. If distributed processing is available and is enabled, the environment default values are used if not specified in the run command arguments. The environmental default values are stored in the .cdsenv file.

Do not use the run command to start the parametric analysis. Instead, use the command that is specific to the analysis.

To start	Use this command
parametric analyses	<u>paramRun</u>

### **OCEAN Reference** Simulation Commands

# **Arguments**

analysisListList of analyses to be run with the run command. $s_analysisType1$ Name of a prespecified analysis to be simulated. $s_analysisTypeN$ Name of another prespecified analysis to be simulated.

The following arguments apply only when the distributed processing mode is enabled:

t_jobName	If the name given is not unique, an integer is appended to create a unique job name.
t_hostName	Name of the host on which to run the analysis. If no host is specified, the system assigns the job to an available host.
t_queueName	Name of the queue. If no queue is defined, the analysis is placed in the default queue.
t_startTime	Desired start time for the job. If dependencies are specified, the job does not start until all dependencies are satisfied.
t_termTime	Termination time for job. If the job has not completed by the specified termination time, the job is aborted.
t_dependentOn	List of jobs on which the specified job is dependent. The job is not started until dependent jobs are completed.
$t_{\tt mailingList}$	List of users to be notified when the analysis is complete.
$s\_block$	When $s\_block$ is not set to nil, the OCEAN script halts until the job is complete.
	Default value: nil
s_notifyFlag	When not set to $\mathtt{nil}$ , the job completion message is echoed to the OCEAN interactive window.
	Default value: t
s_lsfResourceStr	An LSF Resource Requirement string to submit a job. It is effective only in the LSF mode.
sgeHardResourceStr	Requirements for hardware resources for the job to be run in the SGE mode.

sgePriority

sgeSoftResourceStr

the SGE mode.

Requirements for software resources for the job to be run in

Priority for the job being submitted in the SGE mode.

#### Simulation Commands

sgeNoOfProcessors Number of processors to be used in the SGE mode.

sgeParallelEnvName Name of the parallel environment to be used in the SGE

mode.

t drmsCmd A DRMS (Distributed Resource Management System)

command, such as a bsub command for LSF or a qsub command for SGE (Sun Grid Engine) used to submit a job. When this argument is used, all other arguments, except ?jobName will be ignored. Moreover, it will not be possible to call the OCEAN function wait on the jobs submitted using

this argument.

To know more about the command option, refer to the section Submitting a Job in the chapter Using the Distributed Processing Option in the Analog Design Environment of the Virtuoso Analog Distributed Processing OptionUser Guide.

#### Value Returned

s	jobName	Returns the jo	ob name of the j	job submitted.	The job name is
---	---------	----------------	------------------	----------------	-----------------

based on the jobName argument. If the job name submitted is not unique, a unique identifier is appended to the job name. This value is returned for nonblocking distributed

mode.

t dirName Returns the name of the directory in which the results are

stored. This value is returned for local and blocking

distributed modes.

nil Returns nil and prints an error message if there is an error

in the simulation. In this case, look at the

yourSimulator.out file for more information. (This file is

typically located in the psf directory.)

### **Example**

```
run(?jobName "job1" ?drmsCmd "bsub -q lnx32")
=> s jobName/nil
```

where 1nx32 is the name of the queue to which the job is submitted.

run() => t

Starts the simulation.

#### Simulation Commands

```
run('tran, 'ac)
```

Runs only the tran and ac analyses.

```
run('dc)
```

Runs only the dc analysis.

```
run( ?jobName ?block "nil")
=> 'reconFilter
```

Returns a job name of reconfilter for the specified job and runs that job if distributed processing is enabled. The job is submitted nonblocking. The job name is returned.

```
run( ?queue "fast" )
```

Submits the current design and enabled analyses as a job on the fast queue, assuming that distributed processing is available and enabled.

```
run( ?jobName "job1" ?queue "fast" ?host "menaka" ?startTime "22:59"
?termTime "23:25" ?mail "preampGroup")
```

Submits the current design and enabled analyses as a jobName job1 on the fast queue host menaka with the job start time as 22:59 and termination time as 23:25. A mail will be sent to preampGroup after the job ends.

```
run( ?jobName "job1" ?queue "fast" ?host "menaka" ?lsfResourceStr "mem>500")
```

Submits the current design and enabled analyses as a jobName job1 on the fast queue host menaka, if the host has at least 500 MB of RAM memory.

#### Simulation Commands

#### save

```
save(
    [?categ s_categ ] s_saveType
    [ t_saveName1 ] ... [ t_saveNameN ]
    )
    => undefined / nil
```

# **Description**

Specifies the outputs to be saved and printed during simulation.

When specifying particular outputs with <code>saveName</code>, you can include as many outputs as you want. If you want to turn off the default of <code>save</code>, <code>'allv</code>, use the <code>delete('save')</code> command.

#### Simulation Commands

#### **Arguments**

Valid values: analog, digital

Default value: analog

Note: digital is not available.

*s saveType* Type of outputs to be saved.

Valid values:

v: Specifies that a list of subsequent net names be saved.

i : Specifies that a list of subsequent currents be saved.

all: Specifies that all nets and all currents are to be saved.

ally: Specifies that all voltages are to be saved.

alli: Specifies that all currents are to be saved.

Default value: allv

t\_saveName1 Name of the net, device, or other object.

t saveNameN Name of another net, device, or object.

#### Value Returned

undefined The return value for this command/function is undefined.

nil Returns nil and prints an error message if there is a

problem saving the outputs.

### **Example**

```
save( 'v "net34" "net45" )
```

Saves the outputs for net34 and net45.

```
save( 'i "R1" "/Q1/b" )
```

Saves the currents for R1 and Q1/b.

```
save('all)
```

Saves all the nets and currents.

```
save( 'i "q1:b" "r1:p" "mn1:d" )
```

#### Simulation Commands

For the spectre simulator, saves the current through the specified devices.

```
save( ?categ 'analog 'v "/vin" "/vout" )
```

Saves the output for vin and vout.

```
save( 'i "i(q1,b)" "i(r1)" "i(mn1,d)" )
```

For the Cadence-SPICE circuit simulator, saves the current through the same devices.

#### Simulation Commands

# saveOpPoint

```
saveOpPoint(
    t_instName
    [?operatingPoints l_operatingPoints]
)
=> t / nil
```

### **Description**

Specifies the operating point parameters to be saved for a given instance.

### **Arguments**

t\_instName Name of the instance for which the parameters are to be saved.

?operatingPoints 1 operatingPoints

List of the operating point parameters.

#### Values Returned

t Returns t when the function runs successfully.

nil Returns nil if there is an error.

### **Example**

```
saveOpPoint( "/I8/Q3" ?operatingPoints "vbe isub betaac gm re" )
```

This example saves the operating point parameters, vbe isub betaac gm re, for the instance /18/Q3.

#### Simulation Commands

# saveOption

```
saveOption(
     [ s_option1 g_optionValue1 ]...[ s_optionN g_optionValueN ]
    )
     => undefined / nil
```

### **Description**

Specifies save options to be used by the simulator.

You can include as many save options as you want. To include a save option, replace  $s\_option1$  with the name of the desired save option and include another argument to specify the value for the option.

When you use the saveOption command without specifying any arguments, the command returns a list of option and value pairs.

Save options vary, depending on the simulator and interface that you are using. If you are using the Spectre® circuit simulator, for example, you can type the following at an OCEAN prompt to see which options you can set with the saveOption command:

```
simulator('spectre)
ocnHelp('saveOption)
```

See the *Virtuoso Spectre Circuit Simulator User Guide* for more information on these options.

**Note:** The saveOption command does not work with socket simulators. If you are using a socket simulator, you must instead specify save options with the save command described in <u>"save"</u> on page 155.

#### Simulation Commands

### **Arguments**

s option1 Save option. The save options that are available depend on

which simulator you use. (See the documentation for your

simulator.)

g\_optionValue1 Value for the save option.

s optionN Any subsequent save option. The save options that are

available depend on which simulator you use. (See the

documentation for your simulator.)

g optionValueN Value for the save option.

#### Value Returned

undefined The return value for this command/function is undefined.

nil Returns nil if there are problems specifying options.

### **Example**

saveOption( 'save "lvl" 'nestlvl 10 'currents "selected" 'useprobes "yes"
'subcktprobelvl 2 ?saveahdlvars "all")

#### Simulation Commands

#### simulator

```
simulator(
    s_simulator
)
=> s simulator / nil
```

### **Description**

Starts an OCEAN session and sets the simulator name for that session. The previous session (if any) is closed and all session information is cleared.

#### **Arguments**

s simulator

Name of the simulator.

#### **Value Returned**

s\_simulator

Returns the name of the simulator.

nil

Returns nil and prints an error message if the simulator is not registered with the Virtuoso® Analog Design Environment through OASIS. If the simulator is not registered, the simulator from the preceding session is retained.

### **Example**

```
simulator( 'spectre )
=> spectre
```

Specifies that the Spectre® circuit simulator be used for the session.

```
simulator()
=> spectre
```

Returns the simulator that you set for the session. If a simulator was not specified, it returns nil.

#### Simulation Commands

#### solver

```
solver(
    s_solver
)
=> s solver / nil
```

### **Description**

Sets a solver for a given AMS OCEAN session. The valid values for solver are Spectre and UltraSim. You select Spectre if you want to use an accurate AMS-Spectre analog engine. You select UltraSim if you want to use the AMS-Ultrasim or FastSPICE(UltraSim) solver for a given AMS simulation.

**Note:** This command is applicable only when ams is the simulator.

#### **Arguments**

s solver

Name of the solver.

#### Value Returned

s solver

Returns the name of the solver.

nil

Returns nil and prints an error message if the specified solver is not registered with the Virtuoso® Analog Design Environment through OASIS. If the solver is not registered,

the solver from the preceding session is retained.

#### **Example**

```
solver( 'spectre )
=> spectre
```

Specifies AMS-Spectre as the solver to be used for the current AMS session.

```
solver( 'ultraSim )
=> ultraSim
```

Specifies AMS-UltraSim (UltraSim FastSPICE) as the solver to be used for the current AMS session.

#### Simulation Commands

### stimulusFile

```
stimulusFile(
    t_fileName [ t_fileName2 ... t_fileNameN ]
    [ ?xlate g_xlate ]
    )
    => 1 fileNames / nil
```

### **Description**

Specifies stimulus files to be used by the simulator.

When the  $g_xlate$  variable is set to t, the schematic net expressions [#net] and instance name expression [\$instance] in the stimulus file are mapped into simulator names before including. When a netlist is specified as the design, this option must be set to nil.

**Note:** This command does not work with socket simulators.

### **Arguments**

t_fileName	The name of the stimulus file to be included.
t_fileName2t_fileNa	nmeN
	The names of the additional stimulus files to be included.
g_xlate	If set to $t$ , net and instance expressions are translated to simulator names. The default value of the $g_xlate$ variable is $t$ .

#### **Value Returned**

l_fileNames	A list of the stimulus file names is the output if the command is successful
nil	Otherwise nil is returned

### **Example**

```
stimulusFile( "tran.stimulus rf.stimulus" ?xlate nil)
=> ("tran.stimulus rf.stimulus")
```

Includes tran.stimulus and rf.stimulus in the simulator input file. No net and instance expressions are translated.

### Simulation Commands

```
stimulusFile()
=> ("tran.stimulus" "rf.stimulus")
```

Returns the stimulusFile, if one was set earlier. Otherwise, it returns nil.

#### Simulation Commands

#### store

```
store(
    s_analysisType
    t_filename
)
    => t filename / nil
```

### **Description**

Requests that the simulator store its node voltages to a file.

You can restore this file in a subsequent simulation to help with convergence or to specify a certain starting point. This command is not available for the Spectre® circuit simulator, with which you can use the store/restore options: readns, readforce, write, or writefinal.

**Note:** store is available for the cdsSpice and hspiceS simulators.

### **Arguments**

s_analysisType	Type of the analysis.
	Valid values: dc or tran
t_filename	Name of the file in which to store the simulator's node voltages.

#### Value Returned

t_filename	Returns the filename.
nil	Returns nil and prints an error message if there are
	problems storing the information to a file.

### **Example**

```
store( 'dc "./storeFile" )
=> ./storefile
```

Stores the simulator's node voltages in a file named storeFile in the current directory.

```
store( 'tran "./tranStoreFile" )
=> ./transtorefile
```

Stores the node voltages for a transient analysis in a file named transforeFile in the netlist (design) directory unless a full path is specified.

#### Simulation Commands

# temp

```
temp(
    f_tempValue
)
=> s tempValue / nil
```

### **Description**

Specifies the circuit temperature.

# **Arguments**

f\_tempValue Temperature for the circuit.

### Value Returned

s_tempValue	Returns the temperature specified.
nil	Returns nil and prints an error message if there are
	problems setting the temperature.

# **Example**

```
temp( 125 )
=> ?125?
atof(temp( 125 ))
=> 125.0
```

Sets the circuit temperature to 125.

```
temp() => 125
```

Gets the value you had set for the circuit temperature. If you have not set a value for the temperature, it returns the default value.

### Simulation Commands

#### tran

```
tran(
    g_fromValue
    g_toValue
    g_byValue
)
    => g_byValue / nil

tran(
    g_toValue
)
    => undefined / nil
```

### **Description**

Specifies a transient analysis with limited options. If other analysis options are needed, use the <u>analysis</u> command.

To know more about this analysis, see the simulator-specific user guide.

**Note:** The second instance of the tran command is valid only with the spectre simulator.

# **Arguments**

g_fromValue	Starting time for the analysis
g_toValue	Ending time
g byValue	Increment at which to step through the analysis

#### Value Returned

undefined	The return value for this command/function is undefined.
nil	Returns nil and prints an error message if the analysis is not specified.

### **Example**

```
tran( 1u) => "1e-06"
```

Specifies a transient analysis to 1u for the Spectre® circuit simulator

```
tran( 0 1u 1n ) => "1e-09"
```

# **OCEAN Reference** Simulation Commands

Specifies a transient analysis from 0 to 1u by increments of 1n.	

### **OCEAN Reference** Simulation Commands

### vcdFile

```
vcdFile(
    t_vcdFileName
)
=> t vcdFileName / nil
```

### **Description**

Sets a VCD file for a given AMS or UltraSim OCEAN session. You also need to specify a VCD info file while using this command. You can specify only one VCD file for a session. You may use ocnDisplay('vcdFile) to view the currently active VCD file.

**Note:** This command is applicable for AMS and UltraSim simulators. For AMS, it works only when UltraSim is the solver.

### **Arguments**

t vcdFileName	The name of the VCD file to be used for session.
0 . 0 0.1 = 1 0 1.0 0	

#### Value Returned

$t\_vcdFileName$	The VCD file name is the output if the command is successful.
nil	Otherwise, nil is returned.

### **Example**

vcdFile("/tmp/vcdFile.dat")

```
=> "/tmp/vcdFile.dat"

Specifies /tmp/vcdFile.dat as the VCD file to be used for current AMS-UltraSim OCEAN session.
```

#### Simulation Commands

### vcdInfoFile

```
vcdInfoFile(
    t_vcdInfoFileName
)
=> t vcdInfoFileName / nil
```

### **Description**

Sets a VCD info file for a given AMS or UltraSim OCEAN session when you have set UltraSim as the solver. You also need to specify a VCD file while using this command. You can specify only one VCD info file for a session. You may use ocnDisplay('vcdInfoFile) to view the currently active VCD info file.

**Note:** This command is applicable for AMS and UltraSim simulators. For AMS, it works only when UltraSim is the solver.

### **Arguments**

t vcdInfoFileName The name of the VCD info file to be included.

#### Value Returned

t\_vcdInfoFileName The VCD info file name is the output if the command is successful.nil Otherwise, nil is returned.

#### **Example**

```
vcdInfoFile("/tmp/vcdInfoFile.dat")
=> "/tmp/vcdInfoFile.dat"
```

Specifies / tmp/vcdInfoFile.dat as the VCD file to be used for current AMS-UltraSim OCEAN session.

#### Simulation Commands

### vecFile

```
vecFile(
    t_vecFile [ t_vecFile1 ... t_vecFileN ]
)
    => t vecFile(s) / nil
```

### **Description**

Sets the vector files to be used in an AMS or UltraSim OCEAN session. You use the vecFile command to specify a list of vector files which go to control file. You may use ocnDisplay('vecFile) to view the currently active vector files in an OCEAN session.

**Note:** This command is applicable for AMS and UltraSim simulators. For AMS, it works only when UltraSim is the solver.

#### **Arguments**

 $t\_vecFile$  The name of the vector file to be included.  $t\_vecFile1...t\_vecFileN$ 

The names of the additional vector files to be included.

#### Value Returned

t\_vecFile

The names of the vector file(s) are listed if the command is successful.

nil

Otherwise, nil is returned.

### **Example**

```
vecFile("/tmp/vec.dat" "/tmp/vec2.dat")
=> ("/tmp/vec1.dat" "/tmp/vec2.dat")
Specifies /tmp/vec.dat and /tmp/vec.dat2 as the vector files to be used for the current AMS-UltraSim OCEAN session.
```

#### Simulation Commands

#### **hlcheck**

### **Description**

Sets or gets the value of the hlcheck option used in the vec\_include statement in a netlist. You may use the ocnDisplay('hlcheck) command to view the current value of hlcheck in an OCEAN session associated with vector files.

**Note:** This command is applicable only when one or more vector files are specified in a given 'spectre' OCEAN session.

### **Arguments**

t_value	Value to be set for the hlcheck option. Possible values
	include "off", "0", and "1". The value "off" disables
	the hlcheck option in the vec_include statement.

#### Value Returned

t	Returns t if the hlcheck option is set with the value supplied as argument
nil	Otherwise, returns nil and an error message is displayed

#### Example

```
hlcheck( "1" )
=> t
Sets the value of the hlcheck option as 1 in the vec_include statement
hlcheck()
=> "1"
Returns the value of the hlcheck option
```

#### Simulation Commands

# ocnAmsSetOSSNetlister

```
ocnAmsSetOSSNetlister(
    )
    => t / nil
```

# **Description**

Sets the netlister mode to OSS-based for a given ams OCEAN session.

# **Arguments**

None

### **Value Returned**

t Returns t if successful Returns nil otherwise

### **Example**

```
ocnAmsSetOSSNetlister()
t
```

Sets the netlister mode to OSS-based instead of Cellview-based for the current ams simulator session.

#### Simulation Commands

### ocnAmsSetUnlNetlister

```
ocnAmsSetUnlNetlister(
    )
    => t / nil
```

# **Description**

Sets the netlister mode to AMS Unified Netlister for a given ams OCEAN session.

# **Arguments**

None.

#### **Value Returned**

t Returns t if successful nil Returns nil otherwise

### **Example**

```
ocnAmsSetUnlNetlister()
t
```

Sets the netlister mode to AMS Unified Netlister instead of Cellview-based for the current ams simulator session.

# **Data Access Commands**

The data access commands let you open results and select different types of results to analyze. You can get the names and values of signals and components in the selected results, and you can print different types of reports.

In this chapter, you can find information on the following data access commands

dataTypes
deleteSubckt
displaySubckt
getData

<u>i</u>

getResult

<u>openResults</u>

<u>outputParams</u>

<u>outputs</u>

<u>phaseNoise</u>

<u>pv</u>

<u>resultParam</u>

<u>results</u>

<u>selectResult</u>

<u>sp</u>

### **Data Access Commands**

<u>sprobeData</u>
<u>sweepNames</u>
sweepValues
sweepVarValues
<u>v</u>
<u>vswr</u>
<u>zm</u>

zref

#### **Data Access Commands**

# dataTypes

```
dataTypes(
    )
    => 1_dataTypes / nil
```

### **Description**

Returns the list of data types that are used in an analysis previously specified with selectResult.

# **Arguments**

None.

#### Value Returned

$1\_dataTypes$	Returns the list of data types.
nil	Returns nil and an error message if the list of datatypes
	cannot be returned.

### **Example**

```
selectResult( 'dcOpInfo )
dataTypes() => ("bjt" "capacitor" "isource" "mos2" "resistor" "vsource")
```

Returns the data types used in the selected file, in this case, dcOpInfo.

#### Example 2:

```
selectResult( 'model )
dataTypes() => ("bjt" "mos2")
```

Returns the data types used in the selected file, in this case, model.

#### **Data Access Commands**

### deleteSubckt

```
deleteSubckt(
    t_name
)
=> t / nil
```

### **Description**

Deletes the specified subcircuit instance saved using the <u>saveSubckt</u> command.

# **Arguments**

*t\_name* The name of the subcircuit instance.

#### Value Returned

t Returns t if the selected subcircuit instances is deleted.

nil Returns nil if the name of the specified instance is not correct.

### **Examples**

The following example deletes the subcircuit instance 10.

```
deleteSubckt("/I0")
=> t
```

#### **Data Access Commands**

# displaySubckt

```
displaySubckt(
    t_args
    t_outPort
)
    => t / nil
```

### **Description**

Prints the subcircuit information to the output file.

### **Arguments**

t args The value of this argument should always be nil.
---

 $t\_outPort$  The name of the file to save the subcircuit information. If you do

not specify the location with the filename, the file is saved in the

current working directory.

#### Value Returned

t Returns t if the subcircuit information is printed in the specified

output file.

nil Returns nil if the name of the output file is not specified, or an

error occurred.

#### **Examples**

The following example prints the subcircuit information in the subckts.txt file:

```
fptr = outfile("/home/krajiv/subckts.txt")
=> port:"/home/krajiv/subckts.txt"
displaySubckt(nil fptr)
=> t
close(fptr)
=> t
```

#### **Data Access Commands**

# getData

```
getData(
    t_name
    [?result s_resultName [?resultsDir t_resultsDir ]])
    => x_number / o_waveform / nil
```

# **Description**

Returns the number or waveform for the signal name specified.

The type of value returned depends on how the command is used.

#### **Data Access Commands**

### **Arguments**

t	name	Name of the signal.

s resultName Results from an analysis. When specified, this argument

will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult

command.

t\_resultsDir Directory containing the PSF files (results). If you supply

this argument, you must also supply the resultName argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the

openResults command.

#### Value Returned

x number Returns an integer simulation result.

o waveform Returns a waveform object. A waveform object

represents simulation results that can be displayed as a series of points on a grid. (A waveform object identifier

looks like this: srrWave: XXXXX.)

nil Returns nil and an error message if the value cannot be

returned.

### Example

```
getData( "/net6" ) => srrWave:25178234
```

Returns the number or waveform for net6. In this example, the return value is equivalent to v("/net6").

```
getData( "/V1" ?result 'ac )
=> srrWave:96879364
```

Returns the number or waveform for V1. In this example, the return value is equivalent to: i( "/V1" ?result 'ac ).

```
selectResult( 'tran ) =>
ocnPrint( getData( "net1" ) ) =>
```

#### **Data Access Commands**

The <code>getData("net1")</code> command passes a waveform to the <code>ocnPrint</code> command. The <code>ocnPrint</code> command then prints the data for the waveform. In this example, the return value is equivalent to:

```
(v( "net1" )).
ocnPrint( getData( "net1" ?result 'tran ?resultsDir "./simulation/testcell/
spectre/schematic/psf")
```

Returns a signal on net1 for the tran result stored in the path "./simulation/testcell/spectre/schematic/psf".

**Note:** To identify the data type of the value returned by the getData command, you can use the type SKILL function. For scalar values, the type function returns the name of data type. For example, integer or flonum. For waveforms, it returns other.

```
x=getData("/net10")
type(x)
```

The example given above returns other.

```
x=ymax(VT("/net10"))
type(x)
```

This will return flonum.

#### **Data Access Commands**

# getResult

```
getResult (
    [ ?result s_resultName [ ?resultsDir t_resultsDir ] ]
    )
    => o results / nil
```

## **Description**

Gets the data object for a specified analysis without overriding the status of any previously executed selectResult() or openResults() commands.

Returns the data object for a particular analysis similar to the selectResult() function does. Unlike the selectResult() function, all subsequent data access commands will not internally use this information.

## **Arguments**

$s\_resultName$	Results from an analysis. When specified, this argument
	will only be used internally and will not alter the current

result which was set by the selectResult command. The default is the current result selected with the selectResult

command.

t resultsDir Directory containing the PSF files (results). If you supply

this argument, you must also supply the resultName argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the

openResults command.

#### Value Returned

o results Returns the object representing the selected results.

nil Returns nil and an error message if there are problems

accessing the analysis.

## **Example**

```
getResult( ?result 'tran )
```

#### **Data Access Commands**

# i

```
i(
    t_component
    [?result s_resultName [?resultsDir t_resultsDir]]
)
=> o waveform / nil
```

# **Description**

Returns the current through the specified component.

## **Arguments**

t_component	Name of the component.
_	Results from an analysis. When specified, this argument will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult command.

Directory containing the PSF files (results). If you supply this argument, you must also supply the resultName argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the openResults command.

### **Value Returned**

t resultsDir

o_waveform	Returns a waveform object. A waveform object represents simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.).
nil	Returns an error message and nil if there is a problem.

## Example

```
selectResult( 'tran )
i( "/R1" )
```

### **Data Access Commands**

Returns the current through the R1 component.

```
ocnPrint( i( "/R1" ) )
```

Prints the current through the R1 component.

```
ocnPrint( i( "/R1" ?result 'dc ) )
```

Prints the current through the R1 component with respect to the dc swept component.

```
ocnPrint( i( "/R1" ?resultsDir "./test2/psf" ?result 'dc ) )
```

Prints the current through the R1 component with respect to dc for the results from a different run (stored in test2/psf).

#### **Data Access Commands**

# ocnHelp

```
ocnHelp(
    [ ?output t_filename | p_port ]
    [ s_command ]
    )
    => t / nil
```

## **Description**

Provides online help for the specified command.

If no command is specified, provides information about how to use help and provides the different categories of information contained in the help library. If you provide a filename as the <code>?output</code> argument, the <code>ocnHelp</code> command opens the file and writes the information to it. If you provide a port (the return value of the SKILL <code>outfile</code> command), the <code>ocnHelp</code> command appends the information to the file that is represented by the port. If you do not specify <code>?output</code>, the output goes to standard out (stdout).

#### **Data Access Commands**

### **Arguments**

t_fi	lename	File	in w	hich	to	write t	the	inform	ation.	The	ocnHelp	Ō
------	--------	------	------	------	----	---------	-----	--------	--------	-----	---------	---

command opens the file, writes to the file, and closes the

file. If you specify the filename without a path, the ocnHelp command creates the file in the directory pointed to by your SKILL Path. To find out what your SKILL path is, type getSkillPath() at the OCEAN

prompt.

p port Port (previously opened with outfile) through which to

append the information to a file. You are responsible for closing the port. See the <u>outfile</u> command for more

information.

s command Command for which you want help.

### Value Returned

t Displays the online help and returns t.

nil Returns nil and an error message if help cannot be

displayed.

### **Example**

```
ocnHelp()
=> t
```

Displays information about using online help.

```
ocnHelp( 'analysis )
=> t
```

Displays help for the analysis command.

```
ocnHelp( ?output "helpInfo" )
=> t
```

Writes information about using online help to a file named helpInfo.

```
simulator('spectre)
ocnHelp('envOptions)
```

Displays a list of environment options that can be set for a simulator. First, set the simulator and then run the ocnHelp command.

### **Data Access Commands**

### ocnResetResults

```
ocnResetResults(
    )
    => t
```

## **Description**

Unsets the results opened by the <code>openResults</code> command. Use this command to return to the state that existed prior to using the <code>openResults</code> command.

# **Arguments**

None.

#### Value Returned

t

Resets the results and returns t.

## **Example**

```
getResult(?result 'tran) Returns nil when no results have been opened.
openResults("./psf") Makes getResult return valid object.
ocnResetResults() Resets the results opened by openResults and makes getResult return nil.
```

#### **Data Access Commands**

# openResults

```
openResults(
    s_jobName | t_dirName
    [ g_enableCalcExpressions ]
    )
    => t_dirName / nil
```

## **Description**

Opens simulation results stored in PSF files or opens the results from a specified job, depending on which parameter is called.

When openResults passes a symbol, it interprets the value as a job name and opens the results for the specified job.  $s_{jobName}$  is a job name and is defined when a run command is issued.

When openResults passes a text string, it opens simulation results stored in PSF files in the specified directory. The results must have been created by a previous simulation run through OCEAN or the Virtuoso® Analog Design Environment. The directory must contain a file called logFile and might contain a file called runObjFile. When you perform tasks in the design environment, the runObjFile is created. Otherwise, only logFile is created.

If you want to find out which results are currently open, you can use openResults with no argument. The directory for the results that are currently open is returned.

**Note:** If you run a successful simulation with distributed processing disabled, the results are automatically opened for you. Also, a job name is generated by every analysis, even if distributed processing is not enabled.

#### **Data Access Commands**

### **Arguments**

 $s\_jobName$  The name of a distributed process job.  $s\_jobName$  is a job name

and is defined when a run command is issued.

t dirName The directory containing the PSF files.

g enableCalcExpressions

An optional argument, which when set to t, allows the evaluation of Calculator expressions. For this argument to work, the directory mentioned in  $t\_dirName$  must be a psf directory and must contain

runObjFile.

The default value for this argument is t.

#### Value Returned

t dirName The directory containing the PSF files.

nil Returns nil and an error message if there are

problems opening the results.

## **Example**

```
openResults( "./simulation/opamp/spectre/schematic/psf" )
=> "./simulation/opamp/spectre/schematic/psf"
```

Opens the results in the psf directory within the specified path.

```
openResults( "./psf" )
=> psf
```

Opens the results in the psf directory in the current working directory.

```
openResults( "./psf" t )
=> psf
```

Opens the results in the psf directory in the current working directory. It also allows the evaluation of the Calculator expression.

### **Data Access Commands**

# **outputParams**

```
outputParams(
    t_compType
    [?result s_resultName [?resultsDir t_resultsDir]]
)
=> 1 outputParams / nil
```

# **Description**

Returns the list of output parameters for the specified component.

You can use the <u>dataTypes</u> command to get the list of components for a particular set of results.

**Note:** You can use any of the parameters in outputParams as the second argument to the <u>pv</u> command.

#### **Data Access Commands**

### **Arguments**

t\_compType Name of a component.

 $?result s\_resultName$  Results from an analysis. When specified, this argument

will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult

command.

?resultsDir t\_resultsDir

Directory containing the PSF files (results). If you supply this argument, you must also supply the resultName argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the openResults command.

#### Value Returned

1 outputParams Returns the list of parameters.

nil Returns nil and an error message if there are no

associated parameters or if the specified component

(compType) does not exist.

## **Example**

```
selectResult( 'dcOpInfo )
dataTypes() => ("bjt" "capacitor" "isource" "mos2" "resistor" "vsource")
outputParams( "bjt" )
```

Selects the dcOpInfo results, returns the list of components for these results, and returns the list of output parameters for the bjt component.

```
outputParams("bjt" ?result 'dcOpInfo ?resultsDir "/VADE615/simulation/ampTest/
spectre/schematic/psf")
```

Returns a list of output parameters for the bit component for dcOpInfo (dc analysis with save dc operating point) results stored at the location ./psf.

### **Data Access Commands**

# outputs

```
outputs(
    [ ?result s_resultName [ ?resultsDir t_resultsDir ] ]
    [ ?type t_signalType ]
    )
    => 1_outputs / nil
```

# **Description**

Returns the names of the outputs whose results are stored for an analysis. You can plot these outputs or use them in calculations.

#### **Data Access Commands**

## **Arguments**

?result  $s_resultName$ 

Results from an analysis. When specified, this argument will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult command.

? resultDir t resultsDir

Directory containing the PSF files (results). If you supply this argument, you must also supply the resultName argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the openResults command.

?type t signalType

Data type of the signal.

#### Value Returned

l\_outputs

Returns the list of outputs.

nil

Returns nil and an error message if there are problems returning the names of the stored outputs.

## **Example**

```
outputs() => ( "net13" "net16" "net18" )
```

Returns the names of the outputs for the PSF file selected with selectResult.

```
outputs ( ?type "V" )
```

Returns all the signal names that are node voltages. The dataType (signal) returns the data type of the signal.

```
outputs(?result "tran" ?resultsDir "./psf")
=> ( "net11" "net15" "net17")
```

Returns the names of the outputs for the tran results stored at the location ./psf.

### **Data Access Commands**

# phaseNoise

```
phaseNoise(
    g_harmonic S_signalResultName
    [ ?result s_noiseResultName [ ?resultsDir t_resultsDir ] ]
    )
    => o waveform / nil
```

# **Description**

Returns the phase noise waveform which is calculated using information from two PSF data files.

This command should be run on the results of the Spectre pss-pnoise analysis.

#### **Data Access Commands**

## **Arguments**

g harmonic

List of harmonic frequencies.

?result S signalResultName

Name of the result that stores the signal waveform. Use the results() command to obtain the list results.

?resultsDir s noiseResultName

Name of the result that stores the "positive output signal" and "negative output signal" noise waveforms. When specified, this argument will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult command.

t resultsDir

Directory containing the PSF files (results). If you supply this argument, you must also supply the S\_noiseResultName argument. Both the S\_signalResultName and S\_noiseResultName arguments are read from this directory. When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the openResults command.

#### Value Returned

o waveform

Waveform representing the phase noise.

nil

Returns nil if there is an error.

## Example

```
plot(phaseNoise(0 "pss-fd.pss"))
phaseNoise(1 "pss fd" ?result "pnoise" ?resultsDir "./PSF")
```

### **Data Access Commands**

# pν

```
pv(
    t_name
    t_param
    [ ?result s_resultName [ ?resultsDir t_resultsDir ] ]
    )
    => g_value / nil
```

# **Description**

Returns the value of the specified component parameter. You can use the <u>outputParams</u> command to get the list of parameters for a particular component.

#### **Data Access Commands**

### **Arguments**

*t\_name* Name of the node or component.

t param Name of the parameter.

?result s resultName

Results from an analysis. When specified, this argument will only be used internally and will not alter the current result that was set using the selectResult command. The default is the current result selected using the selectResult command.

**Note:** To get the correct value of the variables while running parametric analysis, use the designParamVals value for the resultName argument.

?resultsDir t resultsDir

Directory containing the PSF files (results). If you supply this argument, you must also supply the resultName argument. When specified, this argument will only be used internally and will not alter the current results directory that was set using the openResults command. The default is the current results directory set using the openResults command.

#### Value Returned

*q value* Returns the requested parameter value.

nil Returns nil and prints an error message.

### **Example**

```
selectResult( 'dcOpInfo )
pv( "/I0/M1" "vds" )
```

Returns the value of the vds parameter for the IO/M1 component.

```
pv( "/I0/M1" "vds" ?resultsDir "/VADE/simulation/ampTest/spectre/schematic/test2/psf" )
```

Returns the value of the vds parameter for the IO/M1 component. These values are read from the results directory, /VADE/simulation/ampTest/spectre/schematic/test2/psf.

```
pv( "/I0/M1" "vds" ?result "dcOpInfo" ?resultDir "/VADE/simulation/ampTest/
spectre/schematic/test1/psf")
```

### **Data Access Commands**

Returns the value of the vds parameter for the IO/M1 component. These values are read from the dcOpInfo results saved in the results directory, VADE/simulation/ampTest/spectre/schematic/test1/psf.

```
pv("top-level" "CAP" ?result "designParamVals")
```

Returns the value of the CAP variable for the top-level hierarchy in the design. These values are read from the default results directory.

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#### **Data Access Commands**

## resultParam

```
resultParam(
    S_propertyName
    [?result s_resultName [ ?resultsDir t_resultsDir ] ]
    )
    => L value / nil
```

## **Description**

Returns the value of a header property from the selected result data.

## **Arguments**

s\_propertyName Name of the parameter
?result s resultName

Results from an analysis. When specified, this argument will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult command.

?resultsDir t resultsDir

Results from an analysis. When specified, this argument will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult command.

#### Value Returned

Value of the parameter. The data type depends on the data type of the parameter.

Nil Returns nil and an error message if there are problems returning the results.

## **Example**

```
resultParam("positive output signal" ?result "pnoise.pss")
=> "pif"
resultParam("negative output signal" ?result "pnoise.pss")
=> "0"
```

#### **Data Access Commands**

Returns the name of the positive and negative output signals from PSS-noise analysis result. In this case, the data type of the returned value is a string.

```
resultParam("port1.r.value" ?result "sp")
=> 40.0
resultParam("port2.r.value" ?result "sp")
=> 40.0
```

Returns the reference impedance of the ports in a two-port network from the S-parameter analysis result. In this case, the data type of the returned value is a floating point number.

```
resultParam("positive output signal" ?result "pnoise.pss" ?resultsDir "./psf")
=> "0"
```

Returns the names of the positive output signals from the PSS-noise analysis results stored at the location ./psf.

#### **Data Access Commands**

### results

```
results(
    [ ?resultsDir t_resultsDir ]
)
=> 1 results / nil
```

## **Description**

Returns a list of the type of results that can be selected.

## **Arguments**

```
?resultsDir t resultsDir
```

Directory containing the PSF files (results). When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the openResults command.

#### Value Returned

l_results	Returns the list of result types.
nil	Returns $\mathtt{nil}$ and an error message if there are problems returning the results.

## **Example**

```
results()
=> ( dc tran ac )
```

Returns the list of results available.

```
results( ?resultsDir "./psf" )
```

Returns a list of results stored at the location ./psf.

### **Data Access Commands**

## saveSubckt

```
saveSubckt(
    t_name
    [ ?voltage g_voltage ]
    [ ?current g_current ]
    [ ?power g_power ]
    [ ?vDepth s_vDepth ]
    [ ?iDepth s_iDepth ]
    [ ?pwrDepth s_pwrDepth ]
    [ ?compress g_compress ]
    [ ?filterRC g_filterRC ]
    [ ?ports g_ports ]
    [ ?userOptions g_userOptions ]
    )
    => t / nil
```

# **Description**

Saves and modifies the specified subcircuit instances and signals.

## **Data Access Commands**

# **Arguments**

t_name	The name of the subcircuit instance.
?voltage <i>g_voltage</i>	Specifies whether you want to save the voltage for the subcircuit.
?current g_current	Specifies whether you want to save the current for the subcircuit.
?power g_power	Specifies whether you want to save the power signals for the subcircuit.
?vDepth s_vDepth	The hierarchy level to which you want to save the voltage signal for the subcircuit. If not specified, voltage for all the levels of hierarchy are saved.
?iDepth s_iDepth	The hierarchy level to which you want to save the current signal for the subcircuit. If not specified, current for all the levels of hierarchy are saved.
?pwrDepth s_pwrDepth	The hierarchy level to which you want to save the power signal for the subcircuit. If not specified, power signals for all the levels of hierarchy are saved.
?compress g_compress	Specifies whether you want to reduce the size of the output file. When enabled, the spectre simulator saves the data for a signal only when the value of that signal changes.
?filterRC g_filterRC	Specifies whether to filter out the nodes that are connected only to parasitic elements from the output signal list.
?ports g_ports	Specifies whether to save the output port information for the specified subcircuit.
?userOptions g_userOptions	Specify the other save options that you want to define for the signal.

## **Value Returned**

t Returns t if the subcircuit instance is saved.

nil Returns nil if the name of the specified instance is not correct.

# **Examples**

#### **Data Access Commands**

# Example 1

The following example saves the voltage for five levels and current for two levels of hierarchy for the subcircuit 10.

```
saveSubckt("/I0" ?voltage t ?current t ?vDepth "5" ?iDepth "2")
=> t
```

## Example 2

The following example saves the voltage for two levels and power signals for one level of hierarchy for the subcircuit I1.

```
saveSubckt("/I1" ?voltage t ?power t ?vDepth "2" ?pwrDepth "1")
=> t
```

## Example 3

The following example saves the voltage for two levels and power signals for one level of hierarchy for the subcircuit I1, along with the port information. The output signals are compressed.

```
saveSubckt("/I1" ?voltage t ?power t ?vDepth "2" ?pwrDepth "1" ?port t ?compress t)
=> t
```

#### **Data Access Commands**

## selectResult

```
selectResult(
    S_resultsName
    [ n_sweepValue ]
    )
    => o results / nil
```

## **Description**

Selects the results from a particular analysis whose data you want to examine.

The argument that you supply to this command is a data type representing the particular type of analysis results you want. All subsequent data access commands use the information specified with selectResult.

**Note:** Refer to the <u>results</u> command to get the list of analysis results that you can select.

## **Arguments**

$s\_resultsN$ ame	Results from an analysis.
n_sweepValue	The sweep value you wish to select for an analysis.

#### Value Returned

o_results	Returns the object representing the selected results.
nil	Returns $\mathtt{nil}$ and an error message if there are problems selecting the analysis.

## **Example**

```
selectResult( 'tran )
```

Selects the results for a transient analysis.

```
sweepValues(3.0 3.333333 3.666667 4.0 4.333333 4.666667 5.0 )
selectResult("tran" "3.333333")
```

The sweepValues command prints a list of sweep values.

The selectResult command selects a specific value for a transient analysis.

```
selectResult( 'tran )
```

### **Data Access Commands**

Selects the results for a transient analysis.

```
paramAnalysis("supply" ?start 3 ?stop 5 ?step 1.0/3)
paramRun("supply")
selectResult(( 'tran car( sweepValues() )
```

Selects the data corresponding to the first parametric run.

**Note:** selectResult('tran) would select the entire family of parametric data.

### **Data Access Commands**

# sp

```
sp(
    x_iIndex
    x_jIndex
    [?result s_resultName [ ?resultsDir t_resultsDir ] ]
    )
    => o waveform / nil
```

# **Description**

Returns S-parameters for N port networks.

This command should be run on the results of the Spectre sp (S-parameter) analysis.

# **Arguments**

$x_iIndex$	The <i>i</i> th index of the coefficient in the scattering matrix.
x_jIndex	The jth index of the coefficient in the scattering matrix.
s_resultName	Results from an analysis. When specified, this argument will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult command.
t_resultsDir	Directory containing the PSF files (results). If you supply this argument, you must also supply the resultName argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the openResults command.

### Value Returned

o_waveform	waveform object representing the S-parameter.
nil	Returns nil if there is an error.

## **Example**

```
s21 = sp(2 1)

s12 = sp(1 2)

plot(s21 s12)
```

### **Data Access Commands**

s11 = sp(1 1 ?result "sp" ?resultsDir "./simResult/psf")

Returns the S-parameter s11 for results of S-parameter(sp) analysis stored at the location ./simResult/psf.

#### **Data Access Commands**

# sprobeData

```
sprobeData(
    t_ana
    t_sprobeInst
    [ ?type t_type ]
    [ ?dataDir t_dataDir ]
)
    => o waveform / nil
```

## **Description**

Returns the waveform for the specified analysis and parameter type of the given sprobe instance.

t ana

Specifies the type of analysis.

Valid values:

- 'sp\_probe Returns the waveforms of "S1", "S2", "Z1" or "Z2" as specified by the t\_type argument.
- 'sp\_sprobe\-StabIndex Returns the waveform of stability index (StabIndex)

t\_sprobeInst

Specifies the name of the sprobe instance.

?type t\_type

One or more strings representing the names of the type.

Valid values:

- "S1", "S2", "Z1" or "Z2" (Scattering parameters)
- "StabIndex" (Stability Index)

?dataDir t\_dataDir Directory containing the PSF file. By default, it uses the current results directory that is set by the openResults SKILL command.

### **Data Access Commands**

### Value Returned

o waveform Waveform object that displays a series of points on a grid.

srrWave:XXXXX is used as the waveform object identifier.

nil Returns nil if there is an error and a waveform cannot be

returned.

# **Example**

The following example gets the waveform of parameter S1 of sprobe instance "I0" sprobeData('sp\_sprobe "I0" ?type "S1")

The following example gets the waveform of parameter StabIndex of sprobe instance "IO" sprobeData('sp sprobe\-StabIndex "IO" ?type "StabIndex")

### **Data Access Commands**

# sweepNames

# **Description**

Returns the names of all the sweep variables for either a supplied waveform, a currently selected result (via selectResult()) or a specified result.

#### **Data Access Commands**

### **Arguments**

o waveForm

Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX). When this argument is used, the t\_resultsDir and s\_resultName arguments are ignored.

?result s resultName

Results from an analysis. When specified, this argument will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult command.

?resultsDir t\_resultsDir

Directory containing the PSF files (results). If you supply this argument, you must also supply the resultName argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the openResults command.

#### Value Returned

l\_sweepName

Returns a list of the sweep names.

nil

Returns nil and prints an error message if the sweep names cannot be returned.

### Example

```
selectResult('tran)
sweepNames()
=> ( "TEMPDC" "time" )
```

Returns a list of sweep variables for the selected results. In this case, the return values indicate that the data was swept over temperature and time.

```
sweepNames(?result 'ac)
=> ("TEMPDC" "freq")
sweepNames()
=> ("TEMPDC" "time")
```

### **Data Access Commands**

```
w = VT("/vout")
sweepNames(w)
=> ("r" "time")
```

Returns the sweep variables for the waveform w.

```
sweepNames(?result 'ac ?resultsDir "./test/psf")
=> ("TEMPDC" "freq")
```

Returns the sweep variables for the results of the ac analysis stored at the location <code>./test/psf.</code>

#### **Data Access Commands**

# sweepValues

```
sweepValues(
     [ o_waveForm ]
    )
    => 1 sweepValues / nil
```

## **Description**

Returns the list of sweep values of the outermost sweep variable of either the selected results or the supplied waveform. This command is particularly useful for parametric analyses.

## **Arguments**

o\_waveForm Waveform object representing simulation results that can be

displayed as a series of points on a grid. (A waveform object

identifier looks like this: srrWave: XXXXX.)

### Value Returned

1 sweepValues Returns the list of sweep values.

nil Returns nil and an error message if the list of sweep values

cannot be returned.

## **Example**

```
sweepValues()
=> ( -50 -15 20 55 90.0 )
```

Returns a list of sweep values for the selected results. In this case, the return values indicate the temperature over which the data was swept.

```
w = VT("/vout")
sweepNames( w )
=> ( "r" "time" )
sweepValues( w )
=> ( 2000 4000 6000 )
```

Returns a list of sweep values for the wave  $\mbox{w}$ . In this case, the return values indicate the resistance over which the data was swept.

#### **Data Access Commands**

# sweepVarValues

```
sweepVarValues(
    [ t_varName ]
    [ ?result s_resultName [ ?resultsDir t_resultsDir ] ]
    )
    => 1 sweepName / nil
```

## **Description**

Returns the list of sweep values for a particular swept variable name. This command is particularly useful for parametric analyses.

## **Arguments**

*t\_varName* Name of the specific variable from which the values are retrieved.

?result s resultName

Results from an analysis. When specified, this argument will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult command.

?resultsDir t resultsDir

Directory containing the PSF files (results). If you supply this argument, you must also supply the resultName argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the openResults command.

#### Value Returned

1\_sweepValuesnilReturns the list of sweep values.nil and an error message if the list of sweep values cannot be returned.

## **Example**

selectResult('tran)

#### **Data Access Commands**

```
sweepNames()
=> ("TEMPDC" "Vsupply" "time")
sweepVarValues("TEMPDC")
=> (0 32)
sweepNames(?result 'ac)
=> ("TEMPDC" "Vsupply" "freq")
sweepVarValues("Vsupply" ?result 'ac)
=> (5 12 15)
sweepNames(?result 'ac ?resultsDir "./simResult/psf")
=> ("TEMPDC" "freq")
sweepVarValues("TEMPDC" ?result 'ac ?resultsDir "./simResult/psf")
=> (-15 20 55)
```

### **Data Access Commands**

### ٧

```
v(
    t_net
    [ ?result s_resultName [ ?resultsDir t_resultsDir ] ]
)
=> o_waveform / nil
```

# **Description**

Returns the voltage of the specified net.

#### **Data Access Commands**

#### **Arguments**

t net Name of the net.

?result s resultName Re

Results from an analysis. When specified, this argument will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult command.

?resultsDir t\_resultsDir

Directory containing the PSF files (results). If you supply this argument, you must also supply the resultName argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the openResults command.

#### Value Returned

o waveform object. A waveform object represents

simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like

this: srrWave: XXXXX.).

nil Returns an error message and nil if there is a problem.

### Example

```
selectResult('tran)
v( "/net56" )
```

Returns the voltage for net56.

```
ocnPrint( v( "/net56" ) )
```

Prints tabular information representing the voltage for net 56.

```
ocnPrint( v( "net5" ?result 'dc ) )
```

Prints the voltage of net5 with respect to the dc swept component.

```
ocnPrint( v( "net5" ?resultsDir "./test2/psf" ?result 'dc ) )
```

Prints the voltage of net5 with respect to dc for the results from a different run (stored in test2/psf).

#### **Data Access Commands**

#### vswr

```
vswr(
    x_index
    [ ?result s_resultName [ ?resultsDir t_resultsDir ] ]
    )
    => o waveform / nil
```

### **Description**

Computes the voltage standing wave ratio.

This function is a higher level wrapper for the OCEAN expression

```
(1 + mag(s(x_index x_index))) / (1 - mag(s(x_index x_index)))
```

This command should be run on the results of the Spectre sp (S-parameter) analysis.

#### **Data Access Commands**

#### **Arguments**

 $x_index$  Index of the port.

?result s resultName Results from an analysis. When specified, this argument

will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult

command.

?resultsDir t resultsDir

Directory containing the PSF files (results). If you supply this argument, you must also supply the resultName argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the openResults command.

#### **Value Returned**

o\_waveform Object representing the voltage standing wave

ratio.

nil Returns an error message or nil if there is a problem.

#### **Example**

```
plot( vswr(2) )
vswr1 = vswr(1 ?result "sp" ?resultsDir "./simResult/psf")
```

Returns the voltage standing wave ratio value at port 1 for the results of S-parameter(sp) analysis stored at the location ./simResult/psf.

#### **Data Access Commands**

#### zm

```
zm(
    x_index
    [ ?result s_resultName [ ?resultsDir t_resultsDir ] ]
    )
    => o waveform / nil
```

### **Description**

Computes the port input impedance.

The zm function is computed in terms of the S-parameters and the reference impedance. This function is a higher level wrapper for the OCEAN expression

This command should be run on the results of the Spectre sp (S-parameter) analysis.

#### **Data Access Commands**

#### **Arguments**

 $x_index$  Index of the port.

 $?result s\_resultName$  Results from an analysis. When specified, this argument

will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult

command.

?resultsDir t resultsDir

Directory containing the PSF files (results). If you supply this argument, you must also supply the resultName argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the openResults command.

#### **Value Returned**

o\_waveform bject representing the port input impedance.

nil Returns an error message and nil if there is a problem.

#### Example

```
plot(zm(2))
zm1 = zm(1 ?result "sp" ?resultsDir "./simResult/psf")
```

Returns input impedance at port 1 for results of S-parameter (sp) analysis stored at the location ./simResult/psf.

#### **Data Access Commands**

#### zref

```
zref(
    x_portIndex
    [ ?result s_resultName [ ?resultsDir t_resultsDir ] ]
   )
    => f_impedance / nil
```

### **Description**

Returns the reference impedance for an N-port network.

This command should be run on the results of the Spectre sp (S-parameter) analysis.

#### **Arguments**

 $x_portIndex$  Index of the port.

?result  $s_resultName$ 

Results from an analysis. When specified, this argument will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult command.

?resultsDir t resultsDir

Directory containing the PSF files (results). If you supply this argument, you must also supply the resultName argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the openResults command.

#### Value Returned

f\_impedance Reference impedance.

nil Returns an error message and nil if there is a problem.

### **Example**

```
Zref = zref(2)
zref1 = zref(1 ?result "sp" ?resultsDir "./simResult/psf")
```

### **Data Access Commands**

Returns the reference impedance at port 1 for the results of S-parameter(sp) analysis stored at the location./simResult/psf.

**Data Access Commands** 

# **Plotting and Printing Commands**

This chapter contains information on the following plotting and printing commands:

- addSubwindow
- addSubwindowTitle
- addTitle
- addWaveLabel
- addWindowLabel
- clearAll
- clearSubwindow
- <u>currentSubwindow</u>
- currentWindow
- dbCompressionPlot
- dcmatchSummary
- <u>deleteSubwindow</u>
- deleteWaveform
- <u>displayMode</u>
- getAsciiWave
- graphicsOff
- graphicsOn
- hardCopy
- hardCopyOptions
- ip3Plot

### Plotting and Printing Commands

- newWindow
- noiseSummary
- ocnGenNoiseSummary
- ocnPrint
- ocnPrintTMIReliabilityResults
- ocnPrintTMIResultTypeList
- ocnSetAttrib
- ocnWriteLsspToFile
- ocnYvsYplot
- plot
- plotStyle
- printGraph
- pzFrequencyAndRealFilter
- pzPlot
- pzSummary
- removeLabel
- report
- saveGraphImage
- xLimit
- yLimit

This chapter also includes a topic, Plotting and Printing SpectreRF Functions in OCEAN.

### Plotting and Printing Commands

### addSubwindow

```
addSubwindow()
=> x_subwindowID / nil
```

### **Description**

Adds a subwindow to the current Waveform window and returns the number for the new subwindow, which is found in the upper right corner.

### **Arguments**

None.

#### Value Returned

 $x\_subwindowID$  Returns the window ID of the new subwindow.

nil Returns nil and an error message if there is no current

Waveform window.

### **Example**

addSubwindow()
=>3

Adds a new subwindow to the Waveform window.

### Plotting and Printing Commands

### addSubwindowTitle

```
addSubwindowTitle(
    x_windowtitle
)
=> t / nil
```

### **Description**

Adds a title to the current subwindow in the active window. The current subwindow is defined using the currentSubwindow command.

### **Arguments**

 $x\_windowtitle$  User-defined title for the subwindow.

#### **Value Returned**

t The user-supplied name of the current subwindow.

nil Returns nil if the title is not created.

### **Example**

```
addSubwindowTitle( "waveform 2")
=> +
```

Adds the title waveform 2 to the selected subwindow.

### Plotting and Printing Commands

### addTitle

```
addTitle(
    x_windowtitle
)
=> t / nil
```

### **Description**

Adds a title to the current active OCEAN window. The current window is defined using the currentWindow command.

### **Arguments**

 $x\_windowtitle$  User-defined title for the window.

#### Value Returned

t The user-supplied name of the current window.

nil Returns nil if the title is not created.

### **Example**

```
addTitle( "waveform 1" )
=> t
```

Adds the title waveform 1 to the selected window.

### Plotting and Printing Commands

### addWaveLabel

```
addWaveLabel(
    x_waveIndex
    l_location
    t_label
    [ ?textOffset g_textOffset ]
    [ ?color x_color ]
    [ ?justify t_justify ]
    [ ?fontStyle t_fontStyle ]
    [ ?height x_height ]
    [ ?orient t_orient ]
    [ ?drafting g_drafting ]
    [ ?overBar g_overbar ])
    => s_labelId / nil
```

# **Description**

Attaches a label to the specified waveform curve in the current subwindow.

# Plotting and Printing Commands

# **Arguments**

x_waveIndex	Integer identifying the waveform curve.
l_location	List of two waveform coordinates that describe the location for the label.
t_label	Label for the waveform.
g_textOffset	Boolean that specifies whether to place a marker or label. If set to ${\tt t}$ , a marker is placed. If set to ${\tt nil}$ , a label is placed. Default value: ${\tt nil}$ .
x_color	Label color specified as an index in the technology file.  Default value: 10
t_justify	Justification, which is specified as "upperLeft", "centerLeft", "lowerLeft", "upperCenter", "centerCenter", "lowerCenter", "upperRight", "centerRight", Or "lowerRight".  Default value: "lowerLeft"
t_fontStyle	Font style, which is specified as "euroStyle", "gothic", "math", "roman", "script", "stick", "fixed", "swedish", "raster", or "milSpec".  Default value: the font style of the current subwindow
x_height	Height of the font.  Default value: the font height of the current subwindow
t_orient	Orientation of the text, specified as either "R0" or "R90".  Default value: "R0"
g_drafting	Boolean that specifies whether the label stays backwards or upside-down. If set to $t$ , a backwards or upside-down label is displayed in a readable form. If set to $nil$ , a backwards or upside-down label stays the way it is. Default value: $t$
g_overbar	Boolean that specifies whether underscores in labels are displayed as overbars. If set to $t$ , underscores in labels are displayed as overbars. If set to $nil$ , underbars are displayed as underbars. Default value: $nil$

### Plotting and Printing Commands

#### Value Returned

 $s\_labelId$  Returns an identification number for the waveform label.

nil Returns nil if there is an error.

#### **Examples**

```
addWaveLabel( 1 list( 0 0.5 ) "R5 = " )
```

Attaches the "R5 = " label to the specified coordinates on waveform curve 1.

```
addWaveLabel( 2 list( 0 0.5 ) "R_6 = " ?textOffset 0:20 ?justify "lowerCenter" ?fontStyle "roman" ?height 10 ?orient "R20" ?drafting t ?overbar t)
```

Attaches the label "R6 = " to the specified coordinates on waveform curve. The label specifications are as follows: Justification - lowerCenter, Font Style - roman, Font Height - lowerCenter, and Orientation - R20.

The label will be displayed in a readable form. The underscore in the label will be displayed as an overbar.

#### **Additional Information**

Note the following points:

- The valid label location ranges between absolute co-ordinates (0, 0) on X-axis and (1,1) on Y-axis (upper and lower bound inclusive).
- The valid marker location ranges between data co-ordinates defined by X-axis and Y-axis limits (upper and lower bound inclusive).

#### Case1:

```
addWaveLabel(1 list( -0.5 -0.5 ) "Label 1" ?textOffset nil)
```

The following error message appears when the specified label location (-0.5 0.5) is outside of the defined boundary limits of label.

The location specified for placing a label on the graph is invalid. Specify a valid label location that ranges between absolute coordinates (0,0) on X-axis and (1,1) on Y-axis (upper and lower bounds inclusive).

#### Case 2:

```
addWaveLabel(1 list( 80MHz -0.5) "Marker 1" ?textOffset t)
```

The following error message appears when the specified marker location (80MHz - 0.5) is outside of the X- and Y-axis limits of the graph to be plotted.

### Plotting and Printing Commands

The location specified for placing a marker on the graph is invalid. Specify a valid marker location that ranges between data coordinates '(0,-1)' on X-axis and '(10000,1)' on Y-axis (upper and lower bounds inclusive).

### Plotting and Printing Commands

### addWindowLabel

### **Description**

Displays a label in the current subwindow. The location for the label is specified with a list of two numbers between 0 and 1.

# **Arguments**

l_location	List of two waveform coordinates that describe the location for the label.
	Valid values: 0 through 1
t_label	Label for the waveform.

### Value Returned

$s\_labelId$	Returns an identification number for the subwindow label.
nil	Returns nil if there is an error.

### **Example**

```
label = addWindowLabel( list( 0.75 0.75 ) "test" )
```

Adds the test label to the current subwindow at the specified coordinates and stores the label identification number in label.

### Plotting and Printing Commands

### clearAll

### **Description**

Erases the contents of the current Waveform window and deletes the waveforms, title, date stamp, and labels stored in internal memory.

### **Arguments**

None.

#### Value Returned

t Returns t if the waveform information is deleted.

nil Returns nil and an error message if there is no current

Waveform window.

### **Example**

```
clearAll()
=> t
```

Erases the contents of the current Waveform window.

### Plotting and Printing Commands

# clearSubwindow

# **Description**

Erases the contents of the current subwindow.

# **Arguments**

None.

#### **Value Returned**

t Returns t if the contents of the subwindow are erased.

nil Returns nil and an error message otherwise.

### **Example**

```
clearSubwindow()
=> +
```

Erases the contents of the current subwindow.

### Plotting and Printing Commands

### currentSubwindow

```
currentSubwindow(
    [ ?x_subwindow x_subwindow ]
    )
    => t / nil
```

### **Description**

Sets *x* subwindow as the current subwindow.

### **Arguments**

```
?x subwindow x subwindow
```

(Optional) Number of the subwindow, found in the upper right corner, that is to become the current subwindow.

#### Value Returned

t	Returns t when the subwindow is set to
	x_subwindow. If you do not specify any argument in
	this function, it returns the current subwindow number.
nil	If no subwindow exists.

### **Example**

```
currentSubwindow( 2 )
```

Sets subwindow 2 as the current subwindow.

### Plotting and Printing Commands

#### currentWindow

```
currentWindow(
    w_windowId
)
    => w windowId / nil
```

### **Description**

Specifies w windowId as the current Waveform window.

### **Arguments**

w\_windowId Waveform window ID.

#### Value Returned

w windowId Returns the current Waveform window ID.

nil Returns nil and an error if the current window cannot be set.

#### **Example**

```
currentWindow( window(2) )
```

This example specifies window 2 as the current Waveform window.

```
currentWindow()
```

This example returns the current waveform window. For example, if the current waveform window is 4, this command returns the following:

window:4

#### Plotting and Printing Commands

# dbCompressionPlot

```
dbCompressionPlot(
    o_wave
    x_harmonic
    x_extrapolationPoint
    [?compression x_compression]
)
    => t / nil
```

### **Description**

Plots the nth compression point plot. The  $x\_compression$  argument is optional and defaults to 1 for 1dB compression, if omitted.

This command should be run on the results of the Spectre swept pss analysis.

### **Arguments**

o_wave	The waveform for which to plot the compression.
x_harmonic	Harmonic frequency index.
$x\_extrapolationPoint$	The extrapolation point.
?compression	The amount of dB compression.
$x\_compression$	Default value: 1

#### Value Returned

t	Returns t if the point is plotted
nil	returns nil if there was an error

#### **Example**

#### Plotting and Printing Commands

### dcmatchSummary

```
dcmatchSummary(
    [ ?resultsDir t_resultsDir ]
    [ ?result S_resultName ]
    [ ?output t_fileName | p_port ]
    [ ?paramValues ln_paramValues ]
    [ ?deviceType ls_deviceType ]
    [ ?variations ls_variations ]
    [ ?includeInst lt_includeInst ]
    [ ?excludeInst lt_excludeInst ]
    [ ?truncateData n_truncateData ]
    [ ?truncateType s_truncateType ]
    [ ?sortType ls_sortType ]
    )
    => t_fileName / p_port/ nil
```

### **Description**

Prints a report showing the mismatch contribution of each component in a circuit. If you specify a directory with resultsDir, it is equivalent to temporarily using the <code>openResults</code> command. The <code>dcmatchSummary</code> command prints the results for that directory and resets the <code>openResults</code> command to its previous setting. If you specify a particular result with <code>resultName</code>, it is equivalent to temporarily using the <code>selectResult</code> command on the specified results. The <code>dcmatchSummary</code> command prints the results and resets the <code>selectResult</code> command to its previous setting.

This command should be run on the results of the Spectre dcmatch analysis.

# Plotting and Printing Commands

# **Arguments**

?resultsDir t_resultsDir	The directory containing the dcmatch-analysis results.
?results S_resultName	Results from an analysis for which you want to print the dcmatchSummary report.
?output t_fileName	File in which to write the information. The dcmatchSummary command opens the file, writes to the file and closes the file. If you specify the filename without a path, the dcmatchSummary command creates the file in the directory pointed to by your SKILL Path. To find out what your SKILL path is, type getSkillPath() at the OCEAN prompt.
?output <i>p_port</i>	Port (previously opened with outfile) through which to append the information to a file. You are responsible for closing the port. See the <u>outfile</u> command for more information.
?paramValues ln_paramValues	List of values for swept parameters at which the dcmatchSummary is to be printed. In case there is just one swept parameter the value can be specified as is.
?deviceType ls_deviceType	List of device type strings to be included. Valid values are a list of strings or 'all or a single device name. Default value is 'all.
?variations ls_variations	An association list containing the device name and the associated variations to print. You can also specify the value 'all to print all available variations for a device. Default value is 'all. For Example: '( ("bsim3v3" ("sigmaOut")) "sigmaVth")) ("resistor" ("sigmaOut"))
?includeInst lt_includeInst	List of instance name strings to definitely include in the dcmatchSummary.
?excludeInst lt_excludeInst	List of instance name strings to exclude in the dcmatchSummary.
?truncateData x_truncateData	Specifies a number that the truncateType argument uses to define the components for which information is to be printed.

#### Plotting and Printing Commands

?t	runcateType
s	truncateType

Specifies the method that is used to limit the data being included in the report

Valid values:

- 'top Saves information for the number of components specified with truncateData. The components with the highest contributions are saved. Sample value for truncateData: 10
- 'relative Saves information for all components that have a higher contribution than truncateData \* maximum. Where maximum is the maximum contribution among all the devices of a given type. Sample value for truncateData: 1.9n
- 'absolute Saves information for all the components in the selected set whose contribution are more than truncateData. Sample value for truncateData: 0.1
- 'none Saves information for all the components.
   Sample value for truncateData: Not required

?sortType
ls sortType

Specifies how the printed results are to be sorted. The valid values are nil, 'name, 'output.

#### Value Returned

t_fileName	Returns the name of the port.
p_port	Returns the name of the file.
nil	Returns nil and an error me

Returns nil and an error message if the summary cannot be printed.

### **Example**

```
dcmatchSummary( ?result 'dcmatch-mine )
```

Prints a report for non-swept DC-Mismatch analysis.

```
dcmatchSummary( ?resultsDir "/usr/simulation/lowpass/spectre/schematic" ?result
'dcmatch)
```

Prints a report for non-swept DC-Mismatch analysis for the results from a different run (stored in the schematic directory).

#### Plotting and Printing Commands

dcmatchSummary( ?resultsDir "/usr/simulation/lowpass/spectre/schematic" ?result
'dcmatch ?paramValues `(25) )

Prints a report for swept DC-Mismatch analysis at swept parameter value of 25.

```
dcmatchSummary( ?result dcmatch-mine ?output "./summary.out")
```

Prints a report for non-swept DC-Mismatch analysis in the output file summary.out.

```
dcmatchSummary( ?paramValues 25 ?deviceType "bsim3v3" ?variations '(("bsim3v3"
("sigmaOut "sigmaVth" )))
```

Prints a report for swept DC-Mismatch analysis at swept parameter value of 25 for bsim3v3 deviceType and sigmaOut and sigmaVth variations.

```
dcmatchSummary( ?paramValues 25 ?truncateType 'top ?truncateData 1)
```

Prints a report for swept DC-Mismatch analysis at swept parameter value of 25 printing only the component having the highest contribution.

```
dcmatchSummary( ?paramValues 25 ?sortType 'name )
```

Prints a report for swept DC-Mismatch analysis at swept parameter value of 25 sorted on name.

### Plotting and Printing Commands

### deleteSubwindow

### **Description**

Deletes the current subwindow from the current Waveform window.

# **Arguments**

None.

### **Value Returned**

t Returns t if the current subwindow is deleted.

nil Returns nil and an error message if there is no current

subwindow.

### **Example**

```
deleteSubwindow()
=> t
```

Deletes the current subwindow from the Waveform window.

### Plotting and Printing Commands

### deleteWaveform

### **Description**

Deletes the specified waveform curve or all the waveform curves from the current subwindow of a Waveform window.

### **Arguments**

x_index	Integer identifying a particular waveform curve.
all_string	The string "all" specifying that all waveform curves are to be

deleted.

#### Value Returned

t Returns t if the curves are deleted.

nil Returns nil and an error message if the curves are not

deleted.

#### **Example**

```
deleteWaveform( '1 )
=> t
```

Deletes waveform 1 from the current subwindow.

```
deleteWaveform( "all" )
=> t
```

Deletes all the curves from the current subwindow.

#### Plotting and Printing Commands

# displayMode

```
displayMode(
    t_mode
)
    => t / nil
```

### **Description**

Sets the display mode of the current subwindow.

### **Arguments**

t mode String representing the display mode for the subwindow.

Valid values: strip, composite, or smith.

**Note:** This also works if a plot is not open.

#### **Value Returned**

t Returns t when the display mode of the subwindow is set.

nil Returns nil and an error message if the display mode cannot

be set.

### **Example**

```
displayMode( "composite" )
=> +
```

Sets the current subwindow to display in composite mode.

### Plotting and Printing Commands

# getAsciiWave

```
getAsciiWave(
    t_filename
    x_xColumn
    x_yColumn
    [?xskip x_x_skip]
    [?yskip x_yskip]
    [?formatFloat g_formatFloat]
    [?xName t_xName]
    [?xUnits t_xUnits]
    [?yName t_yName]
    [?yUnits t_yUnits]
)
    => o_wave / nil
```

### **Description**

Reads in an ASCII file of data and generates a waveform object from the specified data. The X-axis data must be real numbers. The Y-axis data can be real or complex values. Complex values are represented as (real imag) or complex (real imag). This function skips blank lines and comment lines. Comments are defined as lines beginning with a semicolon.

### Plotting and Printing Commands

### **Arguments**

t_filename	The name of the Ascii file to be read in.
$x\_xColumn$	The column in the data file that contains the X-axis data.
x_yColumn	The column in the data file that contains the Y-axis data.
?xskip x_xskip	The number of lines to skip in the X column.
?yskip x_yskip	The number of lines to skip in the Y column.
?formatFloat g_formatFloat	A boolean when set to t converts the integer data into float values. Default value: $nil$ .
?xName t_xName	The name of the X-axis.
?xUnits t_xUnits	The units for the X-axis.
?yName t_yName	The name of the Y-axis.
?yUnits t_yUnits	The units for the Y-axis.

#### Value Returned

o wave	The waveforn	າ object.
--------	--------------	-----------

nil Returns nil if the function fails.

#### **Example**

```
getAsciiWave("~/mydatafile.txt " 1 2 )
=> srrWave: 32538648
```

Reads in an ascii file ~/mydatafile.txt, which has x-axis data in the first column and y-axis data in the second column, and returns a waveform object.

```
getAsciiWave("~/mydatafile.txt " 1 2 ?xskip 1 ?yskip 2)
=> srrWave:32538656
```

Reads in an ascii file ~/mydatafile.txt, which has x-axis data in the first column and y-axis data in the second column and skips 1 line in the x\_xcolumn and 2 lines in the y ycolumn, and returns a waveform object.

```
getAsciiWave("/vout_tran_xy.csv" 1 2 ?xskip 1 ?yskip 1 ?formatFloat nil ?xName
"Time" ?xUnits "Secs" ?yName "V(Out)" ?yUnits "V")
```

Reads in an ascii file, vout\_tran\_xy.csv, which has x-axis data in the first column and y-axis data in the second column, skips 1 line in the x xcolumn and 2 lines in the y ycolumn,

# Plotting and Printing Commands

x- and y-axis names are  $\mathtt{Time}$  and  $\mathtt{V}(\mathtt{out})$  , and x-and y-axis units are  $\mathtt{Secs}$  and  $\mathtt{V}$  respectively.

### Plotting and Printing Commands

# graphicsOff

### **Description**

Disables the redrawing of the current Waveform window.

You might use this command to freeze the Waveform window display, send several plots to the window, and then unfreeze the window to display all the plots at once.

### **Arguments**

None.

#### Value Returned

t Returns t if redrawing is disabled.

nil Returns nil if there is an error, such as there is no current

Waveform window.

### **Example**

```
graphicsOff()
=> t
```

Disables the redrawing of the Waveform window.

## Plotting and Printing Commands

# graphicsOn

# **Description**

Enables the redrawing of the current Waveform window.

# **Arguments**

None.

### **Value Returned**

t Returns t if redrawing is enabled.

nil Returns nil if there is an error, such as there is no current

Waveform window.

# **Example**

```
graphicsOn()
=> t
```

Enables the redrawing of the current Waveform window.

### Plotting and Printing Commands

# hardCopy

```
hardCopy(
    w_windowId
)
    => t / nil
```

## **Description**

Sends a Waveform window plot to a printer or a file. To plot to a printer specify a printer name using the ?hcPrinterName argument of the hardCopyOptions command. To plot to a file, specify a file name using the ?hcOutputFile argument of the hardCopyOptions command.

**Note:** You must first set any plotting options with the <u>hardCopyOptions</u> command.

## **Arguments**

w windowId

The window ID of the waveform window whose plot is to be sent to a printer or a file. The default value is the window ID of the current window.

#### Value Returned

t Returns t if successful.

nil Returns nil if there is an error.

# **Example**

```
hardCopy()
=> t
```

Sends a waveform plot to the printer or to a file.

```
w = newWindow()
plot(v("/vout"))
hardCopy(w)
```

Sends the waveform plot of w to the printer or to a file.

### Plotting and Printing Commands

# hardCopyOptions

```
hardCopyOptions(
     [ ?hcCopyNum x hcCopyNum ]
     [ ?hcOffsetHeight x hcOffsetHeight ]
     [ ?hcOffsetWidth x hcOffsetWidth ]
     [ ?hcOrientation s hcOrientation ]
     [ ?hcOutputFile g hcOutputFile ]
     [ ?hcPrinterName s hcPrinterName ]
     [ ?hcTmpDir t_hcTmpDir ]
     [ ?hcPaperSize s hcPaperSize ]
     [ ?hcMakeExactCopy g hcMakeExactCopy ]
     [ ?hcQuality x hcQuality ]
     [ ?hcOptimizeForWindows g hcOptimizeForWindows ]
     [ ?hcImageWidth x hcImageWidth ]
     [ ?hcImageHeight x hcImageHeight ]
     [ ?hcImageSizeUnits s hcImageSizeUnits ]
     [ ?hcImageResolution x ImageResolution ]
     [ ?hcResolutionUnits s hcResolutionUnits ]
     [ ?hcImageAspectRatio x hcImageAspectRatio ]
     [ ?hcUseExistingBackground g hcUseExistingBackground ]
     [ ?hcDisplayTitle g hcDisplayTitle ]
     [ ?hcDisplayLegend g hcDisplayLegend ]
     [ ?hcDisplayAxes q hcDisplayAxes ]
     [ ?hcDisplayGrids g hcDisplayGrids ]
     [ ?hcSaveEachSubwindowSeparately g hcSaveEachSubwindowSeparately ]
    => g value / nil
```

## **Description**

Sets the graph window hardcopy plotting options.

The option takes effect for any graph window or subwindow that is opened after the option is set.

## **Arguments**

?hcCopyNum  $x_hcCopyNum$  The number of copies to plot.

Valid values: any positive integer

Default value: 1

### Plotting and Printing Commands

?hcOffsetHeight
x hcOffsetHeight

The vertical margin.

Valid values: any positive integer

Default value: 0

?hcOffsetWidth
x hcOffsetWidth

The horizontal margin.

Valid values: any positive integer

Default value: 0

?hcOrientation
s\_hcOrientation

The plot orientation.

**Note:** This option works only when you print a graph window and does not work when you save the window to a graph file.

Valid values: 'portrait, 'landscape,
'automatic

Default value: 'automatic

?hcOutputFile
g hcOutputFile

Name of the output file. The output file can be created in one of the following file formats:

- BMP Windows Device Independent Bitmap (.bmp)
- PNG Portable Network Graphics(.png)
- PS PostScript (.ps)
- TIFF Tagged Image File Format (.tif)
- TIFF Tagged Image File Format (.tif)
- EPS Encapsulated Post Script (.eps)
- PDF Portable Document Format (.pdf)
- PPM Portable PixMap File (.ppm)
- JPG Joint Photographic Experts Group (.jpg)
- SVG Scalable Vector Graphics (.svg)
- XPM X PixMap (.xpm)

Valid values: a string or nil

Default value: nil

### Plotting and Printing Commands

?hcPrinterName The name of the printer.
s hcPrinterName

Valid values: a string or nil

Default value: nil

?hcTmpDir t hcTmpDir The name of a temporary directory to be used for

scratch space.

Valid values: name of a temporary directory

Default value: "/usr/tmp"

?hcPaperSize The paper size. The available paper sizes are—
s hcPaperSize letter, legal, executive, folio, ledger,

letter, legal, executive, folio, ledger, tabloid, a0, a1, a2, a3, a4, a5, a6, a7, a8, a9, b0,

b1, b2, b3, b4, b5, b6, b7, b8, b9, b10, c5e,

comm10e, dle.

Default value: a4

?hcMakeExactCopy Saves the exact copy of all subwindows. Only g hcMakeExactCopy ?hcQuality and ?hcOutputFile argument

?hcQuality and ?hcOutputFile arguments work with this option. This option does not work for the eps

file format.

Valid values: t or nil

Default value: nil

?hcQuality  $x_hcQuality$  Modifies the quality of the image. This option works

only for the .jpeg file format. This option does not

work for the eps file format.

Valid values: 20 to 100%

Default value: 85

?hcOptimizeForWindows
g hcOptimizeForWindows

Enables the image to be imported in the Microsoft office application. This option is available when you select the image type as Encapsulated PostScript (\*.eps). This option simplifies the image output so that it can be ready by Microsoft Office 2003 and 2007

applications

Valid values: t or nil

Default value: t

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### Plotting and Printing Commands

?hcImageWidthSets the width of the image.x\_hcImageWidthValid values: Any positive integer value.Default value: 800 pixels?hcImageHeightSets the height of the image

 $x\_hcImageHeight$  Valid values: Any positive integer value.

Default value: 600 pixels

?hcImageSizeUnits Specifies the unit for image size (height and width)

s\_hcImageSizeUnits Valid values: inch, cm, mm, picas, pixels, and

valid values. Then, elli, illili, preas, prixers, and

points

Default value: pixels

?hcImageResolution Sets the image resolution. This option works only for  $x_hcImageResolution$  the bmp, jpeg, png, ppm, tif, and xpm file formats. It does not work for eps, pdf, and svg file formats.

Valid values: Any positive integer value.

Default value: 96

?hcResolutioUnits

Sets the units for image resolution. This option works
only for the bmp, jpeg, png, ppm, tif, and xpm file
formats. It does not work for eps, pdf, and svg file

formats.

Valid values: pixels/cm and pixels/in

Default value: pixels/in

?hcImageAspectRatio Enables the aspect ratio, which is the ratio of the width of the image to its height.

Valid values: t or nil

Default value: nil

?hcUseExistingBackgroun Enables to use the existing background in the graph image.

g\_hcUseExistingBackgrou
nd
Valid values: t or nil

Default value: nil

### Plotting and Printing Commands

?hcDisplayTitle Displays the trace title in the graph image.

g\_hcDisplayTitle Valid values: t or nil

Default value: t

?hcDisplayLegend Displays the trace legend in the graph image.

Default value: t

?hcDisplayAxes Displays the axes in the graph image.

g\_hcDisplayAxes Valid values: t or nil

Default value: t

?hcDisplayGrids Displays the grids in the graph image.

g\_hcDisplayGrids Valid values: t or nil

Default value: t

?hcSaveEachSubWindowSep Saves all subwindows in a graph to a single file or

erately multiple files.

 $g_hcSaveEachSubwindowSe$  parately Valid values: t (multiple files) or nil (single file)

Default value: t

Value Returned

g value Returns the new value of the option.

nil Returns nil if there is an error.

**Example** 

hardCopyOptions( ?hcCopyNum 1 )

Plots one copy of the window or subwindow.

hardCopyOptions(?hcCopyNum 3 ?hcOutputFile "myOutFile.bmp")

Plots three copies of the window or subwindow and sends them to the file myOutFile.bmp.

# Plotting and Printing Commands

# ip3Plot

```
ip3Plot(
    o_wave
    x_sigHarmonic
    x_refHarmonic
    x_extrapolationPoint
)
    => t / nil
```

# **Description**

Plots the IP3 curves.

This command should be run on the results of the Spectre swept pss and pac analysis.

Refer to the "Simulating Mixers" chapter of the *Virtuoso Spectre Circuit Simulator RF Analysis User Guide* for more information on ip3Plot.

# **Arguments**

o_wave	Waveform for which to plot the ip3.
$x\_sigHarmonic$	Index of the third order harmonic.
$x\_refHarmonic$	Index of the first order (fundamental) harmonic.
x_extrapolationPoint	Extrapolation point.

### Value Returned

t	Returns t if the curves are plotted.
nil	Returns nil if there is an error.

# **Example**

```
ip3Plot(v("/net28") 47 45 -25)
```

## Plotting and Printing Commands

# newWindow

# **Description**

Creates a new Waveform window and returns the window ID.

# **Arguments**

None.

### Value Returned

w windowId Returns the window ID of the new	v Waveform window.
---	--------------------

nil Returns nil and an error message if the new Waveform

window cannot be created.

# **Example**

```
newWindow()
=> window:3
```

Creates a new Waveform window that is numbered 3 in the upper right corner.

### Plotting and Printing Commands

# noiseSummary

```
noiseSummary(
    s_type
     [ ?result s_resultName [ ?resultsDir t_resultsDir ] ]
     [ ?frequency f frequency ]
     [ ?weight f weight ]
     [ ?output t_fileName | p_port ]
     [ ?noiseUnit t noiseUnit ]
     [ ?truncateData x_truncateData ]
     [ ?truncateType s truncateType ]
     [ ?digits x digits ]
     [ ?percentDecimals x percentDecimals ]
     [ ?from f from ]
     [ ?to f_to ]
     [ ?deviceType ls_deviceType ]
     [ ?weightFile t weightFile ]
     [ ?paramValues ls paramValues ]
     [ ?hierLevel s_hierLevel ]
     [ ?sort ls sort ]
     [ ?includeInsts ls_includeInsts ]
     [ ?excludeInsts ls excludeInsts ]
     [ ?combineIteratedInsts g combineIteratedInsts ]
     [ ?suffixNotation t suffixNotation ]
    => t / nil
```

### **Description**

Prints a report showing the noise contribution of each component in a circuit.

This command should be run on the results of the Spectre noise analysis.

### Plotting and Printing Commands

### **Arguments**

s type Type of noise-analysis results for which to print the report.

Valid values: spot, to specify noise at a particular frequency, or integrated, to specify noise integrated

over a frequency range.

?result s\_resultName Results from an analysis. When specified, this

argumentwill only be used internally and will not alter the current result which was set by the selectResult command.

The default is the current result selected with the

selectResult command.

?resultsDir Directory containing the PSF files (results). If you supply  $t\_resultsDir$  this argument, you must also supply the resultName

argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the openResults

command.

?frequency Frequency value of interest. f frequency

?weight f\_weight Waveform representing the function with which the integral

is weighted.

Default value: 1.0

?output  $t_fileName$  | The type of output:

p port

■ File in which to write the information. The noiseSummary command opens the file, writes to the file, and closes the file. If you specify the filename without a path, the noiseSummary command creates the file in the directory pointed to by your SKILL Path. To find out what your SKILL path is, type getSkillPath() at the OCEAN prompt.

Port (previously opened with outfile) through which to append the information to a file. You are responsible for closing the port. See the <u>outfile</u> command for more information.

### Plotting and Printing Commands

?r	noiseUnit	
t	noiseUnit	-

Specifies the type of noise unit to be saved.

Valid values:

?truncateData x\_truncateData

?truncateType
s truncateType

Specifies a number that the truncateType argument uses to define the components for which information is to be printed.

Specifies the method that is used to limit the data being included in the report.

Valid values:

- 'top Saves information for the number of components specified with truncateData. The components with the highest contributions are saved. Sample value: 10
- 'relative Prints components which have noise contribution (percent) higher than that specified by ?truncateData. Sample value: .1
- 'none Saves information for all the components.

?digits x digits

?percentDecimals
x\_percentDecimals

?from f from

?to f to

?deviceType
ls deviceType

Number of significant digits with which the contributors are printed.

Number of decimals printed for any relative contribution.

For integrated noise, the start value for frequency.

For integrated noise, the end value for frequency.

List of device type strings to be included.

Valid values: a list of strings or 'all

### Plotting and Printing Commands

?weightFile t weightFile Absolute or relative path of the file that contains information about weights. This data is used to compute weighted noise. If the values are provided for both parameters, weight and weightFile, the value for weight gets precedence.

?paramValues *ls paramValues* 

List of values where each value co-relates to a specific sweep variable name. This field must be used when the data is parametric. The order of this list must coincide with the list returned by the sweepNames function excluding the frequency variable.

?hierLevel s hierlevel Specifies the hierarchy level up to which the noise summary needs to be computed.

?sort ls sort

Determines the order of the report according to the specified symbol categories. Valid values:

- 'individual: Sorts the report from the largest noise contributor to the smallest.
- 'composite: Sorts the report by the total noise contribution of each device. Each device entry contains the percentage of the noise contribution from this device and the noise contribution from each of its contributors.
- 'name: Produces the same format as composite noise but sorts it alphabetically by device instance name.

?includeInsts ls includeInsts

List of strings containing instances to be included.

?excludeInsts ls excludeInsts List of strings containing instances to be excluded.

Valid values: a list of strings or 'none.

Valid values: a list of strings or 'all.

g\_combineIteratedInst one contributor.

?combineIteratedInsts Combines the noise contributions of iterated instances to

Valid values: t or nil

Default value: nil

### Plotting and Printing Commands

?suffixNotation		
t_	suffixNotation	

Changes the notation of the noise contribution value for the data output to the specified value. For example,

3.36916e-05, is shown as 33.6916u.

Default value: nil

#### Value Returned

t Returns t if noise summary is successfully printed to the

file or port.

nil Returns nil and an error message is shown if the

summary cannot be printed.

## **Example**

```
noiseSummary( 'integrated ?result 'noiseSweep-noise )
```

## Prints a report for an integrated noise analysis.

```
noiseSummary( 'integrated ?resultsDir
   "/usr/simulation/lowpass/spectre/schematic"
   ?result 'noise)
```

Prints a report for an integrated noise analysis for the results from a different run (stored in the schematic directory).

```
noiseSummary( 'spot ?resultsDir
   "/usr/simulation/lowpass/spectre/schematic"
   ?result 'noise ?frequency 100M )
```

Prints a report for a spot noise analysis at a frequency of 100M.

```
noiseSummary('integrated ?truncateType 'none ?digits 10
?weightFile "./weights.dat")
```

Prints the weighted noise for an integrated noise analysis using information in the weight file weights.dat.

```
noiseSummary('integrated ?output "./NoiseSum1" ?noiseUnit "V" ?truncateData 20
?truncateType 'top ?from 10 ?to 10M ?deviceType list("bjt" "mos" "resistor"))
```

Prints a report for an integrated noise analysis in the frequency range 10-10M for 20 components with deviceType bjt, mos or resistor.

```
noiseSummary('integrated ?from 1 ?to 100M ?truncateType 'top ?truncateData 20
?deviceType 'all ?noiseUnit "V^2" ?output "./filename.ns" ?paramValues list(2.47e-
9))
```

Prints a report for an integrated noise analysis at a specific swept value.

### Plotting and Printing Commands

# ocnGenNoiseSummary

```
ocnGenNoiseSummary(
    x_hierLevel
    s_resultName
    t_resultsDir
)
=> t / nil
```

## **Description**

Generates a noise summary report at the end of a simulation run. This report contains a summary of noise contribution of each component for the specified hierarchy level.

### **Arguments**

$x\_hierLevel$	Specifies the level of hierarchy upto which the current
	data is to be saved for all the components.

s resultName Results from an analysis.

When specified, this argument will only be used internally and will not alter the current result which was set by the <u>selectResult</u> command. The default is the current result selected with the <code>selectResult</code>

command.

t\_resultsDir Directory containing the PSF files (results).

When specified, this argument will only be used internally and will not alter the current results directory that is set by the <u>openResults</u> command. The default is the current results directory set by the <u>openResults</u> command.

**Note:** If you specify this argument, you must also specify the  $s\_resultName$  argument.

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# Plotting and Printing Commands

### **Value Returned**

t Creates a noise summary file within the results

directory.

nil Returns nil otherwise.

# Example

```
ocnGenNoiseSummary(1 ?result 'pnoise)
=> t
```

Generates a noise summary report for all the components up to the first level of the hierarchy.

### Plotting and Printing Commands

### ocnPrint

```
ocnPrint(
    [ ?output t_filename | p_port ]
    [ ?precision x_precision ]
    [ ?numberNotation s_numberNotation ]
    [ ?numSpaces x_numSpaces ]
    [ ?width x_width ]
    [ ?from x_from ]
    [ ?to x_to ]
    [ ?step x_step ]
    [ ?linLog t_linLog ]
    o_waveform1 [ o_waveform2 ... ]
    )
    => t / nil
```

## **Description**

Prints the text data of the waveforms specified in the list of waveforms.

If you provide a filename as the <code>?output</code> argument, the <code>ocnPrint</code> command opens the file and writes the information to it. If you provide a port (the return value of the SKILL <code>outfile</code> command), the <code>ocnPrint</code> command appends the information to the file that is represented by the port. There is a limitation of <code>ocnPrint</code> for precision. It works upto 30 digits for the Solaris port and 18 digits for HP and AIX.

**Note:** ocnPrint() prints z-state as "HiZ" for digital signals.

## Plotting and Printing Commands

### **Arguments**

?output t filename

File in which to write the information. The ocnPrint command opens the file, writes to the file, and closes the file. If you specify the filename without a path, the OCEAN environment creates the file in the directory pointed to by your SKILL Path. To find out what your SKILL path is, type getSkillPath() at the OCEAN prompt.

?output p port

Port (previously opened with outfile) through which to append the information to a file. You are responsible for closing the port. See the <u>outfile</u> command for more information.

?precision x precision

The number of significant digits to print. This value overrides any global precision value set with the setup command.

Valid values: 1 through 16

Default value: 6

**Note:** To print the specified significant number of digits, ensure that the value of the  $x\_width$  argument is the same or greater than the value of the  $x\_precision$  argument.

?numberNotation s\_numberNotation

The notation for print ed information. This value overrides any global format value set with the setup command.

Valid values: 'suffix, 'engineering, 'scientific, 'none

Default value: 'suffix

The format for each value is 'suffix: 1m, 1u, 1n, etc.; 'engineering: 1e-3, 1e-6, 1e-9, etc.; 'scientific: 1.0e-2, 1.768e-5, etc.; 'none.

The value 'none is provided so that you can turn off formatting and therefore greatly speed up printing for large data files. For the fastest printing, use the 'none value and set the 'output argument to a filename or a port, so that output does not go to the CIW.

### Plotting and Printing Commands

?numSpace  $x_numSpaces$  The number of spaces between columns.

Valid values: 1 or greater

Default value: 4

?width x width The width of each column.

Valid values: 4 or greater

Default value: 14

?from x from The start value at x axis for the waveform to be printed.

?to x to The end value at x axis for the waveform to be printed.

?step x step The step by which text data to be printed is

incremented.

?linLog t linLog The scale to be used for printing.

Valid values: Linear, Log

Default value: Linear

o waveform bject representing simulation results that

can be displayed as a series of points on a grid. (A

waveform object identifier looks like this:

srrWave:XXXXX.)

o waveform2 Additional waveform object.

#### Value Returned

t Returns t if the text for the waveforms is printed.

nil Returns nil and an error message if the text for the

waveforms cannot be printed.

### **Example**

```
ocnPrint( v( "/net56" ) )
=> t
```

Prints the text for the waveform for the voltage of net56.

```
ocnPrint( vm( "/net56" ) vp( "/net56" ) ) => t
```

Prints the text for the waveforms for the magnitude of the voltage of net56 and the phase of the voltage of net56.

## Plotting and Printing Commands

```
ocnPrint( ?output "myFile" v( "net55" ) )
=> t.
```

Prints the text for the specified waveform to a file named myFile.

```
ocnPrint( ?output "./myOutputFile" v("net1") ?from 0 ?to 0.5n ?step 0.1n )
```

Prints the text for the specified waveform from 0 to 0 . 5n on the x axis in the incremental steps of 0 . 1n.

## Plotting and Printing Commands

# ocnPrintTMIReliabilityResults

```
ocnPrintTMIReliabilityResults(
    t_psfdir
    [ ?resultType t_resultType ]
    [ ?threshold t_threshold ]
    [ ?exportFile t_exportFilePath ]
    [ ?sortRule t_sortRule ]
    ) t / nil
```

# **Description**

Prints simulation results for reliability analysis run in the OCEAN mode.

# **Arguments**

t_psfdir	Path to the PSF directory in which simulation results of reliability analysis for the test are saved.
?resultType t_resultType	Name of the result type for which you need to view results.
	Default value: nil
?threshold t_threshold	Threshold value. A result value greater than this threshold will be printed. Other values are ignored.
	Default value: nil
<pre>?exportFile t_exportFilePath</pre>	Path to the file where you need to export the printed results in CSV format. If you do not specify this path, the results are displayed on the terminal.
	Default value: nil
?sortRule t_sortRule	Rule to identify the column to be used for ranking of data.
	Default value: nil

### Value Returned

t	When the results of reliability simulation are successfully printed.
nil	Returns nil otherwise.

## Plotting and Printing Commands

### **Examples**

### Example 1

The following example code prints the reliability results from the given psf directory on the terminal:

```
psfdir = "./simulation/bertlink/test/maestro/results/maestro/Interactive.0/3/
bertlink:osc13:1/psf"
=> "./simulation/bertlink/test/maestro/results/maestro/Interactive.0/3/
bertlink:osc13:1/psf"
ocnPrintTMIReliabilityResults(psfdir)
=> t
```

### Example 2

The following example code checks the available result types in the reliability results and exports those to a file:

```
psfdir = "./simulation/bertlink/test/maestro/results/maestro/Interactive.0/3/
bertlink:osc13:1/psf"
=> "./simulation/bertlink/test/maestro/results/maestro/Interactive.0/3/
bertlink:osc13:1/psf"

ocnPrintTMIResultTypeList(psfdir)
=> ("DeltaTemp" "didlinBti" "didlinHci" "didlinHciBti" "didsatBti"
   "didsatHci" "didsatHciBti" "dvtlinBti" "dvtlinHci" "dvtlinHciBti"
   "lifetimeBti" "lifetimeHci" "lifetimeHciBti"
)
ocnPrintTMIReliabilityResults(psfdir ?resultType "DeltaTemp" ?threshold "3"
?exportFile "./TMIresults.csv")
; prints results on the terminal
=> t
```

# Example 3

The following example code sorts the results on the basis of column didlinHciBti:

```
ocnPrintTMIReliabilityResults(psfdir ?sortRule "didlinHciBti" ?exportFile "./1")
Rank InstanceName Model DeltaTemp didsatHciBti didlinHciBti dvtlinHciBti
                                     8.90E+00
                                               4.69E+00
  IL6.M1
          pch svt mac.12
                           3.68E+01
                                                          3.34E-02
          pch svt mac.12
  IL5.M1
                           3.81E+01
                                     8.11E+00
                                               4.27E+00
                                                          3.04E-02
  IL4.M1 pch svt mac.12
                          3.81E+01 8.11E+00 4.27E+00
                                                         3.04E-02
  IL7.M1 pch svt mac.12 2.18E+01 8.04E+00 4.24E+00
                                                         3.02E-02
  I1.M1 pch_svt_mac.12 1.25E+01 7.16E+00 3.77E+00 2.69E-02
  IL1.M1 pch svt mac.12 3.67E+01 6.49E+00 3.42E+00 2.44E-02
  IL3.M1 pch_svt_mac.12 3.67E+01 6.49E+00 3.42E+00 2.44E-02
  IL2.M1 pch_svt_mac.12 3.67E+01 6.49E+00 3.42E+00 2.44E-02 I4.M1 pch_svt_mac.12 7.13E+00 6.32E+00 3.33E+00 2.37E-02
7
  I9.M1 pch svt mac.12 5.45E+00 6.21E+00 3.27E+00 2.33E-02
=> t
;; note the ranking done in the results printed above.
```

### Plotting and Printing Commands

# ocnPrintTMIResultTypeList

## **Description**

Returns a list of valid result types.

## **Arguments**

t psfdir

?netlistFilePreName

t prefixOfNetlistFile

Path to the PSF directory in which simulation results of reliability analysis for the test are saved.

Prefix of the name of the netlist file. The function uses this prefix to search for the relevant mapping file saved by simulator.

For example, if the specified prefix is "input", the netlist name is "input.scs", and the name of mapping file is ".input\_tmi\_deg.mapping". If the prefix is "amsControlSpectre", the netlist file is "amsControlSpectre.scs", and the name of mapping file is

".amsControlSpectre\_tmi\_deg.mapping".

Default value: "input"

#### Value Returned

1\_resultTypesnilList of result types in the simulation results.Returns nil otherwise.

### **Example**

The following example code checks the available result types in the reliability results and exports those to a file:

```
psfdir = "./simulation/bertlink/test/maestro/results/maestro/Interactive.0/3/
bertlink:osc13:1/psf"
=> "./simulation/bertlink/test/maestro/results/maestro/Interactive.0/3/
bertlink:osc13:1/psf"
```

## Plotting and Printing Commands

```
ocnPrintTMIResultTypeList(psfdir)
=> ("DeltaTemp" "didlinBti" "didlinHci" "didlinHciBti" "didsatBti"
  "didsatHci" "didsatHciBti" "dvtlinBti" "dvtlinHci" "dvtlinHciBti"
  "lifetimeBti" "lifetimeHci" "lifetimeHciBti"
)
ocnPrintTMIReliabilityResults(psfdir ?resultType "DeltaTemp" ?threshold "3"
?exportFile "./TMIresults.csv")
=> t
```

## Plotting and Printing Commands

# ocnSetAttrib

```
ocnSetAttrib(
   [?XAxisLabel xLabel]
   [?YAxisLabel yLabel]
   [?XScale xscale]
   [?YScale yscale]
   [?XLimit xlimit]
   [?YLimit ylimit]
   [?YRange yrange]
   [?Origin origin]
)
   => t / nil
```

# **Description**

Sets the waveform window plotting attributes.

# Plotting and Printing Commands

# **Arguments**

?XAxisLabel xLabel	Label (symbol or string) for the X axis in the waveform window.
?YAxisLabel yLabel	Label (symbol or string) for the Y axis associated with the $stripNumber$ in the waveform window.
?XScale xScale	Scale of the X axis in the waveform window.
	Valid values (symbols): 'auto, 'log, and 'linear
?YScale <i>yScale</i>	Scale of the Y axis associated with the $stripNumber$ in the waveform window.
	Valid values (symbols): 'log and 'linear
?XLimit <i>xLimit</i>	Displays limits of the X axis in the waveform window.
	Valid values: List of two numbers or 'auto (symbol).
	The first number in the list indicates the minimum limit and the second indicates the maximum limit.
	' auto sets the limit to autoscale.
?YLimit <i>yLimit</i>	Displays limits of the Y axis associated with the $stripNumber$ in the waveform window.
	Valid values: List of two numbers or 'auto (symbol).
	The first number in the list indicates the minimum limit and the second indicates the maximum limit.
	' auto sets the limit to autoscale.
?YRange <i>yRange</i>	Y range (integer) of the waveforms associated with the $stripNumber$ in the waveform window.
?Origin origin	Axes origin of the waveform window.
	Valid values: List of two numbers.

**Note:** The valid range for *stripNumber* is 1-20.

### Plotting and Printing Commands

### Value Returned

t Returns t if the valus of all arguments are set

successfully.

nil Returns nil if one or more arguments fail to set as

specified.

## **Example**

```
ocnSetAttrib( ?XAxisLabel 'XMylabel ?YAxisLabel 'YMyLabelt ?stripNumber 2
     )
=> t
```

Sets the X and Y axis labels to XMylabel and YMyLabel, respectively.

```
ocnSetAttrib(?XScale 'log ?YScale 'linear ?stripNumber 2 )
=> t
```

Sets the scale of X and Y axis to log and linear, respectively.

```
ocnSetAttrib(?XScale 'auto ?XLimit '(3 7) ?YLimit 'auto ?stripNumber 2 )
=> t
```

Sets the scale of X axis to autoscale. Sets the Y display limits to autoscale.

### Plotting and Printing Commands

# ocnWriteLsspToFile

```
ocnWriteLsspToFile(
    filename t_filename
    net1 input_node_name
    term1 input_src_terminal
    net2 output_node_name
    term2 output_src_terminal
    [?format t_format]
    [?datafmt t_data_format]
    [?port1 port1_name]
    [?port2 port2_name]
    [?result1 result1_name]
    [?result2 result2_name]
)
=> nil
```

# **Description**

Writes the large signal S-Parameter results to a file in Touchstone or Spectre format.

### Plotting and Printing Commands

### **Arguments**

t filename Name of the file in which results are to be written.

input\_node\_name
Name of the input node.

input src terminal Name of the input source terminal.

output node name Name of the output node.

output src terminal Name of the output source terminal.

?format t format Format of file in which results are to be written.

Possible values: touchstone, spectre

Default value: 'touchstone

?data fmt t data format Format of data being written.

Possible values: magphase, dbphase, realimag

Default value: realimag

?port1 port1\_name Name of the first port.

Default value: 50

?port2 port2 name Name of the second port.

Default value: 50

?result1 result1 name Name of the first pss result.

Default value: sweeplssp1 lssp1 fd-sweep

?result2 result2\_name Name of the second pss result.

Default value: sweeplssp2 lssp2 fd-sweep

### **Value Returned**

t Specifies that the results are written to the specified

file successfully.

nil Returns nil if the results are not written.

### Example

ocnWriteLsspToFile "lssp.sp2" "/net026" "/PORT1/PLUS" "/RFOUT"
"/PORT2/PLUS" ?format "touchstone" ?datafmt "realimag" ?port1 50 ?port2 50
?result1 "sweeplssp1 lssp1 fd-sweep" ?result2 "sweeplssp2 lssp2 fd-sweep")

### Plotting and Printing Commands

# ocnYvsYplot

```
ocnYvsYplot(
    [ ?wavex o_wavex ?wavey o_wavey ]
    [ ?exprx o_exprx ?expry o_expry ]
    [ ?titleList l_titleList ]
    [ ?colorList l_colorList ]
    )
    => wave / nil
```

# **Description**

Plots a wave against another wave or an expression against another expression.

This is currently supported for a family of waveforms generated from simple parametric simulation results data. It is not supported for a family of waveforms generated from parametric simulation with parametr, Corners or MonteCarlo results data.

## Plotting and Printing Commands

# **Arguments**

?wavex o_wavex	Reference wave against which the wave provided needs to be plotted.
?wavey o_wavey	Wave to be plotted against the reference wave.
?exprx o_exprx	Reference expression against which the expression provided needs to be plotted.
?expry o_expry	Expression to be plotted against the reference expression.
?titleList l_titleList	List of waveform titles. If the waveform is simple, only one label will be required. If the waveform is param, a list of labels needs to be provided.
?colorList l_colorList	List specifying the colors for the waveforms. If you do not supply this argument, the default colors are used. The colors that are available are defined in your technology file.
	Valid Values: "y1" through "y66".

### Value Returned

wave Returns the waveform specified.

nil Returns nil if the plot could not be generated.

### Example

```
wy = VT("/vout")
wx = VT("/vin")
ex = "VT('/vin')"
ey = "VT('/vout')"
ocnYvsYplot(?wavex wx ?wavey wy ?titleList '("simpleWave") ?colorList '(y1))
```

Plots wave wy against wave wx with the title being simpleWave and the color being y1.

```
ocnYvsYplot(?exprx ex ?expry ey ?titleList '("simpleWave") ?colorList '(y2))
```

Plots expression ey against expression ex with the title being simpleWave and the color being y2.

### Plotting and Printing Commands

# plot

```
plot(
    o_waveform1 [ o_waveform2 ... ]
    [ ?expr l_exprList ]
    [ ?strip x_stripNumber ]
    )
    => t / nil
```

## **Description**

Plots waveforms in the current subwindow. If there is no Virtuoso Visualization and Analysis XL window, this command opens one.

**Note:** plot is implemented as a macro and not as a SKILL function. Therefore, the functions that expect a function name as an argument will not accept plot as a valid argument. For example, the following call to the function apply is not valid:

```
apply('plot)
```

## **Arguments**

o_waveform1	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave:XXXXX.)
o_waveform2	Additional waveform object.
?expr l_exprList	List of strings used to give names to the waveform objects.
?strip x_stripNumber	An integer using which you can plot waveforms selectively on different strips and subwindows. If you specify an integer, it is used as the strip for all waveforms. To use the strip option for multiple waveforms, you can specify a list of strip numbers

## Plotting and Printing Commands

#### Value Returned

t Returns t if the waveforms are plotted.

nil Returns nil and an error message if the waveforms

cannot be plotted.

### **Additional Information**

Following are the scenarios that show how the plot and displayMode functions work together:

**Case 1**: When no waveform plot is open and you plot a waveform, w1, and then plot another waveform, w2, both the waveforms are plotted in one strip. Now if you set the displayMode ('strip) function, the waveforms are plotted in two different strips.

- **1.** plot (w1)
- **2.** plot (w2)

w1 and w2 are plotted in one strip.

**3.** displayMode('strip)

w1 and w2 are plotted in two strips.

**4.** plot (w3)

w3 is plotted in a new strip. The ?strip argument is not required in this case.

Case 2: When no waveform plot is open and you set displayMode ('strip):

- displayMode('strip)
- **2.** plot (w1)
- **3.** plot (w2)

w2 is plotted in a new strip. The explicit ?strip argument is not required in this case.

**Case 3**: When no waveform plot is open:

- **1.** plot (w1)
- **2.** plot (w2)

w1 and w2 are plotted in one strip.

**3.** plot(w3 ?strip 2)

### Plotting and Printing Commands

w3 is plotted in a new strip because the plot function contains the ?strip 2 argument.

**4.** plot (w4)

w4 is plotted in the same strip as in which w3 is plotted.

### Case 4:When no waveform plot is open:

- **1.** plot(w1 ?strip 1)
- **2.** plot(w2 ?strip 2)

w2 is plotted in strip 2.

**3.** plot(w3 ?strip 1)

w3 is plotted in strip 1.

**4.** plot(w4 ?strip 2)

w4 is plotted in strip 2, which now becomes an active strip.

**5.** displayMode('strip)

This divides the traces contained in strip 2 into individual strips, because strip 2 is the active strip now. Now, if you plot new waveforms in this strip, they are plotted in new strips. However, strip 1continues to have two signals, w1 and w3.

**6.** plot (w5)

w5 is plotted in a new strip.

**7.** plot (w6)

w6 is plotted in a new strip.

### Case 5: When no waveform plot is open:

- 1. window=awvCreatePlotWindow()
- **2.** awvSetDisplayMode(window "strip")
- **3.** plot (w1)
- **4.** plot (w2)
- **5.** plot (w3)
- **6.** plot (w4)

w1, w2, w3, and w4 are plotted in four different strips.

### Plotting and Printing Commands

**Case 6**: Plot digital and analog data and set <code>displayMode(`composite)</code>. This combines all analog signals into a single strip. Now, if you set the <code>displayMode(`strip)</code>, the analog signals are divided into individual strips. These operations are applicable only on the active strip.

### **Example**

```
plot(v( "/net56" ) )
```

Plots the waveform for the voltage of net56.

```
plot( vm( "/net56" ) vp( "/net56" ) )
```

Plots the waveforms for the magnitude of the voltage of net56 and the phase of the voltage of net56.

```
plot( v( "OUT" ) i( "VFB" ) ?expr list( "voltage" "current" ) )
```

Plots the waveforms, but changes one legend label from v("OUT") to voltage and changes the other legend label from i("VFB") to current.

```
plot( v( "OUT" ) i( "VFB" ) )
```

Plots the waveforms v ( "OUT" ) and i ( "VFB" ) on the Y axes 1 and 2, respectively. plot (wave1 wave2 wave3 ?strip list(1 2 2))

Plots wave1 to strip 1, and wave2 and wave3 to strip 2.

## Plotting and Printing Commands

# plotStyle

```
plotStyle(
     S_style
)
     => t / nil
```

# **Description**

Sets the plotting style for all the waveforms in the current subwindow.

If the plotting style is bar and the display mode is smith, the plotting style is ignored until the display mode is set to strip or composite.

# Plotting and Printing Commands

# **Arguments**

 $S_style$ 

Plotting style for the subwindow. Valid values: auto, scatterplot, bar, joined

Argument	Description
auto	The appropriate plotting style is automatically chosen.
scatterplot	Data points are not joined.
bar	Vertical bars are drawn at each data point that extend from the point to the bottom of the graph.
joined	Each data point is joined to adjacent data points by straight-line segments.

#### **Value Returned**

t

Returns t if the plotting style is set.

nil

Returns nil and an error message if the plotting style is not set.

# **Example**

Sets the plot style to auto.

## Plotting and Printing Commands

# printGraph

```
printGraph(
     [ ?window x window ]
     [ ?printerName s hcPrinterName ]
     [ ?horizontalMargin x horizontalMargin ]
     [ ?verticalMargin x verticalMargin ]
     [ ?numCopy x numCopy ]
     [ ?paperSize x paperSize ]
     [ ?orientation s_orientation ]
     [ ?fileName s fileName ]
     [ ?tempDir s tempDir ]
     [ ?matchWindow g matchWindow ]
     [ ?numGraphsPerPage x numGraphsPerPage ]
     [ ?printMarkerTable g_printMarkerTable ]
     [ ?markerTableLocation s markerTableLocation ]
     [ ?enableHeader g enableHeader ]
     [ ?enableFooter g enableFooter ]
     [ ?headerLeftText s_headerLeftText ]
     [ ?headerCenterText s headerCenterText ]
     [ ?headerRightText s headerRightText ]
     [ ?footerLeftText s_footerLeftText ]
     [ ?footerCenterText s footerCenterText ]
     [ ?footerRightText s footerRightText ]
     [ ?printColor g printColor ]
     [ ?doubleSidedPrint g doubleSidedPrint ]
     [ ?duplexMode s duplexMode ]
     [ ?pageOrder s pageOrder ]
     => t / nil
```

# Description

Prints the graph plotted in the specified window.

## **Arguments**

?window X\_window Window ID of the waveform window whose plot is to be sent to a printer or a file. The default value is the window ID of the current window.

```
?printerName s_printerName
Name of the printer to be used for printing.
Valid values: a string or nil
Default value: nil
?horizontalMargin x_horizontalMargin
Horizontal margin.
Valid values: any positive integer
```

## Plotting and Printing Commands

Default value: 18

?verticalMargin x verticalMargin

Vertical margin

Valid values: any positive integer

Default value: 18

?numCopy x numCopyNumber of copies to be printed.

Valid values: any positive integer

Default value: 1

?paperSize x paperSize

Size of paper used for printing.

Valid values: letter, legal, executive, folio, ledger, tabloid, a0, a1, a2, a3, a4, a5, a6, a7, a8, a9, b0, b1, b2, b3,b4, b5, b6, b7, b8, b9, b10, c5e, comm10e, and dle.

Default value: letter

?orientation s orientation

Paper orientation. This option works only when you print a graph window and does not work when you save the window to a graph file.

Valid values: potrait, landscape, and automatic

Default value: landscape

?fileName s fileName

Name of the output file. The output file can be created in one of the following file formats:

PS - PostScript (.ps)

PDF - Portable Document Format (.pdf)

Valid values: any string value or nil

Default value: nil

?tempDir  $s_tempDirName$  of a temporary directory to be used for scratch space.

Valid values: name of a temporary directory

Default value: "/usr/tmp"

?matchWindow g matchWindow

Specifies whether the print output is identical to the current graph window. This option is used if you want to print all the subwindows in a PDF file in the same order in which they are arranged in the graph.

Valid values: t or nil Default value: nil

?numGraphsPerPage x\_numGraphsPerPage Specifies how many graphs are to be printed per page. Valid values: Integer values 1, 2, 3, 4, 8, 12, 16, 20

Default value: 1

?printMarkerTable g printMarkerTable

## Plotting and Printing Commands

Specifies whether the marker table is to be printed.

Valid values: t or nil.

Default value: nil.

?MarkerTableLocation s MarkerTableLocation

Specifies the location of the marker table on the page to be printed.

Valid values: belowGraph and separatePage.

Default value: belowGraph.

?enableHeader g\_enableHeader

Specifies whether the page contains a header.

Valid values: t or nil

Default value: t

?enableFooter g enableFooter

Specifies whether the page contains a footer.

Valid values: t or nil

Default value: t

 $? header Left Text \ s\_header Left Text \\$ 

Sets the text to be printed to the left of header.

Valid values: any string value

Default value: " "

?headerCenterText s headerCenterText

Sets the text to be printed in the center of the header.

Valid values: Any string value and the following macros—\$TOTALPAGES, \$TITLE,

\$USERID, \$PRINTER, \$PAGE, \$DATE, \$DATETIME, \$AUTHOR, \$TIME.

Default value: \$TITLE

 $\verb|?headerRightText| s_headerRightText|$ 

Sets the text to be printed to the right of the header.

Valid values: Any string value and the following macros—\$TOTALPAGES, \$TITLE,

\$USERID, \$PRINTER, \$PAGE, \$DATE, \$DATETIME, \$AUTHOR, \$TIME.

Default value: \$DATETIME

?footerLeftText s footerLeftText

Sets the text to be printed to the left of footer.

Valid values: Any string value and the following macros—\$TOTALPAGES, \$TITLE,

\$USERID, \$PRINTER, \$PAGE, \$DATE, \$DATETIME, \$AUTHOR, \$TIME.

Default value: Printed on \$PRINTER by \$USERID

?footerCenterText s footerCenterText

Sets the text to be printed in the center of the footer.

Valid values: Any string value and the following macros—\$TOTALPAGES, \$TITLE,

SUSERID, SPRINTER, SPAGE, SDATE, SDATETIME, SAUTHOR, STIME.

Default value: " "

#### Plotting and Printing Commands

?footerRightText  $s_footerRightText$ 

Sets the text to be printed on the right of the footer.

Valid values: Any string value and the following macros—\$TOTALPAGES, \$TITLE,

\$USERID, \$PRINTER, \$PAGE, \$DATE, \$DATETIME, \$AUTHOR, \$TIME.

Default value: Page \$PAGE of \$TOTALPAGES

?printColor g printColor

Specifies whether the print is to be colored

Valid values: t or nil

Default value: t

?doubleSidedPrint g doubleSidedPrint

Specifies whether both the sides of paper is used for printing.

Valid values: t or nil

?duplexMode s\_duplexMode Specifies the duplex printing mode.

Valid values: none, auto, shortSide, longSide.

Default value: auto

?pageOrder s pageOrder

Specifies the order in which pages are printed.

Valid values: collate and reverse

Default value: collate

#### Value Returned

t Returns t if the

function runs successfully.

nil Returns nil if

there is an error.

#### **Examples**

printGraph()

Prints the current graph window with the default printing options.

printGraph(?printerName "ind001" ?paperSize "a4" ?orientation
'portrait)

# Plotting and Printing Commands

Prints the current graph window by using the printer, ind001, with paper size a4 and orientation portrait.

#### Plotting and Printing Commands

# pzFrequencyAndRealFilter

```
pzFrequencyAndRealFilter(
    o_wave
    [?freqfilter f_fval]
    [?realfilter f_rval])
=> o waveform / nil
```

# **Description**

Returns a filtered Pole or Zero waveform from the pole zero simulation data. Filtering is done on the basis of given maximum frequency and minimum real value.

**Note:** This command also works for the parametric or sweep data.

# **Arguments**

o_wave	Input Pole or Zero waveform (complex points) from the simulation data of PZ analysis.
?freqfFilter f_val	Maximum pole and zero frequency value to filter out poles and zeros that are outside the frequency band of interest (FBOI) and that do not influence the transfer function in the FBOI.
?realFilter <i>f_rval</i>	Minimum real value which is used to filter out poles and zeros whose real value are less than or equal to the value specified.

#### **Values Returned**

o_waveform	Returns a Pole or Zero waveform.
nil	Returns nil if there is an error.

#### **Examples**

```
pzFrequencyAndRealFilter(wave ?freqfilter 1e+24 ?realfilter 2e+10)
=> srrWave:175051584
```

Returns a filtered Pole or Zero waveform, which is filtered on the basis of given maximum frequency and minimum real value.

## Plotting and Printing Commands

# pzPlot

```
pzPlot(
    [ ?resultsDir t_resultsDir ]
    [ ?result S_resultName ]
    [ ?plot S_toPlot ]
    [ ?freqfilter f_fval ]
    [ ?realfilter f_rval ])
    => t / nil
```

#### **Description**

Plots a report showing the poles and zeros of the network. If you specify a directory with resultsDir, the pzPlot command plots the results for that directory. The  $S_toPlot$  option can be used to plot only poles, only zeros or both poles and zeros information.

This command should be run on the results of the Spectre pz (pole-zero) analysis.

**Note:** This command also works for the parametric or sweep data.

## **Arguments**

- t\_resultsDirDirectory containing the results. If you specify a directory with resultsDir, the pzPlot command plots the results for that directory.
- S resultNamePointer to results from the analysis for which you want to plot the report.
- $S_toPlot$ Use this option to plot only poles, only zeros or both poles and zeros information. Valid values: 'poles, 'zeros, 'polesZeros.
- $f_{Val}$  Maximum pole and zero frequency value to filter out poles and zeros that are outside the frequency band of interest (FBOI) and that do not influence the transfer function in the FBOI.
- $f_rval$ Real value which is used to filter out poles and zeros whose real value are less than or equal to the value specified.

#### Value Returned

t Returns t if it plots a report.

nilReturns nil otherwise.

#### **Example**

```
pzPlot(?resultsDir "/usr/simulation/lowpass/spectre/schematic" ?result 'pz)
```

#### Plotting and Printing Commands

Plots a report for all the poles and zeros for the specified results.

```
pzPlot(?resultsDir "/usr/simulation/lowpass/spectre/schematic" ?plot 'poles)
```

Plots a report containing only poles for the specified results.

```
pzPlot( ?plot 'zeros ?realfilter -1.69e-01)
```

Plots a report for all those zeros whose real values are greater than the real value specified.

```
pzPlot( ?plot 'polesZeros ?freqfilter 2.6e-01 )
```

Plots a report for all those poles and zeros whose frequency is within the frequency band of interest (2.6e-01).

#### Plotting and Printing Commands

# pzSummary

```
pzSummary(
    [?resultsDir t_resultsDir]
    [?result S_resultName]
    [?print S_toPrint]
    [?freqfilter f_fval]
    [?realfilter f_rval]
    [?output t_output]
)
    => t / nil
```

#### **Description**

Prints a report with the poles and zeros of the network. If you specify a directory with resultsDir, the pzSummary command prints the results for that directory. Use the  $S\_toPrint$  option to print only poles, only zeros or both poles and zeros information.

This command should be run on the results of the Spectre pz (pole-zero) analysis.

**Note:** This command also works for the parametric or sweep data.

## **Arguments**

- t\_resultsDirDirectory containing the results. If you specify a directory with resultsDir, the pzSummary command plots the results for that directory.
- S\_resultNamePointer to results from the analysis for which you want to print the report.
- $S_toPlot$ Use this option to plot only poles, only zeros or both poles and zeros information. Valid values: 'poles, 'zeros, 'polesZeros.
- $f_{val}$  Maximum pole and zero frequency value to filter out poles and zeros that are outside the frequency band of interest (FBOI) and that do not influence the transfer function in the FBOI.
- $f_rval$ Real value which is used to filter out poles and zeros whose real value are less than or equal to the value specified.
- $t\_output$ Provides an option to write the output to a file. The possible values can be a file name or a port name.

#### Value Returned

t Returns t if it prints a report.

nilReturns nil otherwise.

## Plotting and Printing Commands

# **Example**

```
pzSummary(?resultsDir "/usr/simulation/lowpass/spectre/schematic" ?result 'pz)
```

Prints a report for all the poles and zeros for the specified results.

```
pzSummary(?resultsDir "/usr/simulation/lowpass/spectre/schematic" ?print 'poles)
```

Prints a report containing only poles for the specified results.

```
pzSummary( ?print 'zeros ?realfilter -1.69e-01)
```

Prints a report for all those zeros whose real values are less than or equal to the real value specified.

```
pzSummary( ?print 'polesZeros ?freqfilter 2.6e-01 )
```

Prints a report for all those poles and zeros whose frequency is within the frequency band of interest (2.6e-01).

```
pzSummary( ?output "/tmp/file")
```

Prints results in the file.

```
pzSummary( ?output "file")
```

Prints results in a file located in the current working directory.

```
pzSummary( ?output oFile)
where, oFile=outfile("/tmp/file")
```

#### Prints to the opened file

```
pzSummary( ?output nil)
pzSummary( ?output t)
pzSummary( ?output 32)
```

Prints results on the CIW or Ocean command-line.

# Plotting and Printing Commands

#### removeLabel

# **Description**

Removes the label, or all the labels identified in a list, from the current subwindow.

#### **Arguments**

1\_id List of labels to remove.

# **Value Returned**

t Returns t when the label or labels are removed.

nil Returns nil if there is an error.

### **Example**

```
label = addWindowLabel( list( 0.75 0.75 ) "test" )
```

Adds the "test" label to the current subwindow at the specified coordinates and stores the label identification number in label.

```
removeLabel( label )
```

Removes the label whose identification number is stored in label. In this case, the "test" label is removed.

#### Plotting and Printing Commands

#### report

```
report(
    [ ?output t_filename | p_port ]
    [ ?type t_type ]
    [ ?name t_name ]
    [ ?param t_param ]
    [ ?format s_reportStyle ]
    [ ?report s_reportStyle ]
    [ ?maxLineWidth charsPerLine ]
    )
    => t / nil
```

# **Description**

Prints a report of the information contained in an analysis previously specified with selectResult.

You can use this command to print operating-point, model, or component information. If you provide a filename as the <code>?output</code> argument, the <code>report</code> command opens the file and writes the information to it. If you provide a port (the return value of the SKILL <code>outfile</code> command), the <code>report</code> command appends the information to the file that is represented by the port.

**Note:** You can use the dataTypes command to see what types of reports you can choose. For Spectre® circuit simulator operating points, be sure to choose dcOpInfo.

#### **Arguments**

t\_filenameFile in which to write the information. The report command opens the file, writes to the file, and closes the file. If you specify the filename without a path, the OCEAN environment creates the file in the directory pointed to by your SKILL Path. To find out what your SKILL path is, type getSkillPath() at the OCEAN prompt.

 $p\_port$ Port (previously opened with outfile) through which to append the information to a file. You are responsible for closing the port. See the <u>outfile</u> command for more information.

t typeType of information to print, such as all bits.

t nameName of the node or component.

t paramName of the parameter to print. It is also a list.

s reportStyleSpecifies the format of the output.

Valid values: spice and paramValPair

Default value: paramValPair

#### Plotting and Printing Commands

The spice format looks like this:

	Param1	Param2	Param3
Name1	value	value	value
Name2	value	value	value
Name3	value	value	value

The paramValPair format looks like this:

Name1

Param1=value Param2=value Param3=value

Name2

Param1=value Param2=value Param3=value

Name3

Param1=value Param2=value Param3=value

charsPerLineNumber of characters to be printed per line.

#### Value Returned

t Returns t if the information is printed.

nilReturns nil and an error message if the information cannot be printed.

#### **Example**

The following example shows how to display a report by using the results of an analysis already run. First, run the results() command to get a list of the type of results that exist in the current results directory.

```
results()
= > ( dcOpInfo tran ac dc)
```

From the list of result types returned by the previous function, select a particular type of results for which you want to print the report.

```
selectResult( 'dcOpInfo )
= > t
```

Use the report function to print the results. The following examples show how to print different details in a report:

```
report() = > t
```

#### Plotting and Printing Commands

Prints all the operating-point parameters.

```
report( ?type "bjt" )
= > t
```

Prints all the bjt operating-point parameters.

```
report( ?type "bjt" ?param "ib" )
= > t
```

Prints the ib parameter for all bits.

```
report( ?type "bjt" ?name "/Q1" ?param "ib" )
= > t.
```

Prints the ib parameter for the bjt named Q1.

```
report( ?output "myFile" )
=> t
```

Prints all the operating-point parameters to a file named myFile.

```
report( ?output myAlreadyOpenedPort )
=> t
```

Prints all the operating-point parameters to a port named myAlreadyOpenedPort.

The report () can also be used by providing the set of parameters as a list as follows:

```
Type : bsim3v3
Params : cdg cgb gm ids
report(?type "bsim3v3" ?param "cdg")
report(?type "bsim3v3" ?param '( "cdg" "cgb"))
report(?type "bsim3v3" ?param '( "cdg" "cgb" "gm" "ids"))
report( ?format 'spice ?maxLineWidth 200 )
=> t
```

Prints the report in spice format wrapping at column 200.

#### Plotting and Printing Commands

# saveGraphImage

```
saveGraphImage(
     [ ?window x_window ]
     [ ?fileName x_fileName ]
     [ ?exactCopy g exactCopy ]
     [ ?quality x quality ]
     [ ?msOptimize g_msOptimize ]
     [ ?width x width ]
     [ ?height x_height ]
     [ ?units s_units ]
     [ ?resolution x resolution ]
     [ ?resolutionUnits s resolutionUnits ]
     [ ?aspectRatio g_aspectRatio ]
     [ ?enableTitle g_enableTitle ]
     [ ?enableLegend g_enableLegend ]
     [ ?enableAxes g_enableAxes ]
     [ ?enableGrids g enableGrids ]
     [ ?backgroundColor s_backgroundColor ]
     [ ?saveAllSubwindows g saveAllSubwindows ]
     [ ?saveEachSubwindowSeparately g\_saveEachSubwindowSeparately ]
    => x fileName / nil
```

# **Description**

Saves the graph as an image.

#### Plotting and Printing Commands

#### **Arguments**

?window x\_window

Window ID of the waveform window whose plot is to be saved in a file. The default value is the window ID of the current window.

?fileName x fileName

Name of the output file to be created. The output file can be created in one of the following file formats:

- BMP Windows Device Independent Bitmap (.bmp)
- PNG Portable Network Graphics(.png)
- TIFF Tagged Image File Format (.tiff)
- EPS Encapsulated Post Script (.eps)
- PDF Portable Document Format (.pdf)
- PPM Portable PixMap File (.ppm)
- JPG Joint Photographic Experts Group (.jpg)
- SVG Scalable Vector Graphics (.svg)
- XPM X PixMap (.xpm)Valid values: any string value or nil

Default value: nil.

**Note:** If *fileName* argument is not specified, the graph image is saved in a image.png file.

?exactCopy g exactCopy

Saves the exact copy of all subwindows. Only <code>?quality</code> and <code>?fileName</code> arguments work with this option. This option does not work for the <code>eps</code> file format.

Valid values: t or nil

Default value: nil

?quality x quality

Modifies the quality of the image. This option works only for the .jpeg file format. This option does not work for the eps file format.

Valid values: 20 to 100%

Default value: 85%

#### Plotting and Printing Commands

?msOptimize  $g_{msOptimize}$  Enables the image to be imported in the Microsoft

office application. This option is available when you select the image type as Encapsulated PostScript (\*.eps). This option simplifies the image output so that it can be ready by Microsoft Office 2003 and

2007 applications

Valid values: t or nil

Default value: t

?width x width Sets the width of the image.

Valid values: Any positive integer value.

Default value: 800 pixels for bmp, png, tiff, ppm and xpm file formats and 8.33 inches for pdf, svg

and eps file formats.

?height x height Sets the height of the imageValid values: Any positive

integer value.

Default value: 600 pixels for bmp, png, tiff, ppm and xpm file formats and 6.25 inches for pdf, svg

and eps file formats.

?units s units Specifies the unit for image size (height and

width) Valid values: inch, cm, mm, picas, pixels,

and points

Default value: pixels for bmp, png, tiff, ppm file formats and xpm and inch for pdf, svg and eps file

formats.

?resolution x resolution Sets the image resolution. This option works only for

the bmp, jpeg, png, ppm, tiff, and xpm file formats. It does not work for eps, pdf, and svg file formats.

Valid values: Any positive integer value.

Default value: 96

?resolutionUnits s resolutionUnits

#### Plotting and Printing Commands

Sets the units for image resolution. This option works only for the bmp, jpeg, png, ppm, tif, and xpm file formats. It does not work for eps, pdf, and svg file formats.

Valid values: pixels/cm and pixels/in

Default value: pixels/in

?aspectRatio g aspectRatio

Enables the aspect ratio, which is the ratio of the width of the image to its height.

Valid values: t or nil

Default value: nil

?enableTitle g enableTitle

Displays the trace title in the graph image.

Valid values: t or nil

Default value: t

?enableLegend g\_enableLegend

Displays the trace legend in the graph image.

Valid values: t or nil

Default value: t

?enableAxes q enableAxes

Displays the axes in the graph image.

Valid values: t or nil

Default value: t

?enableGrids g enableGrids

Displays the grids in the graph image.

Valid values: t or nil

Default value: t

**Note:** You cannot set this argument to t if ?enableAxes is set to nil. If you want to display the grids in the graph image, ensure that ?enableAxes is set to t.

#### Plotting and Printing Commands

?backgroundColor g backgroundColor

Specify the background color.

Default value: nil, which means graph image is saved with the current background colorV

alid values: All the valid color values are defined at the following location: <a href="http://www.w3.org/TR/SVG/types.html#ColorKeywords">http://www.w3.org/TR/SVG/types.html#ColorKeywords</a>

For example, red, blue, green, black, white, gray, cyan, magenta, yellow, and so on.

?saveAllSubwindows g saveAllSubwindows

Saves all subwindows or the current subwindow.

Default value: t

Valid values: t (current window is saved) or nil (all windows are saved)

?saveEachSubwindowSeparately g saveEachSubwindowSeparately

Specifies whether to save each subwindow in a separate image file or in the same image file.

Valid values: t or nil

Default value: t

#### Value Returned

 $x_fileName$  Returns the name of the output file.

nil Returns nil if there is an error.

#### **Examples**

■ saveGraphImage()

Saves the current graph window with the default saving options.

saveGraphImage(?fileName "ViVA.jpg" ?enableTitle t ?enableLegend
nil)

Saves the current graph window in the ViVA. jpg file with only trace legend enabled.

#### Plotting and Printing Commands

saveGraphImage(?window currentWindow() ?fileName "ViVA.jpg"
?backgroundColor "light gray")

Saves the current graph window in the ViVA. jpg file with background color as light gray.

#### **Additional Information**

Following are the guidelines supported by the saveGraphImage function:

- Arguments exactCopy, quality, resolution, and resolutionUnits are ignored for eps file format.
- Only fileName and quality arguments can be used with exactCopy argument. All other arguments are ignored.
- Argument *quality* can be used only with jpeg file format. It is ignored for other formats.
- Arguments resolution and resolutionUnits cannot be set for eps, pdf, and svg file formats.
- Argument msOptimize can be set to nil only for eps file format.
- Argument enableGrids cannot be set to true when enableAxes is nil.

## Plotting and Printing Commands

# **xLimit**

# **Description**

Sets the X axis display limits for the current subwindow. This command does not take effect if the display mode is set to smith.

#### **Arguments**

l minMax

List of two numbers in waveform coordinates that describe the limits for the display. The first number is the minimum and the second is the maximum. If this argument is set to nil, the limit is set to auto.

#### Value Returned

t Returns t when the X axis display limits are set.

nil Returns nil and an error message if the X axis display limits

are not set.

#### **Example**

```
xLimit( list( 1 100 ) )
=> t
```

Sets the X axis to display between 1 and 100.

# Plotting and Printing Commands

# yLimit

# **Description**

Sets the Y axis display limits for the waveforms associated with a particular Y axis and strip in the current subwindow.

If you do not specify  $x\_stripNumber$ , the limits are applied when the subwindow is in composite mode.

# **Arguments**

l_minMax	List of two numbers in waveform coordinates that describe the limits for the display. The first number is the minimum and the second is the maximum. If this argument is set to $\mathtt{nil}$ , the limit is set to $\mathtt{auto}$ .
?stripNumber x_stripNumber	Specifies the strip in which the y display is to be limited in the range specified by $l\_minMax$ .
	Valid values: 1 through 20

#### Value Returned

t	Returns t if the Y axis display limits are set.
nil	Returns $\mathtt{nil}$ and an error message if the Y axis display limits cannot be set.

#### **Example**

```
yLimit( list( 4.5 7.5 ) )
=> t
```

Sets Y axis 1 to display from 4.5 to 7.5.

# Plotting and Printing Commands

# Plotting and Printing SpectreRF Functions in OCEAN

You can access SpectreRF functions in OCEAN by using the getData function and then plot or print them in OCEAN using the ocnPrint and plot functions.

To take an example, after performing a spectre sp analysis in the Analog Design Environment, click Results - Direct Plot - Main Form. In the S-Parameter Results form, select the function and other options that you want to plot. Also, select the Add to Outputs option under the *Plot* button. Then, click *OK*. The expression will be added to the *Outputs* pane of the ADE window. When all the desired expressions are created in the *Outputs* pane, use the ADE – Session – Save Ocean Script command to create the OCEAN script for these plots.

To plot the expression in OCEAN, use the following command:

```
plot(<expression in Output pane>)
For example,
   plot(Gmax())
                     for Gmax in S-parameter analysis
```

You can print the functions using the ocnPrint command. For example:

```
ocnPrint( Gmax() Kf() )
```

After a spectre sp noise analysis, use the following command to access the sp noise data.

```
selectResult("sp noise")
```

A sample OCEAN script to help you print or plot NFmin (minimum noise figure), N F (noise figure), and RN (noise resistance) results follows. Plotting NNR (normalized noise resistance) is similar to plotting RN.

```
; start ocean with Spectre as the as the simulator.
simulator( 'spectre )
; specify design and model path
design( "/usr1/mnt4/myhome/simulation/myckt/schematic/netlist/myckt.c")
path( "/usr1/mnt4/myhome/models" )
; specify analysis used: sp with noise
analysis('sp ?start "100M" ?stop "10G" ?donoise "yes"
?oprobe "/PORT1" ?iprobe "/PORT0" )
;set design variables
         "r2" 37)
desVar(
        "r1" 150)
desVar(
;set temperature
temp(25)
;run sp noise analysis with the above desVar list.
run()
```

#### Plotting and Printing Commands

```
printf("\n simulation has finished.")
printf("\n selecting sp noise results")
selectResult("sp_noise")
printf("\n print NFmin and plot NF")
NFmin = getData("NFmin")
NF = getData("NF")
ocnPrint( NFmin )
plot( NF )
printf("\n plot Rn")
Rn = getData("RN" ?result "sp_noise")
plot( Rn ?expr '( "Rn" ) )
exit
```

For more information, see the section *Periodic Noise Analysis* and the appendix *Plotting Spectre S-Parameter Simulation Data* in the *Virtuoso Spectre Circuit Simulator RF Analysis User Guide*.

For more information on these functions, click these links: getData, sp, ocnPrint, and plot.

# Plotting and Printing Commands

# **OCEAN Aliases**

The aliases in this chapter provide you with shortcuts to commonly used pairs of commands. By default, these aliases operate on results previously selected with <u>selectResult</u>. However, you can also use an alias on a different set of results. For example, to specify a different set of results for the vm alias, use the following syntax.

```
vm( t net [?result s resultName] )
```

where  $s_{resultName}$  is the name of the datatype for the particular analysis you want.

You can use the vm alias on results stored in a different directory as follows:

```
vm( t net [?resultsDir t resultsDir] [?result s resultName] )
```

where  $t\_resultsDir$  is the name of a different directory containing PSF results, and  $s\_resultName$  is the name of a datatype contained in that directory. (If you specify another directory with  $t\_resultsDir$ , you must also specify the particular results with  $s\_resultName$ .)

#### List of Aliases

Alias	Syntax	Description
vm	<pre>vm( t_net [?resultsDir t_resultsDir] [?result s_resultname]) =&gt; o_waveform/ nil</pre>	Aliased to $mag(v())$ . Gets the magnitude of the voltage of a net.
vdb	<pre>vdb(t_net [?resultsDir t_resultsDir] [?result s_resultname]) =&gt; o_waveform/ nil</pre>	Aliased to $db20(v())$ . Gets the power gain in decibels from net in to net out.
vp	<pre>vp(t_net [?resultsDir t_resultsDir] [?result s_resultname]) =&gt; o_waveform/ nil</pre>	Aliased to $phase(v())$ . Gets the phase of the voltage of a net.

# **OCEAN Aliases**

# List of Aliases, continued

vr	<pre>vr(t_net [?resultsDir t_resultsDir] [?result s_resultname]) =&gt; o_waveform/ nil</pre>	Aliased to $real(v())$ . Gets the real part of a complex number representing the voltage of a net.
vim	<pre>vim(t_net [?resultsDir t_resultsDir] [?result s_resultname]) =&gt; o_waveform/ nil</pre>	Aliased to $imag(v())$ . Gets the imaginary part of a complex number representing the voltage of a net.
im	<pre>im(t_component [?resultsDir t_resultsDir] [?result s_resultName]) =&gt; 0_waveform/ nil</pre>	Aliased to $mag(i())$ . Gets the magnitude of the AC current through a component.
ip	<pre>ip(t_component [?resultsDir t_resultsDir] [?result s_resultName]) =&gt; 0_waveform/ nil</pre>	Aliased to $phase(i())$ . Gets the phase of the AC current through a component.
ir	<pre>ir(t_component [?resultsDir t_resultsDir] [?result s_resultName]) =&gt; 0_waveform/ nil</pre>	Aliased to real (i()). Gets the real part of a complex number representing the AC current through a component.
iim	<pre>iim(t_component [?resultsDir t_resultsDir] [?result s_resultName]) =&gt; 0_waveform/ nil</pre>	Aliased to imag(i()). Gets the imaginary part of a complex number representing the AC current through a component.

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# Predefined and Waveform (Calculator) Functions

This chapter contains information about the following functions. Some additional predefined data access commands are described in the <u>Virtuoso Analog Design Environment L SKILL Language Reference</u>.

#### Predefined Arithmetic Functions

abs

<u>acos</u>

add1

<u>asin</u>

<u>atan</u>

<u>cos</u>

<u>exp</u>

int

<u>linRg</u>

<u>log</u>

logRg

max

<u>min</u>

mod

<u>random</u>

round

# Predefined and Waveform (Calculator) Functions

<u>sin</u> <u>sqrt</u> <u>srandom</u> sub1 <u>tan</u> <u>xor</u> Waveform (Calculator) Functions <u>average</u> abs\_jitter analog2Digital <u>awvCreateBus</u> <u>awvPlaceXMarker</u> <u>awvPlaceYMarker</u> <u>awvRefreshOutputPlotWindows</u> <u>b1f</u> bandwidth clip <u>clipX</u> closeResults compare compression compressionVRI compressionVRICurves complex complexp conjugate convolve

# Predefined and Waveform (Calculator) Functions

<u>cPwrContour</u> cReflContour <u>cross</u> <u>db10</u> db20 <u>dbm</u> <u>delay</u> <u>delayMeasure</u> <u>deriv</u> dft dftbb <u>dnl</u> <u>dutyCycle</u> <u>evmQAM</u> <u>evmQpsk</u> <u>eyeDiagram</u> <u>eyeAperture</u> <u>eyeMeasurement</u> edgeTriggeredEyeDiagram <u>flip</u> <u>fourEval</u> freq freq\_jitter **frequency** <u>ga</u> gac gainBwProd

# Predefined and Waveform (Calculator) Functions

<u>gainMargin</u>
gmax
<u>gmin</u>
gmsg
gmux
<u>gp</u>
<u>gpc</u>
groupDelay
<u>gt</u>
harmonic
<u>harmonicFreqList</u>
harmonicList
<u>histo</u>
histogram2D
<u>iinteg</u>
imag
<u>inl</u>
integ
intersect
i <u>pn</u>
<u>ipnVRI</u>
<u>ipnVRICurves</u>
<u>kf</u>
<u>ln</u>
<u>log10</u>
<u>lsb</u>
<u>lshift</u>

# Predefined and Waveform (Calculator) Functions

<u>mag</u>
<u>nc</u>
normalQQ
<u>overshoot</u>
pavg
<u>peak</u>
<u>peakToPeak</u>
period_jitter
<u>phase</u>
phaseDeg
<u>phaseDegUnwrapped</u>
<u>phaseMargin</u>
phaseRad
phaseRadUnwrapped
<u>PN</u>
<u>pow</u>
<u>prms</u>
<u>psd</u>
<u>psdbb</u>
<u>pstddev</u>
<u>pzbode</u>
pzfilter
rapidIPNCurves
<u>rapidIIPN</u>
<u>real</u>
<u>riseTime</u>
<u>rms</u>

# Predefined and Waveform (Calculator) Functions

<u>rmsNoise</u>
<u>rmsVoltage</u>
<u>rmsTerminalVoltage</u>
<u>root</u>
<u>rshift</u>
<u>sample</u>
<u>settlingTime</u>
<u>slewRate</u>
<u>smithType</u>
<u>spectralPower</u>
<u>spectrumMeas</u>
<u>spectrumMeasurement</u>
<u>ssb</u>
<u>stddev</u>
tangent
<u>thd</u>
<u>thd_fd</u>
<u>unityGainFreq</u>
<u>value</u>
<u>xmax</u>
<u>xmin</u>
<u>xval</u>
<u>ymax</u>
<u>ymin</u>
RF Functions

# Predefined and Waveform (Calculator) Functions

# **Predefined Arithmetic Functions**

Several functions are predefined in the Virtuoso<sup>®</sup> SKILL language. The full syntax and brief definitions for these functions follows the table.

#### **Predefined Arithmetic Functions**

Synopsis	Result	
General Functions		
add1(n)	n + 1	
abs	n	
sub1(n)	n-1	
exp(n)	e raised to the power n	
$linRg(n_from, n_to, n_by)$	Returns list of numbers in linear range from $n\_from$ to $n\_to$ in $n\_by$ steps	
log(n)	Natural logarithm of n	
<pre>logRg(n_from, n_to, n_by)</pre>	Returns list of numbers in log10 range from $n\_from$ to $n\_to$ in $n\_by$ steps	
max(n1 n2)	Maximum of the given arguments	
min( <i>n1 n2</i> )	Minimum of the given arguments	
mod(x1 x2)	x1 modulo $x2$ , that is, the integer remainder of dividing $x1$ by $x2$	
round(n)	Integer whose value is closest to n	
sqrt(n)	Square root of n	
Trigonometric Functions		
sin(n)	sine, argument $n$ is in radians	
$\cos(n)$	cosine	

tan(n)

asin(n)

arc cosine arc tangent

tangent

arc sine, result is in radians

# Predefined and Waveform (Calculator) Functions

# **Predefined Arithmetic Functions**

Synopsis	Result
Random Number Generator	
random(x)	Returns a random integer between 0 and $x$ -1. If random is called with no arguments, it returns an integer that has all of its bits randomly set.
srandom(x)	Sets the initial state of the random number generator to $\boldsymbol{x}$ .

# Predefined and Waveform (Calculator) Functions

### abs

# **Description**

Returns the absolute value of a floating-point number or integer.

# **Arguments**

n number

Floating-point number or integer.

### **Value Returned**

 $n_result$ 

The absolute value of  $n_number$ .

```
abs( -209.625)
=> 209.625
abs( -23)
=> 23
```

# Predefined and Waveform (Calculator) Functions

#### acos

# **Description**

Returns the arc cosine of a floating-point number or integer.

# **Arguments**

n number

Floating-point number or integer.

### Value Returned

f result

Returns the arc cosine of  $n_number$ .

### **Example**

acos(0.3) => 1.266104

# Predefined and Waveform (Calculator) Functions

# add1

# **Description**

Adds 1 to a floating-point number or integer.

# **Arguments**

n number

Floating-point number or integer to increase by 1.

### **Value Returned**

 $n_result$ 

 $n_number plus 1.$ 

# **Example**

Adds 1 to 59.

# Predefined and Waveform (Calculator) Functions

# asin

```
asin(
    n_number
)
=> f_result
```

# **Description**

Returns the arc sine of a floating-point number or integer.

# **Arguments**

n number

Floating-point number or integer.

### **Value Returned**

f\_result

The arc sine of  $n_number$ .

```
asin(0.3) => 0.3046927
```

# Predefined and Waveform (Calculator) Functions

### atan

# **Description**

Returns the arc tangent of a floating-point number or integer.

# **Arguments**

n number

Floating-point number or integer.

### **Value Returned**

f\_result

The arc tangent of  $n_number$ .

# **Example**

atan(0.3) => 0.2914568

# Predefined and Waveform (Calculator) Functions

#### cos

```
cos(
    n_number
)
=> f_result
```

# **Description**

Returns the cosine of a floating-point number or integer.

# **Arguments**

n number

Floating-point number or integer.

### **Value Returned**

 $f_result$ 

The cosine of  $n_number$ .

```
cos(0.3)
=> 0.9553365
cos(3.14/2)
=> 0.0007963
```

# Predefined and Waveform (Calculator) Functions

# exp

# **Description**

Raises e to a given power.

# **Arguments**

n number

Power to raise e to.

### **Value Returned**

f\_result

The value of e raised to the power  $n_number$ .

```
exp(1)
=> 2.718282
exp(3.0)
=> 20.08554
```

### Predefined and Waveform (Calculator) Functions

### int

### **Description**

Returns the largest integer not larger than the given argument.

**Note:** This function works on vector as well as waveform data. The function is applied to individual elements of the vector and waveform data.

### **Arguments**

n\_arg

A numeric value (which can be integer or floating point number).

#### Value Returned

x result

The value of the largest integer not larger than the value specified by  $n_{arg}$ .

```
int( 3.01 )
=> 3
int( 3.99 )
=> 3
```

# Predefined and Waveform (Calculator) Functions

# linRg

```
linRg(
    n_from n_to n_by
)
    => l_range / nil
```

# **Description**

Returns a list of numbers in the linear range from  $n_from$  to  $n_to$  incremented by  $n_by$ .

# **Arguments**

n_from	Smaller number in the linear range.
n_to	Larger number in the linear range.
n_by	Increment value when stepping through the range.

#### Value Returned

l_range	List of numbers in the linear range.
nil	Returned if error.

```
range = linRg(-30 30 5)
(-30 -25 -20 -15 -10 -5 0 5 10 15 20 25 30)
```

# Predefined and Waveform (Calculator) Functions

# log

```
log(
     n_number
)
     => f_result
```

# **Description**

Returns the natural logarithm of a floating-point number or integer.

# **Arguments**

n number

Floating-point number or integer.

### **Value Returned**

f\_result

The natural logarithm of  $n_number$ .

### Predefined and Waveform (Calculator) Functions

# logRg

```
logRg(
    n_from n_to n_by
)
=> l_range / nil
```

### **Description**

Returns a list of numbers in the log10 range from  $n_from$  to  $n_to$  advanced by  $n_by$ .

The list is a geometric progression where the multiplier is 10 raised to the  $1/n_by$  power. For example if  $n_by$  is 0.5, the multiplier is 10 raised to the 2nd power or 100.

### **Arguments**

n_from	Smaller number in the linear range.
n_to	Larger number in the linear range.
n_by	Increment value when stepping through the range.

#### Value Returned

l_range	List of numbers in the linear range.
nil	Returned if error.

```
logRg(1 1M 0.5)
(1.0 100.0 10000.0 1000000.0)
```

# Predefined and Waveform (Calculator) Functions

#### max

# **Description**

Returns the maximum of the values passed in. Requires a minimum of two arguments.

# **Arguments**

n_num1	First value to check.
n_num2	Next value to check.
[n_num3]	Additional values to check.

#### Value Returned

n\_result The maximum of the values passed in.

```
\max (3 \ 2 \ 1)
=> 3
\max (-3 \ -2 \ -1)
=> -1
```

# Predefined and Waveform (Calculator) Functions

# min

# **Description**

Returns the minimum of the values passed in. Requires a minimum of two arguments.

# **Arguments**

n_num1	First value to check.
n_num2	Next value to check.
[n_num3]	Additional values to check.

#### Value Returned

n result The minimum of the values passed in.

### **Example**

```
\min (1 \ 2 \ 3)
=> 1
\min (-1 \ -2.0 \ -3)
=> -3.0
```

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# Predefined and Waveform (Calculator) Functions

### mod

```
mod(
    x_integer1
    x_integer2
)
    => x_result
```

# **Description**

Returns the integer remainder of dividing two integers. The remainder is either zero or has the sign of the dividend.

# **Arguments**

x_integer1	Dividend.
x_integer2	Divisor.

#### Value Returned

x_result	The integer remainder of the division. The sign is determined by
	the dividend.

```
mod(4 3) => 1
```

### Predefined and Waveform (Calculator) Functions

### random

### **Description**

Returns a random integer between 0 and  $x_number$  minus 1.

If you call random with no arguments, it returns an integer that has all of its bits randomly set.

### **Arguments**

x\_number An integer.

### **Value Returned**

 $x\_result$  Returns a random integer between 0 and  $x\_number$  minus 1.

```
random( 93 ) => 26
```

# Predefined and Waveform (Calculator) Functions

### round

# **Description**

Rounds a floating-point number to its closest integer value.

# **Arguments**

n\_arg

Floating-point number.

### **Value Returned**

 $x_result$ 

The integer whose value is closest to n arg.

```
round(1.5)
=> 2
round(-1.49)
=> -1
round(1.49)
=> 1
```

### Predefined and Waveform (Calculator) Functions

### sin

```
sin(
     n_number
)
=> f result
```

### **Description**

Returns the sine of a floating-point number or integer.

# **Arguments**

n number

Floating-point number or integer.

#### Value Returned

f result

The sine of  $n_number$ .

# **Example**

```
sin(3.14/2)
=> 0.9999997
sin(3.14159/2)
=> 1.0
```

Floating-point results from evaluating the same expressions might be machine-dependent.

# Predefined and Waveform (Calculator) Functions

# sqrt

```
sqrt(
    n_number
)
=> f result
```

# **Description**

Returns the square root of a floating-point number or integer.

# **Arguments**

n number

Floating-point number or integer.

### **Value Returned**

 $f_result$ 

The square root of  $n_number$ .

```
sqrt(49)
=> 7.0
sqrt(43942)
=> 209.6235
```

# Predefined and Waveform (Calculator) Functions

### srandom

# **Description**

Sets the seed of the random number generator to a given number.

# **Arguments**

x number

An integer.

# **Value Returned**

t

This function always returns t.

```
srandom( 89 )
=> t
```

# Predefined and Waveform (Calculator) Functions

# sub1

```
sub1(
    n_number
)
=> n result
```

# **Description**

Subtracts 1 from a floating-point number or integer.

# **Arguments**

n number

Floating-point number or integer.

### **Value Returned**

 $n_result$ 

Returns n\_number minus 1.

# **Example**

Subtracts 1 from 59.

# Predefined and Waveform (Calculator) Functions

### tan

```
tan(
     n_number
)
=> f_result
```

# **Description**

Returns the tangent of a floating-point number or integer.

# **Arguments**

n number

Floating-point number or integer.

### **Value Returned**

f\_result

The tangent of  $n_number$ .

```
tan(3.0) => -0.1425465
```

### Predefined and Waveform (Calculator) Functions

#### xor

# **Description**

Returns the XOR value of the boolean inputs.

### **Arguments**

g in1 The first boolean input.

 $g_in2$  The second boolean input.

#### Value Returned

g\_res The resultant XOR value.

# Predefined and Waveform (Calculator) Functions

# **Waveform (Calculator) Functions**

The calculator commands are described in this section.

### Predefined and Waveform (Calculator) Functions

# average

```
average(
    o_waveform
    [ ?overall t_overall ]
    )
    => n average / o waveformAverage / nil
```

### **Description**

Computes the average of a waveform over its entire range.

Average is defined as the integral of the expression f(x) over the range of x, divided by the range of x.

For example, if y=f(x), average (y) =

$$\int_{0}^{\infty} f(x)dx$$

$$\frac{from}{to - from}$$

where from is the initial value for x and to is the final value.

### Predefined and Waveform (Calculator) Functions

### **Arguments**

be displayed as a series of points on a grid. (A waveform

object identifier looks like this: srrWave: XXXXX.)

?overall t overall

#### Value Returned

n average Returns a number representing the average value of the

input waveform.

o\_waveformAverage Returns a waveform (or family of waveforms) representing

the average value if the input is a family of waveforms.

nil Returns nil and an error message otherwise.

### **Example**

average( v( "/net9" ) )

Gets the average voltage (Y-axis value) of /net9 over the entire time range specified in the simulation analysis.

### Predefined and Waveform (Calculator) Functions

# abs\_jitter

```
abs_jitter(
    o_waveform
    t_crossType
    n_threshold
    [?xUnit t_xUnit]
    [?yUnit t_yUnit]
    [?Tnom n_Tnom])
=> o_waveform / nil
```

# **Description**

Calculates the absolute jitter values in the intput waveform for the given threshold. The output waveform can be expressed in degrees, radians, or unit intervals (UI). The absolute jitter can be plotted as a function of cycle number, crossing time, or reference clock time.

### Predefined and Waveform (Calculator) Functions

### **Arguments**

o waveform bject representing simulation results that

can be displayed as a series of points on a grid. (A

waveform object identifier looks like this:

srrWave:XXXXX.)

t crossType The points at which the curves of the waveform

intersect with the threshold. While intersecting, the

curve may be either rising or falling.

Valid values: rising and falling, respectively.

Default crossType is rising.

n threshold The threshold value against which the at which the

input waveform intersects to calculate the absolute

jitter.

?xUnit t xUnit The unit defined for X-axis of the output waveform.

Valid values: s (time) and cycle.

Default: s

Cycle numbers refer to the n'th occurrence where the

waveform crosses the given threshold.

?yUnit t yUnit The unit defined for Y-axis of the output waveform.

Valid values: rad (radians), UI (unit intervals), and S

(degrees)

Default value: rad.

? Tnom *n* Tnom The nominal time period of the input waveform. The

waveform is expected to be a periodic waveform that contains noise. If *Tnom* is nil, the abs\_jitter function finds the approximate average time period of

the input waveform.

Default value: nil.

#### Value Returned

o waveform Returns a waveform representing the absolute jitter

value for the given threshold.

nil Returns nil and an error message otherwise.

# Predefined and Waveform (Calculator) Functions

# Example

```
abs_jitter(v("net9" "rising" 1.0 ?xUnit "cycle" ?yUnit "UI" )
```

Gets the absolute jitter /net9 for the threshold value 1.0. Thom value is selected as nil.

### Predefined and Waveform (Calculator) Functions

# analog2Digital

```
analog2Digital(
    o_wave
    t_thresholdType
    [?vhi n_vhi]
    [?vlo n_vlo]
    [?vc n_vc]
    [?timeX n_timex]
)
    => o_digWave / n_digval / nil
```

# **Description**

Returns the digital form of the analog input, which can be a waveform, list or family of waveforms, or a string representation of expression(s).

# Predefined and Waveform (Calculator) Functions

# **Arguments**

o_wave	Input waveform.
t_thresholdType	Can take the values hilo or centre. If $t\_thresholdType$ is centre, it is a high state (1) unless its value is less than $n\_vc$ , in which case it is a low state (0). If $t\_thresholdType$ is hilo, any value less than $n\_vlo$ is a low state (0), any value greater than $n\_vhi$ is a high state (1) and the rest is treated as unknown based on the value of $n\_timex$ .
?vhi <i>o_vhi</i>	High threshold value (used only when $t\_thresholdType$ is hilo). If you do not specify this value, it is calculated internally as:
	<pre>vHigh = (topLine - bottomLine)*0.6 + bottomline</pre>
?vlo o_vlo	Low threshold value (used only when $t\_thresholdType$ is hilo). If you do not specify this value, it is calculated internally as:
	<pre>vLow = (topLine - bottomLine)*0.4 + bottomline</pre>
?vc <i>o_vc</i>	Central threshold value (used only when $t\_thresholdType$ is centre). If you do not specify this value, it is calculated internally using vCenter = average (wave).
?timeX n_timeX	The value that determines logic X. The timeX value is used to decide the state X. When threshold is $hilo$ , the analog signal will be converted to logic X if:
	■ analog signal value lies between vHigh and vLow
	■ lapse time between vHigh and vLow is larger than timeX

# **Value Returned**

o_digWave	A waveform (or a list of waveforms) is returned if the analog input specified was o_wave.
o_digVal	A scalar value is returned if the analog input specified was $o_{val}$ .
nil	Returns nil if the specified Waveform window does not exist.

# Predefined and Waveform (Calculator) Functions

# Example

# analog2Digital(v

### Predefined and Waveform (Calculator) Functions

### **awvCreateBus**

```
awvCreateBus(
    w_bus
    l_wavelist
    r_radix
)
```

#### **Definition**

Creates a bus with the given digital signals and radix.

### **Arguments**

w_bus	Name of the digital waveform representing a bus.
l_wavelist	List of the digital waveforms in the bus.
r_radix	Radix of the bus.

#### Value Returned

None.

### **Example**

Following are the examples to create a digital binary bus with name bus.

```
awvCreateBus("bus" list( awvAnalog2Digital( v("/data<0> " ?result
"tran-tran") nil nil 0.5 nil "centre")

awvAnalog2Digital( v("/datab<1> " ?result "tran-tran") nil nil 0.5
nil "centre")

awvAnalog2Digital( v("/data<1> " ?result "tran-tran") nil nil 0.5 nil
"centre")

awvAnalog2Digital( v("/datab<0> " ?result "tran-tran") nil nil 0.5
nil "centre") ) "Binary")
```

### Predefined and Waveform (Calculator) Functions

### awvPlaceXMarker

```
awvPlaceXMarker(
    w_windowId
    n_xLoc
    [ ?subwindow x_subwindowId ]
    [ ?label t_label ]
)
    => t xLoc / t / nil
```

# **Description**

Places a vertical marker at a specific x-coordinate in the optionally specified subwindow of the specified window.

# **Arguments**

w_windowId	Waveform window ID.
$n\_xLoc$	The x-coordinate at which to place the marker.
?subwindow	$x\_subwindowId$
	Waveform subwindow ID.

```
?label t label
```

#### **Value Returned**

t_xLoc	Returns a string of x-coordinates if the command is successful and the vertical marker info form is opened.
t	Returns this when the command is successful but the vertical marker info form is not opened.
nil	Returns nil or an error message.

### **Example**

```
awvPlaceXMarker( window 5) => "5"
```

Vertical marker info form is opened when the command is executed.

```
awvPlaceXMarker( window 6 ?subwindow 2) => t
```

# Predefined and Waveform (Calculator) Functions

Vertical marker info form is not opened.

### Predefined and Waveform (Calculator) Functions

### awvPlaceYMarker

```
awvPlaceYMarker(
    w_windowId
    n_yLoc
    [ ?subwindow x_subwindowId ]
    [ ?label t_label ]
    )
    => t_yLoc / t / nil
```

### **Description**

Places a horizontal marker at a specific y-coordinate in the optionally specified subwindow of the specified window.

# **Arguments**

w_windowId	Waveform window ID.
n_yLoc	The y-coordinate at which to place the marker.
?subwindow x_subwindow	Id
	Waveform subwindow ID.
?label t_label	

#### Value Returned

t_yLoc	Returns a string of y-coordinates if the command is successful and the horizontal marker info form is opened.
t	Returns this when the command is successful but the horizontal marker info form is not opened.
nil	Returns nil or an error message.

# **Example**

■ Paces a horizontal marker with a label, myHorizontalMarker1, at Y= 2.0 in the strip number 1 of the current subwindow of the current graph window:

```
awvPlaceYMarker(gw 2.0 ?label "myHorizontalMarker1" ?subwindow gsw ?stripNum
1)
```

### Predefined and Waveform (Calculator) Functions

Where, gw= awvGetCurrentWindow() and gsw= awvGetCurrentSubwindow(gw)

■ Paces a horizontal marker with a label, myHorizontalMarker2, at Y= 2.0 in the strip number 2 of the current subwindow of the current graph window:

awvPlaceYMarker(gw 2.0 ?label "myHorizontalMarker2" ?subwindow gsw ?stripNum
2)

Where, gw= awvGetCurrentWindow() and gsw= awvGetCurrentSubwindow(gw)

# Predefined and Waveform (Calculator) Functions

# awv Refresh Output Plot Windows

```
awvRefreshOutputPlotWindows(
    s_session
)
```

# **Description**

Refreshes all existing plot windows (with new simulation data, if any) attached with the session *s\_session*.

# **Arguments**

s session

Currently active environment variable.

### **Value Returned**

None.

# Predefined and Waveform (Calculator) Functions

# b1f

# **Description**

Returns the alternative stability factor in terms of the supplied parameters.

# **Arguments**

o_s11	Waveform object representing s11.
o_s12	Waveform object representing s12.
o_s21	Waveform object representing s21.
o_s22	Waveform object representing s22.

### Value Returned

o_waveform	Waveform object representing the alternative stability factor.
nil	Returns nil and an error message otherwise.

```
s11 = sp(1 1)
s12 = sp(1 2)
s21 = sp(2 1)
s22 = sp(2 2)
plot(blf(s11 s12 s21 s22))
```

## Predefined and Waveform (Calculator) Functions

## bandwidth

```
bandwidth(
    o_waveform
    n_db
    t_type
)
    => n_value / o_waveform / nil
```

# **Description**

Calculates the bandwidth of a waveform.

# **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
$n\_db$	Positive number that defines the bandwidth.
t_type	Type of input filter.
	Valid values: "low", "high" or "band".

### Value Returned

n_value	Returns a number representing the value of the bandwidth if the input argument is a single waveform.
o_waveform	Returns a waveform (or family of waveforms) representing the bandwidth if the input argument is a family of waveforms.
nil	Returns nil and an error message otherwise.

# **Example**

```
bandwidth( v( "/OUT" ) 3 "low")
```

Gets the 3 dB bandwidth of a low-pass filter.

```
bandwidth( v( "/OUT" ) 4 "band" )
```

Gets the 4 dB bandwidth of a band-pass filter.

### Predefined and Waveform (Calculator) Functions

# clip

```
clip(
    o_waveform
    n_from
    n_to
)
    => o waveform / nil
```

# **Description**

Restricts the waveform to the range defined by  $n_from$  and  $n_to$ .

You can use the clip function to restrict the range of action of other commands. If  $n\_from$  is  $nil, n\_from$  is taken to be the first X value of the waveform, and if  $n\_to$  is  $nil, n\_to$  is taken to be the last X value of the waveform. If both  $n\_to$  and  $n\_from$  are nil, the original waveform is returned.

# **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_from	Starting value for the range on the X axis.
n_to	Ending value for the range on the X axis.

#### Value Returned

o_waveform	Returns a waveform object if the input argument is a waveform object or returns a family of waveforms if the input argument is a family of waveforms.
nil	Returns nil and an error message otherwise.

## **Example**

```
x = clip(v("/net9") 2m 4m)
plot(x)
```

Plots the portion of a waveform that ranges from 2 ms to 4 ms.

```
plot( clip( v( "/net9" ) nil nil ) )
```

# Predefined and Waveform (Calculator) Functions

Plots the original waveform.

```
plot( clip( v( "/net9" ) nil 3m ) )
```

Plots the portion of a waveform that ranges from 0 to 3 ms.

# Predefined and Waveform (Calculator) Functions

# clipX

```
clipX(
    o_waveform
    n_from
    n_to
)
    => o waveform / nil
```

## **Description**

Restricts the waveform to the range defined by  $n_from$  and  $n_to$ .

The clipX function works in the same manner as the clip function, with an exception that clipX restricts the waveform within the range [from, to] without interpolating or extrapolating any values. In other words, clipX returns a waveform that consists only data points from the input waveform.

**Note:** The clipX function snaps the from and/or to values to a data point in the input waveform if these values are less than 1e-6 \* stepSize, where stepSize is the average step of the input waveform in range [from,to].

# **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_from	Starting value for the range on the X axis.
n_to	Ending value for the range on the X axis.

### Value Returned

o_waveform	Returns a waveform object if the input argument is a waveform object or returns a family of waveforms if the input argument is a family of waveforms.
nil	Returns nil and an error message otherwise.

# Predefined and Waveform (Calculator) Functions

### closeResults

```
closeResults(
    t_dirName
)
    => t / nil
```

### **Definition**

Closes the simulation results stored in the input results directory. The function closes all the internal resources opened by the tool that are related to the results directory. It is recommended that you must call this function before deleting a results directory, moving the directory to any other location, or renaming a results directory.

After calling the closeResults function, the OCEAN commands, such as selectResults, getData, pv, which can also be called without passing the resultsDir argument and run based upon previously called openResults call, stops working if called without passing the resultsDir argument.

# **Arguments**

t_dirName	Name of the directory which was earlier used in the
	Description

openResults function.

### **Values Returned**

t	If the results database has been closed successfully.
nil	If the results database has not been closed successfully.

### Predefined and Waveform (Calculator) Functions

# compare

```
compare(
    o_waveform1
    o_waveform1
    [ f_abstol [ f_reltol ] ]
    )
    => o comparisonWaveform / nil
```

### **Description**

Compares the two given waveforms based on the specified values for absolute and relative tolerances. This function compares only the sections of the two waveforms where the X or independent axes overlap.

The following situations are possible:

- If neither relative nor absolute tolerance is specified, the function returns the difference of the two waveforms (o waveform1 o waveform2).
- If only the absolute tolerance is specified, the function returns the difference of the two waveforms only when the absolute value of the difference is greater than the absolute tolerance (|o\_waveform1 o\_waveform2| > f\_abstol); otherwise it returns a zero waveform.
- If only the relative tolerance is specified, the function returns the difference of the two waveforms only when the absolute value of the difference is greater than the product of the relative tolerance and the larger of the absolute values of the two waveforms (|o\_waveform1 o\_waveform2| > f\_reltol\* max(|o\_waveform1|, |o\_waveform2|)); otherwise it returns a zero waveform.
- If both relative and absolute tolerances are specified, the function returns the difference of the two waveforms only when the absolute value of the difference is greater than the sum of the separately calculated tolerance components (|o\_waveform1 o\_waveform2| > f\_abstol + f\_reltol \* max(|o\_waveform1|, |o waveform2|)); otherwise it returns a zero waveform.

**Note:** The function also compares parametric waveforms. However, for a successful comparison of parametric waveforms, the family tree structures of the two input waveforms should be the same. For both the input waveforms, the number of child waveforms at each level should also be the same, except at the leaf level where the elements are simple scalars. This is an obvious condition to obtain a meaningful comparison.

## Predefined and Waveform (Calculator) Functions

## **Arguments**

o\_waveform1 Waveform 1.

o\_waveform2 Waveform 2.

f\_abstol Absolute tolerance.
 Default value: 0.0

f\_reltol Relative tolerance.

#### Value Returned

o comparisonWaveform

Returns the difference of the two given waveforms based on the specified values of the relative and absolute tolerances.

nil Returns nil and an error message otherwise.

Default value: 0.0

## **Example**

```
compare( wave1 wave2 2.2 0.4 )
=> srrWave:175051528
```

Returns the difference of the waveforms wave1 and wave2 based on the specified absolute and relative tolerances of 2.2 and 0.4, respectively.

## Predefined and Waveform (Calculator) Functions

# compression

```
compression(
   o_waveform
   [?x f_x]
   [?y f_y]
   [?compression f_compression]
   [?io s_measure]
   [?tanSlope t_tanSlope]
)
   => f compPoint / nil
```

## **Description**

Performs an *n*th compression point measurement on a power waveform.

The compression function uses the power waveform to extrapolate a line of constant slope (dB/dB) according to a specified input or output power level. This line represents constant small-signal power gain (ideal gain). The function finds the point where the power waveform drops n dB from the constant slope line and returns either the X coordinate (input referred) value or the Y coordinate (output referred) value.

# Predefined and Waveform (Calculator) Functions

# **Arguments**

o_waveform	Waveform object representing output power (in dBm) versus input power (in dBm).
?x f_x	The X coordinate value (in dBm) used to indicate the point on the output power waveform where the constant-slope power line begins. This point should be in the linear region of operation.
	Default value: Unless $f_y$ is specified, defaults to the X coordinate of the first point of the $o_{waveform}$ wave.
?y f_y	The Y coordinate value (in dBm) used to indicate the point on the output power waveform where the constant-slope power line begins. This point should be in the linear region of operation.
	Default value: Unless $f_x$ is specified, defaults to the Y coordinate of the first point of the $o_{waveform}$ wave.
?compression f_compression	The delta (in dB) between the power waveform and the ideal gain line that marks the compression point.
	Default value: 1
?io s_measure	Symbol indicating whether the measurement is to be input referred ('input) or output referred ('output).
	Default value: input
?tanSlope tranSlope	Default value: 1

### **Value Returned**

f_compPoint	Depending on the setting of $s_{measure}$ , returns either input referred or output referred compression point.
nil	Returns nil and an error message otherwise.

```
xloc = compression( wave ?x -25 ?compress 1)
yloc = compression( wave ?x -25 ?measure "Output")
; Each of following returns a compression measurement:
compression(dB20(harmonic(v("/Pif" ?result "pss_fd") 2)))
```

## Predefined and Waveform (Calculator) Functions

### Predefined and Waveform (Calculator) Functions

# compressionVRI

```
compressionVRI(
    o_vport
    x_harm
    [?iport o_iport]
    [?rport f_rport]
    [?epoint f_epoint]
    [?gcomp f_gcomp]
    [?measure s_measure]
    [?format format]
)
    => o waveform / n number / nil
```

# Description

Performs an *n*th compression point measurement on a power waveform.

Use this function to simplify the declaration of a compression measurement. This function extracts the specified harmonic from the input waveform(s), and uses  ${\tt dBm(spectralPower((i\ or\ v/r),v))} \ \ to\ calculate\ a\ power\ waveform. \ The\ function\ passes\ this\ power\ curve\ and\ the\ remaining\ arguments\ to\ the\ {\tt compression}\ function\ to\ complete\ the\ measurement.$ 

The compression function uses the power waveform to extrapolate a line of constant slope (dB/dB) according to a specified input or output power level. This line represents constant small-signal power gain (ideal gain). The function finds the point where the power waveform drops n dB from the constant slope line and returns either the X coordinate (input referred) value or the Y coordinate (output referred) value.

### Predefined and Waveform (Calculator) Functions

### **Arguments**

o vport Voltage across the output port. This argument must be a

family of spectrum waveforms (1 point per harmonic) created by parametrically sweeping an input power (in

dBm) of the circuit.

x harm Harmonic index of the voltage wave contained in

o vport. When o iport is specified, also applies to a

current waveform contained in o iport.

?iport o iport Current into the output port. This argument must be a

family of spectrum waveforms (1 point per harmonic) created by parametrically sweeping an input power (in dBm) of the circuit. When specified, the output power is

calculated using voltage and current.

Default value: nil

?rport f rport Resistance into the output port. When specified and

o iport is nil, the output power is calculated using

voltage and resistance.

Default value: 50

?epoint f epoint The X coordinate value (in dBm) used to indicate the point

on the output power waveform where the constant-slope power line begins. This point should be in the linear region

of operation.

Default value: the X coordinate of the first point of the

o\_waveform wave

?gcomp f gcomp The delta (in dB) between the power waveform and the

ideal gain line that marks the compression point.

Default value: 1

?measure s measure Symbol indicating if measurement is to be input referred

('input) or output referred ('output).

Default value: input

?format format Default Value: power

#### Value Returned

o\_waveform Returns a waveform when o\_waveform1 is a family of

waveforms.

# Predefined and Waveform (Calculator) Functions

f\_number Returns a number when o\_waveform1 is a waveform.

Returns nil and an error message otherwise.

### Predefined and Waveform (Calculator) Functions

# compressionVRICurves

```
compressionVRICurves(
   o_vport
   x_harm
   [ ?iport o_iport ]
   [ ?rport f_rport ]
   [ ?epoint f_epoint ]
   [ ?gcomp f_gcomp ]
   [ ?format format ]
   )
   => o waveform / nil
```

# Description

Constructs the waveforms associated with an *n*th compression measurement.

Use this function to simplify the creation of waveforms associated with a compression measurement. This function extracts the specified harmonic from the input waveform(s), and uses dBm(spectralPower((i or v/r), v)) to calculate a power waveform.

The compressionVRICurves function uses the power waveform to extrapolate a line of constant slope (1dB/1dB) according to a specified input or output power level. This line represents constant small-signal power gain (ideal gain). The function shifts the line down by n dB and returns it, along with the power waveform, as a family of waveforms.

This function only creates waveforms and neither performs a compression measurement nor includes labels with the waveforms. Use the compression or compression VRI function for making measurements.

### Predefined and Waveform (Calculator) Functions

### **Arguments**

family of spectrum waveforms (1 point per harmonic) created by parametrically sweeping an input power (in

dBm) of the circuit.

x harm Harmonic index of the wave contained in o vport.

When o\_iport is specified, also applies to a current

waveform contained in o iport.

?iport o iport Current into the output port. This argument must be a

family of spectrum waveforms (1 point per harmonic) created by parametrically sweeping an input power (in dBm) of the circuit. When specified, the output power is

calculated using voltage and current.

Default value: nil

?rport f rport Resistance into the output port. When specified and

o iport is nil, the output power is calculated using

voltage and resistance.

Default value: 50

?epoint f epoint The X coordinate value (in dBm) used to indicate the

point on the output power waveform where the constantslope power line begins. This point should be in the linear

region of operation.

Default value: the X coordinate of the first point of the

o\_waveform wave

?gcomp f gcomp The delta (in dB) between the power waveform and the

ideal gain line that marks the compression point.

Default value: 1

?format format Default Value: power

#### Value Returned

o waveform Returns a family of waveforms containing the output

power and tangent line.

nil Returns nil and an error message otherwise.

# Predefined and Waveform (Calculator) Functions

## Predefined and Waveform (Calculator) Functions

# complex

```
complex(
   f_real
   f_imaginary
)
   => o complex
```

# **Description**

Creates a complex number of which the real part is equal to the real argument, and the imaginary part is equal to the imaginary argument.

## **Arguments**

f real The real part of the complex number.

f\_imaginary The imaginary part of the complex number.

### **Value Returned**

o complex Returns the complex number.

```
complex( 1.0 2.0 )
=> complex( 1, 2 )
```

## Predefined and Waveform (Calculator) Functions

# complexp

```
complexp(
    g_value
)
    => t / nil
```

# **Description**

Checks if an object is a complex number. The suffix  ${\tt p}$  is added to the name of a function to indicate that it is a predicate function.

## **Arguments**

g value

A skill object.

## **Value Returned**

Returns t when g value is a complex number.

nil

t

Returns nil if there is an error.

```
complexp( (complex 0 1) )
=> t
complexp( 1.0 )
=> nil
```

## Predefined and Waveform (Calculator) Functions

# conjugate

```
conjugate(
    { o_waveform | n_x }
)
=> o waveform / n y / nil
```

# **Description**

Returns the conjugate of a waveform or number.

# **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_x	Complex or imaginary number.

## **Value Returned**

o_waveform	Returns the conjugate of a waveform if the input argument is a waveform.
n_y	Returns the result of $n_x$ being mirrored against the real axis (X axis) if the input argument is a number.
nil	Returns nil and an error message otherwise.

# **Example**

For this example, assume that the first three statements are true for the conjugate function that follows them.

```
x=complex(-1 -2)

real(x) = -1.0

imag(x) = -2.0

conjugate(x) = complex(-1, 2)
```

Returns the conjugate of the input complex number.

## Predefined and Waveform (Calculator) Functions

## convolve

```
convolve(
    o_waveform1
    o_waveform2
    n_from
    n_to
    t_type
    n_by
)
    => o_waveform /n_number /nil
```

# **Description**

Computes the convolution of two waveforms.

Convolution is defined as

to
$$\int f1(s)f2(t-s)ds$$
from

£1 and £2 are the functions defined by the first and second waveforms.

**Note:** The convolve function is numerically intensive and might take longer than the other functions to compute.

## Predefined and Waveform (Calculator) Functions

# **Arguments**

Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
Additional waveform object.
Starting point (X-axis value) of the integration range.
Ending point (X-axis value) of the integration range.
Type of interpolation.
Valid values: "linear" or "log".
Increment.

### Value Returned

o_waveform	Returns a waveform object representing the convolution if one of the input arguments is a waveform. Returns a family of waveforms if either of the input arguments is a family of waveforms.
n_number	Returns a value representing the convolution if both of the input arguments are numbers.
nil	Returns nil and an error message otherwise.

# **Example**

Gets the waveform from the convolution of the sine waveform and triangle waveform within the range of 0 to 10.

# Predefined and Waveform (Calculator) Functions

### **cPwrContour**

```
cPwrContour(
    o_iwave
    o_vwave
    x_harm
    [ ?iwaveLoad o_iwaveLoad ]
    [ ?vwaveLoad o_vwaveLoad ]
    [ ?maxPower f_maxPower ]
    [ ?minPower f_minPower ]
    [ ?numCont x_numCont ]
    [ ?refImp f_refImp ]
    [ ?closeCont g_closeCont ]
    [ ?modifier s_modifier ]
    [ ?ifam ifam ]
    [ ?vfam vfam ]
    )
    => o_waveform / nil
```

## **Description**

Constructs constant power contours for Z-Smith plotting. The trace of each contour correlates to reference reflection coefficients that all result in the same power level.

The  $x_harm$  harmonic is extracted from all the input waveforms. Power is calculated using the spectral Power function. The reference reflection coefficients are calculated using voltage, current, and a reference resistance.

# Predefined and Waveform (Calculator) Functions

# **Arguments**

o_iwave	Current used to calculate power, expected to be a two- dimensional family of harmonic waveforms.
o_vwave	Voltage used to calculate power, expected to be a two- dimensional family of harmonic waveforms.
x_harm	Harmonic index of the waves contained in $o_iwave$ and $o_vwave$ .
?iwaveLoad o_iwaveLoad	Current used to calculate reflection coefficient, expected to be a two-dimensional family of harmonic waveforms.
	Default value: o_iwave
?vwaveLoad o_vwaveLoad	Voltage used to calculate reflection coefficient, expected to be a two-dimensional family of harmonic waveforms.
	Default value: o_vwave
?maxPower f_maxPower	Maximum power magnitude value for contours.
	Default value: automatic
?minPower f_minPower	Minimum power magnitude value for contours.
	Default value: automatic
?closeCont x_numCont	Total number of contours returned.
	Default value: 8
?refImp f_refImp	Reference resistance used to calculate reflection coefficients.
	Default value: 50
?closeCont g_closeCont	Boolean indicating when to close the contours. When nil, largest segment of each contour is left open.
	Default value: nil
?modifier s_modifier	Symbol indicating the modifier function to apply to the calculated power. The modifier function can be any single argument OCEAN function such as db10 or dBm.
	Default value: dbm

?ifam ifam

?vfam *vfam* 

# Predefined and Waveform (Calculator) Functions

#### Value Returned

o\_waveformnilReturns a family of waveforms (contours) for Z-Smith plotting.nilReturns nil and an error message otherwise.

### **Example**

The following example plots constant output power contours according to output:

The following example plots constant output power contours according to output reflection coefficients:

The following example plots constant input power contours according to output reflection coefficients:

# Predefined and Waveform (Calculator) Functions

### cReflContour

```
cReflContour(
    o_iwave
    o_vwave
    x_harm
    [ ?iwaveLoad o_iwaveLoad ]
    [ ?vwaveLoad o_vwaveLoad ]
    [ ?maxRefl f_maxRefl ]
    [ ?minRefl f_minRefl ]
    [ ?numCont x_numCont ]
    [ ?refImp f_refImp ]
    [ ?closeCont g_closeCont ]
    )
    => o waveform / nil
```

# Description

Constructs constant reflection coefficient magnitude contours for Z-Smith plotting. The trace of each contour correlates to reference reflection coefficients that all result in the same reflection coefficient magnitude.

The  $x_harm$  harmonic is extracted from all the input waveforms. Reflection coefficient magnitude is calculated using voltage, current, reference resistance, and the mag function. The reference reflection coefficients are calculated separately by using voltage, current, and a reference resistance.

# Predefined and Waveform (Calculator) Functions

# **Arguments**

o_iwave	Current used to calculate reflection coefficient magnitude, expected to be a two-dimensional family of spectrum waveforms.
o_vwave	Voltage used to calculate reflection coefficient magnitude, expected to be a two-dimensional family of spectrum waveforms.
x_harm	Harmonic index of the waves contained in $o_iwave$ and $o_vwave$ .
?iwaveLoad o_iwaveLoad	Current used to calculate reference reflection coefficient, expected to be a two-dimensional family of harmonic waveforms.
	Default value: o_iwave
?vwaveLoad o_vwaveLoad	Voltage used to calculate reference reflection coefficient, expected to be a two-dimensional family of spectrum waveforms.
	Default value: o_vwave
?maxRefl f_maxRefl	Maximum reflection coefficient magnitude value for contours.
	Default value: automatic
?minRefl f_minRefl	Minimum reflection coefficient magnitude value for contours.
	Default value: automatic
?numCont x_numCont	Total number of contours returned.
	Default value: 8
?refImp f_refImp	Reference resistance used to calculate reflection coefficients.
	Default value: 50
?closeCont g_closeCont	Boolean indicating when to close the contours. When $\min$ , the largest segment of each contour is left open.
	Default value: nil

## Predefined and Waveform (Calculator) Functions

#### Value Returned

o\_waveformnilReturns a family of waveforms (contours) for Z-Smith plotting.nil and an error message otherwise.

# **Example**

The following example plots constant output reflection coefficient contours according to output reflection coefficients:

The following example plots constant output reflection coefficient contours according to output reflection coefficients:

The following example plots constant output reflection coefficient contours according to output reflection coefficients:

```
cReflContour(i("/C25/PLUS" ?result "pss_fd")
   v("/net30" ?result "pss_fd") 1
   ?iwaveLoad i("/I8/out" ?result "pss_fd")
   ?vwaveLoad v("/net28" ?result "pss_fd") ?refImp 50.0
   ?numCont 9)
```

# Predefined and Waveform (Calculator) Functions

#### cross

```
cross(
    o_waveform
    n_crossVal
    x_n
    s_crossType
    [ g_multiple [ s_Xname ] ]
    )
    => o_waveform / g_value / nil
```

# **Description**

Computes the X-axis value at which a particular crossing of the specified edge type of the threshold value occurs.

# Predefined and Waveform (Calculator) Functions

# **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_crossVal	Y-axis value at which the corresponding values of X are calculated.
x_n	Number that specifies which X value to return. If $x_n$ equals 1, the first X value with a crossing is returned. If $x_n$ equals 2, the second X value with a crossing is returned, and so on. If you specify a negative integer for $x_n$ , the X values with crossings are counted from right to left (from maximum to minimum). If you specify $x_n$ equals to 0, it returns all occurrences of the crossing events.
s_crossType	Type of the crossing.
	Valid values: 'rising, 'falling, 'either.
g multiple	
<u>.</u>	An optional boolean argument that takes the value $nil$ by default. If set to $t$ , the value specified for the $x_n$ argument is ignored and the function returns all occurrences of the crossing event.
s_xName	default. If set to $t$ , the value specified for the $x_n$ argument is ignored and the function returns all occurrences of the crossing

# **Value Returned**

o_waveform	Returns a waveform if the input argument is a family of waveforms.
g_value	Returns the X-axis value of the crossing point if the input argument is a single waveform.
nil	Returns nil and an error message otherwise.

```
cross( v( "/net9" ) 2.5 2 'rising )
```

# Predefined and Waveform (Calculator) Functions

Gets the time value (X axis) corresponding to specified voltage "/net9"=2.5V (Y axis) for the second rising edge.

```
cross( v( "/net9" ) 1.2 1 'either )
```

Gets the time value (X axis) corresponding to specified voltage "/net9"=1.2V (Y axis) for the first edge, which can be a rising or falling edge.

```
cross(VT("/out") 2.5 0 0 t "time") (s)
```

Returns multiple occurrences of crossing events specified against time-points at which each crossing event occurs.

```
cross(VT("/out") 2.5 0 0 t "cycle") (s)
```

Returns multiple occurrences of crossing events specified against cycle numbers, where a cycle number refers to the n'th occurrence of the crossing event in the input waveform.

## Predefined and Waveform (Calculator) Functions

# db10

# **Description**

Returns 10 times the log10 of the specified waveform object or number. This function can also be written as dB10.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_number	Number.

### Value Returned

o_waveform	Returns a waveform object if the input argument is a waveform object or returns a family of waveforms if the input argument is a family of waveforms.
n_number	Returns a number if the input argument is a number.
nil	Returns nil and an error message otherwise.

## **Example**

```
db10( ymax( v( "/net9" ) ) )
```

Returns a waveform representing log10 (ymax (v("/net9")) multiplied by 10.

```
db10(1000)
=> 30.0
```

Gets the value log10 (1000) multiplied by 10, or 30.

## Predefined and Waveform (Calculator) Functions

## db20

```
db20(
     {o_waveform | n_number}
)
=> o waveform / n number / nil
```

# **Description**

Returns 20 times the log10 of the specified waveform object or number. This function can also be written as dB20.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_number	Number.

### Value Returned

o_waveform	Returns a waveform object if the input argument is a waveform object or returns a family of waveforms if the input argument is a family of waveforms.
n_number	Returns a number if the input argument is a number.
nil	Returns nil and an error message otherwise.

## **Example**

```
db20( ymax( v( "/net9" ) ) )
```

Returns a waveform representing 20 times log10 (ymax(v("/net9")). db20( 1000 ) => 60.0

Returns the value of 20 times log10 (1000), or 60.

## Predefined and Waveform (Calculator) Functions

## dbm

# **Description**

Returns 10 times the log10 of the specified waveform object plus 30. This function can also be written as dBm.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_number	Number.

### Value Returned

o_waveform	Returns a waveform object if the input argument is a waveform object or returns a family of waveforms if the input argument is a family of waveforms.
n_number	Returns a number if the input argument is a number.
nil	Returns nil and an error message otherwise.

## **Example**

```
dbm( ymax( v( "/net9" ) ) )
```

Returns a waveform representing 10 times log10 (ymax (v("/net9")) plus 30.

## Predefined and Waveform (Calculator) Functions

# delay

```
delay(
     [ ?wf1 o_waveform1 ]
     [ ?value1 n_value1 ]
     [ ?edge1 s_edge1 ]
     [ ?nth1 x nth1 ]
     [ ?td1 n_td1 ]
     [ ?wf2 o waveform2 ]
     [ ?value2 n_value2 ]
     [ ?edge2 s_edge2 ]
     [ ?nth2 x nth2
     [ ?td2 n td2 ]
     [ ?td2r0 n_td2r0 ]}
     [ ?stop n stop ]
     [ ?histoDisplay g_histoDisplay ]
     [ ?noOfHistoBins x noOfHistoBins ]
     [ ?period1 period1 ]
     [ ?period2 period2 ]
     [ ?multiple multiple ]
     [ ?xName xName ]
    @rest args
    => o waveform / n value / nil
```

# **Description**

Calculates the delay between a trigger event and a target event.

The delay command computes the delay between two points using the cross command.

# Predefined and Waveform (Calculator) Functions

# **Arguments**

?wf1 o_waveform1	First waveform object.
?value1 n_value1	Value at which the crossing is significant for the first waveform object.
?edge1 s_edge1	Type of the edge that must cross $n_value1$ . Valid values: rising, falling, either
	Default Value: either
?nth1	Number that specifies which crossing is to be the trigger event. For example, if $x_nth1$ is 2, the trigger event is the second edge of the first waveform with the specified type that crosses $n_value1$ .
	Default Value: 1
?td1	Time at which to start the delay measurement. The simulator begins looking for the trigger event, as defined by $o\_waveform1$ , $n\_value1$ , $t\_edge1$ , and $x\_nth1$ , only after the $n\_td1$ time is reached.
	Default Value: 0
?wf2 o_waveform2	Second waveform object.
?value2 <i>n_value2</i>	Value at which the crossing is significant for the second waveform.
?edge2 <i>s_edge2</i>	Type of the edge for the second waveform.
	Valid values: rising, falling, either
	Default Value: either
?nth2 <i>x_nth2</i>	Number that specifies which crossing is to be the target event. For example, if $x_nth2$ is 2, the target event is the second edge of the second waveform with the specified type that crosses $n_value2$ .
	Default Value: 1

#### Predefined and Waveform (Calculator) Functions

Time to start observing the target event.  $n_td2$  is specified relative to the trigger event. This parameter cannot be specified at the same time as  $n_td2r0$ .

The simulator begins looking for the target event, as defined by  $o\_waveform2$ ,  $n\_value2$ ,  $t\_edge2$ , and x nth2, only after the n td2 time is reached

If you specify neither  $n_t d2$  nor  $n_t d2r0$ , the simulator begins looking for the target event at t = 0.

**Note**: For non- multiple, If td2 is specified, find the cross point of wf1 at edge nth. Use this as trigger point for target(wf2) and ignore wf2 before wf1 trigger event to find its cross point. Calculate the delay between two cross points. If td2 is not specified, find the cross point at edge nth1 of wf1 and cross point of edge nth2 of wf2 with target at time at t=0 and calculate the delay between two cross points.

Time to start observing the target event, relative to t = 0. Only applicable if both  $o\_waveform1$  and  $o\_waveform2$  are specified. This parameter cannot be specified at the same time with  $n\_td2....$ The simulator begins looking for the target event, as defined by  $o\_waveform2$ ,  $n\_value2$ ,  $t\_edge2$ , and x nth2, only after the n tdr0 time is reached.

f you specify neither  $n_td2$  nor  $n_td2r0$ , the simulator begins looking for the target event at t = 0.

?td2 and ?td2r0 take precedence over other options.

?stop n\_stop

Time to stop observing the target event.

?histoDisplay g histoDisplay

When set to t, returns a waveform that represents the statistical distribution of the riseTime data in the form of a histogram. The height of the bars (bins) in the histogram represents the frequency of the occurrence of values within the range of riseTime data.

Valid values: t nil
Default value: nil

#### Predefined and Waveform (Calculator) Functions

?noOfHistoBins x noOfHistoBins

Denotes the number of bins represented in the

histogram representation.

Valid values: Any positive integer

Default value: 1

?period1 period1 Periodic interval for the first waveform.

?period2 *period2* Periodic interval for the second waveform.

?multiple multiple

Finds all the cross points of wf1. If td2 is specified, finds cross points of target (wf2) staring at trigger event(first cross point of wf1) and ignore wf2 before wf1 trigger event. If td2 is not specified, finds cross points of target (wf2) at time at t=0. Calculate the delay between each of the cross points falling at interval of period1 and period2 of wf1 and wf2 respectively.

?xName xName Specifies whether you want to retrieve delay data

against *trigger* time, *target* time (or another X-axis parameter for non-transient data) or *cycle*. Cycle numbers refer to the n'th occurrence of the delay event

in the input waveform.

The value in this field is ignored when you specify

Number of Occurences as single.

@Rest args Variable list of arguments passed to the delay function

(as created from the Calculator UI). These variables also include support for multiple occurrences of the

delay event.

**Note:**  $g_histoDisplay$  and  $x_noOfHistoBins$  are added for backward compatibility only. It will be deprecated in future releases. Use the histo function for plotting the histogram of the resulting function.

#### Predefined and Waveform (Calculator) Functions

#### Value Returned

o_waveform	Returns a waveform representing the delay if the input argument is a family of waveforms.
n_value	Returns the delay value if the input argument is a single waveform.
nil	Returns nil and an error message otherwise.

### **Example**

```
delay(?wf1 v("16" ?result "tran"), ?value1 2.5, ?edge1 "rising", ?nth1 1, ?td1 0.0,
?wf2 v("1" ?result "tran"), ?value2 2.5, ?edge2 "rising", ?nth2 1, ?td2 nil , ?stop
nil, ?period1 1 ?period2 1 ?multiple t ?xName "trigger" )
```

Calculates the delay between two waveforms wf1 and wf2 with the argument values specified as above.

```
delay( ?wf1 wf1 ?value1 2.5 ?nth1 2 ?edge1 'either ?wf2 wf2 ?value2 2.5 ?nth2 1
?edge2 'falling )
```

Calculates the delay starting from the time when the second edge of the first waveform reaches the value of 2.5 to the time when the first falling edge of the second waveform crosses 2.5.

```
delay( ?td1 5 ?wf2 wf2 ?value2 2.5 ?nth2 1 ?edge2 'rising ?td2 5)
```

Calculates the delay starting when the time equals 5 seconds and stopping when the value of the second waveform reaches 2.5 on the first rising edge 5 seconds after the trigger.

```
delay( ?wf1 wf1 ?value1 2.5 ?nth1 1 ?edge1 'rising ?td1 5 ?wf2 wf2 ?value2 2.5 ?nth2
1 ?edge2 'rising ?td2 0)
```

Waits until after the time equals 5 seconds, and calculates the delay between the first and the second rising edges of wf2 when the voltage values reach 2.5.

```
delay(VT("/out"), 2.5, 1, 'rising, VT("/in"), 2.5, 1, 'rising', 1, 1, t)
```

Computes the delay between the rising edges of VT("/out") and VT("/in") when the waveforms cross their respective threshold values (that is, 2.5).

```
delay(VT("/out") 1.5 1 "rising" VT("/out") 1.5 2 "rising" 1 1 t "trigger") (s)
```

Returns multiple occurrences of delay specified against trigger time-points at which each delay event occurs.

```
delay(VT("/out") 1.5 1 "rising" VT("/out") 1.5 2 "rising" 1 1 t "target") (s)
```

Returns multiple occurrences of delay specified against target time-points at which each delay event occurs.

```
delay(VT("/out") 1.5 1 "rising" VT("/out") 1.5 2 "rising" 1 1 t "cycle") (s)
```

# Predefined and Waveform (Calculator) Functions

Returns multiple occurrences of delay specified against cycle numbers, where a cycle number refers to the n'th occurrence of the delay event in the input waveform.

# Predefined and Waveform (Calculator) Functions

# delayMeasure

```
delayMeasure(
    o_waveform1
    o_waveform2
    [ ?edge1 s_edge1 ]
    [ ?nth1 x_nth1 ]
    [ ?value1 n_value1 ]
    [ ?edge2 s_edge2 ]
    [ ?nth2 x_nth2
    [ ?value2 n_value2 ]
    => n_value / nil
```

## **Description**

Calculates the delay between a trigger event and a target event.

## Predefined and Waveform (Calculator) Functions

# **Arguments**

?wf1 o_waveform1	First waveform object.
?wf2 o_waveform2	Second waveform object.
?edge1 <i>s_edge1</i>	Type of the edge that must cross $n_value1.$ Valid values: rising, falling, either
	Default Value: either
?nth1	Number that specifies which crossing is to be the trigger event. For example, if $x_nth1$ is 2, the trigger event is the second edge of the first waveform with the specified type that crosses $n_value1$ .
	Default Value: 1
?value1 <i>n_value1</i>	Threshold value at which the crossing is significant for the first waveform object. If this value is nil or blank, threshold is calculated internally using average (wavel)
?edge2 <i>s_edge2</i>	Type of the edge for the second waveform.
	Valid values: rising, falling, either
	Default Value: either
?value2 <i>n_value2</i>	Threshold value at which the crossing is significant for the second waveform. If this value is nil or blank, threshold is calculated internally using average (wave2).
?nth2 x_nth2	Number that specifies which crossing is to be the target event. For example, if $x\_nth2$ is 2, the target event is the second edge of the second waveform with the specified type that crosses $n\_value2$ .
	Default Value: 1

## Example

delayMeasure(wave1 wave2)

Calculates the delay between the two waveforms, wave1 and wave2.

delayMeasure( wave1 wave2 ?value1 2.5 ?nth1 2 ?edge1 'either ?wf2 wf2 ?value2 2.5
?nth2 1 ?edge2 'falling )

## Predefined and Waveform (Calculator) Functions

Calculates the delay starting from the time when the second edge of the first waveform reaches the value of 2.5 to the time when the first falling edge of the second waveform crosses 2.5.

### Predefined and Waveform (Calculator) Functions

#### deriv

```
deriv(
    o_waveform
)
=> o waveform / nil
```

# **Description**

Computes the derivative of a waveform with respect to the X axis.

Note the following:

- After the second derivative, the results become inaccurate because the derivative is obtained numerically.
- Use the magnitude value instead of dB in frequency domain.

### **Arguments**

o_waveform	Waveform object representing simulation results that can be
	displayed as a series of points on a grid. (A waveform object
	identifier looks like this: srrWave: XXXXX.)

#### **Value Returned**

o_waveform	Returns a waveform object representing the derivative with respect to the X axis of the input waveform. Returns a family of waveforms if the input argument is a family of waveforms.
nil	Returns nil and an error message otherwise.

## **Example**

```
plot( deriv( VT( "/net8" ) ) )
```

Plots the waveform representing the derivative of the voltage of "/net8".

```
plot( deriv( mag(VF( "/OUT" ) ) )
```

Plots the waveform representing the derivative of the frequency of "/OUT".

## Predefined and Waveform (Calculator) Functions

#### dft

## **Description**

Computes the discrete Fourier transform and fast Fourier transform of the input waveform.

The waveform is sampled at the following n timepoints:

```
from, from + deltaT, from + 2 * deltaT,...,
from + (N - 1) * deltaT
```

The output of dft is a frequency waveform, W(f), which has (N/2 + 1) complex values—the DC term, the fundamental, and (N/2 - 1) harmonics.

**Note:** The last time point, (from + (N - 1) \* deltaT), is (to - deltaT) rather than to. The dft command assumes that w(from) equals w(to).

The DFT function assumes that w(from) equals to w(to). A warning message appears when w(from) is not equal to w(to) in the following situation:

window function is Rectangular or not specified, and

```
|w(from) -w(to)| >1e-3*range(w),
where, range(w) is max(w) - min(w) in [from,to]
```

#### Predefined and Waveform (Calculator) Functions

### **Arguments**

o\_waveform Waveform object representing simulation results that can be

displayed as a series of points on a grid. (A waveform object

identifier looks like this: srrWave: XXXXX.)

 $n_from$  Starting value for the dft computation.

*n to* Ending value for the dft computation.

 $x_num$  Number of timepoints.

t windowName Variable representing different methods for taking a dft

computation.

Valid values: Rectangular, ExtCosBell, HalfCycleSine,

Hanning Or Cosine2, Triangle Or Triangular,

Half3CycleSine Or HalfCycleSine3, Hamming, Cosine4, Parzen, Half6CycleSine Or HalfCycleSine6, Blackman,

Kaiser, or Nuttall.

For more information about windowName, see the information about Discrete Fourier Transform (dft) in the <u>Virtuoso Analog</u>

Design Environment L User Guide.

n\_param1 Smoothing parameter.

Applies only if the t windowName argument is set to Kaiser.

n adcSpan Specifies the peak saturation level of the FFT waveform. When

specified the magnitude of the input waveform is divided by adc

span value before computing FFT. This is full-scale span

ignoring any DC offsets.

Valid values: Any floating point number

Default Value: 1.0

#### Value Returned

waveform
 Returns a waveform representing the magnitude of the various

harmonics for the specified range of frequencies. Returns a family of waveforms if the input argument is a family of

waveforms.

nil Returns nil and an error message otherwise.

### **Example**

## Predefined and Waveform (Calculator) Functions

```
plot( dft( v( "/net8" ) 10u 20m 64 "rectangular" ) )
```

Computes the discrete Fourier transform, fast Fourier transform, of the waveform representing the voltage of "/net8". The computation is done from 10u to 20m with 64 timepoints. The resulting waveform is plotted.

# Predefined and Waveform (Calculator) Functions

## dftbb

```
dftbb(
    o_waveform1
    o_waveform2
    f_timeStart
    f_timeEnd
    x_num
    [ ?windowName t_windowName ]
    [ ?smooth x_smooth ]
    [ ?cohGain f_cohGain ]
    [ ?spectrumType s_spectrumType ]
    )
    => o_waveformComplex / nil
```

# **Description**

Computes the discrete Fourier transform (fast Fourier transform) of a complex signal.

# Predefined and Waveform (Calculator) Functions

# **Arguments**

o_waveform1	Time domain waveform object with units of volts or amps.
o_waveform2	Time domain waveform object with units of volts or amps.
f_timeStart	Start time for the spectral analysis interval. Use this parameter and $f_{timeEnd}$ to exclude part of the interval. For example, you might set these values to discard initial transient data.
$f\_$ time $End$	End time for the spectral analysis interval.
x_num	The number of time domain points to use. The maximum frequency in the Fourier analysis is directly proportionate to $x_{num}$ and inversely proportional to the difference between $f_{timeStart}$ and $f_{timeEnd}$ .
<pre>?windowName t_windowName</pre>	The window to be used for applying the moving window FFT.
	Valid values: Rectangular, ExtCosBell, HalfCycleSine, Hanning, Cosine2, Triangle Or Triangular, Half3CycleSine Or HalfCycleSine3, Hamming, Cosine3, Cosine4, Parzen, Half6CycleSine Or HalfCycleSine6, Blackman, Kaiser, Or Nuttall.
	Default value: Rectangular.
?smooth x_smooth	The Kaiser window smoothing parameter. If there are no requests, there is no smoothing.
	Valid values: 0 <= x_smooth <= 15
	Default value: 1
?cohGain <i>f_cohGain</i>	A scaling parameter. A non-zero value scales the power spectral density by 1/(f_cohGain).
	Valid values: 0 <= f_cohGain <= 1. You can use 1 if you do not want the scaling parameter to be used.
	Default value: 1

### Predefined and Waveform (Calculator) Functions

?spectrumType
t\_spectrumType

A string that can be either singleSided or doubleSided. When this option is single-sided, the resultant waveform is only on one side of the y axis starting from 0 to N-1. When it is double-sided, it is symmetric to the Y axis from -N/2 to (N/2) -1.

Default value: SingleSided

#### Value Returned

o\_waveformComplex The discrete Fourier transform for baseband signals of

the two waveforms returned when the command is

successful.

nil Returns nil and an error message otherwise.

# **Example**

dftbb(VT("/net32") VT("/net11") , 0, 16m, 12000, ?windowName 'Hanning,?smooth 1,
?cohGain 1, ?spectrumType "SingleSided")

### Predefined and Waveform (Calculator) Functions

## dnl

```
dnl(
    o_dacSignal
    o_sample | o_pointList | n_interval
    [ ?mode t_mode ]
    [ ?threshold n_threshold ]
    [ ?crossType t_crossType ]
    [ ?delay f_delay ]
    [ ?method t_method ]
    [ ?units x_units ]
    [ ?nbsamples n_nbsamples ]
    )
    => n_dnl / nil
```

## **Description**

Computes the differential non-linearity of a transient simple or parametric waveform.

# Predefined and Waveform (Calculator) Functions

# **Arguments**

o_dacSignal	Waveform for which the differential non-linearity is to be calculated.
o_sample	Waveform used to obtain the points for sampling the dacSignal. These are the points at which the waveform crosses the threshold while either rising or falling (defined by the crossType argument) with the delay added to them.
n_pointList	List of domain values at which the sample points are obtained from the dacSignal.
n_interval	The sampling interval.
?mode t_mode	The mode for calculating the threshold.
	Valid values: auto and user.Default value: auto.
	If set to user, an n_threshold value needs to be provided.
	If set to auto, n_threshold is calculated internally.
?threshold $n\_threshold$	The threshold value against which the differential non-linearity is to be calculated. It needs to be specified only when the <i>mode</i> is selected as user.
?crossType t_crossType	The points at which the curves of the waveform intersect with the threshold. While intersecting, the curve may be either rising or falling.
	Valid values: rising and falling, respectively.
	Default crossType is rising.
?delay f_delay	The delay time after which the sampling begins.
	Valid values: Any valid time value.
	Default value: 0.
?method t_method	The method to be used for calculation.
	Valid values: end (end-to-end) and fit (straight line).
	Default value: end.

#### Predefined and Waveform (Calculator) Functions

?units x units Unit for expressing the output waveform.

Valid values: abs (absolute) and 1sb (multiples of least

significant bit).

Default value: abs.

?nbsamples n nbsamples Number of samples used for calculating the non-linearity.

If not specified, the samples are taken against the entire

data window.

**Note:** For each of the three ways in which the sample points can be specified, only a few of the other optional arguments are meaningful, as indicated below:

- For o\_sample, the arguments  $t_mode$ ,  $n_threshold$ ,  $t_crossType$ ,  $f_delay$ ,  $t_method$ , and  $x_mits$  are meaningful.
- **For** n pointList, the arguments t method and x units are meaningful.
- For n\_interval, the arguments t\_method, x\_units, and n\_nbsamples are meaningful.

#### Value Returned

 $n_dn1$  Returns the differential waveform.

nil Returns nil and an error message otherwise.

#### **Example**

```
dnl( wave1 wave2 ?crossType "rising" ?delay 0.4 )
=> srrWave:175051544
```

Returns the differential non-linearity for wave1 by taking the points at which wave2 crosses the internally calculated threshold while rising as the sample points and adding a delay of 0.4 to them.

## Predefined and Waveform (Calculator) Functions

# dutyCycle

```
dutyCycle(
    o_waveform
    [?threshold n_threshold]
    [?xName t_xName]
    [?outputType t_outputType]
    [?mode mode]
)
    => o_waveform / f_average / nil
```

## **Description**

Computes the duty cycle for a given waveform as a function of time or cycle.

**Note:** Duty cycle is the ratio of the time for which the signal remains 'high' and the time period of the signal.

## **Arguments**

o_waveform	Waveform, expression, or a family of waveforms.
?threshold n_threshold	The threshold value. It needs to be specified only when the $mode$ is selected as user.
?xName t_xName	The X-axis of the output waveform.
	Valid values: time and cycle.
	Default value: time.
<pre>?outputType t_outputType</pre>	Type of output. Valid values: average and plot.
	If set to average, the output is an average value. If set to plot, the output is a waveform.
	In both the cases, the output is expressed in terms of a percentage.
	Default value: plot.

### Predefined and Waveform (Calculator) Functions

?mode t mode

The mode used to calculate the threshold value.

Valid values: auto and user.

Default value: auto.

If you want to specify the threshold value, set the variable to user. If you want Virtuoso Visualization and Analysis XL to calculate the threshold value, set the variable to auto.

The Auto Threshold Value is calculated as the average which is integral of the waveform divided by the X range.

## Predefined and Waveform (Calculator) Functions

## **Value Returned**

o_waveform	Returns a waveform that represents duty cycle as a function of time.
f_average	Returns the average duty cycle value as a percentage.
nil	Returns nil if the duty cycle cannot be calculated.

# **Example**

```
dutyCycle( wave1 )
=> srrWave:175051552
```

Returns the duty cycle as a function of time for the wave wave1.

```
dutyCycle( wave1 ?outputType "average" )
=> 52.1066
```

Returns the average (in percentage) of the duty cycle values for the wave wave1.

#### Predefined and Waveform (Calculator) Functions

#### evmQAM

```
evmQAM(
    o_waveformI
    o_waveformQ
    n_tDelay
    n_samplingT
    x_levels
    g_normalize
    [ ?percent d_percent ]
)
    => o waveform / nil
```

## Description

Processes the I and Q waveform outputs from the transient simulation run to calculate the Error Vector Magnitude (EVM) for multi-mode modulations. The function plots the I versus Q scatterplot. EVM is a useful measurement to describe the overall signal amplitude and phase modulated signal quality. It is based on a statistical error distribution normalized from an ideal digital modulation. Quadrature Amplitude Modulation (QAM) is a typical modulation scheme where EVM is useful. The EVM is calculated by detecting the I and Q signal levels corresponding to the four possible I and Q symbol combinations and calculating the difference between the actual signal level and the ideal signal level.

Note: This function is not supported for families of waveforms.

#### Predefined and Waveform (Calculator) Functions

#### **Arguments**

o\_waveformI The waveform for the I signal.

o waveformQ The waveform for the Q signal.

n tDelay The start time (a numerical value) for the first valid

symbol. This can be obtained from the Waveform Viewer window by recording the time of the first

minimum or first maximum (whichever is earlier) on the

selected signal stream.

n samplingT A sampling time (a numerical value) for the symbol.

Each period is represented by a data rate. The data rate at the output is determined by the particular

modulation scheme being used.

*x* levels The modulation levels.

Valid values: 4, 16, 64, 256

Default value: 4

g normalize An option to see the scatter plot normalized to the ideal

values +1 and -1 (for example, when superimposing scatter plots from different stages in the signal flow, where the levels may be quite different but you want to see relative degradation or improvement in the scatter). This option does not affect the calculation of the EVM

number.

Valid values: nil, t

Default value: t

?percent d percent

#### Value Returned

o waveform Returns a waveform object representing the EVM value

computed from the input waveforms.

nil Returns nil and an error message if the function is

unsuccessful.

#### Example

evmQAM( v("samp\_out\_Q"), v("samp\_out\_I") 1.5u, 181.81n, 4, t )

# Predefined and Waveform (Calculator) Functions

Calculates the EVM value for the modulation level 4 in normalized form.

## Predefined and Waveform (Calculator) Functions

## evmQpsk

```
evmQpsk(
    o_waveform1
    o_waveform2
    n_tDelay
    n_sampling
    g_autoLevelDetect
    n_voltage
    n_offset
    g_normalize
    [?percent d_percent]
)
    => o waveform / nil
```

### **Description**

Processes the I and Q waveform outputs from the transient simulation run to calculate the Error Vector Magnitude (EVM) and plot the I versus Q scatterplot. EVM is a useful measurement to describe the overall signal amplitude and phase modulated signal quality. It is based on a statistical error distribution normalized from an ideal digital modulation. Quadrature Phase Shift Keying (QPSK) is a typical modulation scheme where EVM is useful. The EVM is calculated by detecting the I and Q signal levels corresponding to the four possible I and Q symbol combinations and calculating the difference between the actual signal level and the ideal signal level.

**Note:** This function is not supported for families of waveforms.

#### Predefined and Waveform (Calculator) Functions

### **Arguments**

o\_waveform1 The waveform for the I signal.

o waveform2 The waveform for the Q signal.

n tDelay The start time for the first valid symbol. This can be

obtained from the Waveform Viewer window by recording the time of the first minimum or first maximum (whichever

is earlier) on the selected signal stream.

n sampling A period for the symbol. Each period is represented by a

data rate. The data rate at the output is determined by

the particular modulation scheme being used.

g autoLevelDetect An option to indicate that you want the amplitude

(n\_voltage) and DC offset (n\_offset) to be automatically calculated. Amplitude is calculated by averaging the rectified voltage level of the signal streams and DC offset by averaging the sum of an equal number of positive and negative symbols in each signal stream. These values are used to determine the EVM value. If this value is set to nil, you must specify values for

n voltage and n offset.

Valid values: 'nil, 't

Default value: 't

n voltage The amplitude of the signal.

n offset The DC offset value.

q normalize An option to see the scatter plot normalized to the ideal

values +1 and -1 (for example, when superimposing scatter plots from different stages in the signal flow, where the levels may be quite different but the you want

to see relative degradation or improvement in the

scatter). This option does not affect the calculation of the

EVM number.

Valid values: nil, t

Default value: nil

?percent d percent

## Predefined and Waveform (Calculator) Functions

#### Value Returned

o waveform Returns a waveform object representing the EVM value

computed from input waveforms.

nil Returns nil and an error message if the function is

unsuccessful.

#### **Example**

```
evmQpsk( v("samp_out_Q"), v("samp_out_I") 1.5u, 181.81n, t, nil, nil, nil)
```

Calculates the EVM value when  $g_{autoLevelDetect}$  is set to t. In this case, no values are specified for  $n_{voltage}$  and  $n_{offset}$ .

```
evmQpsk( v("samp out Q"), v("samp out I") 1.5u, 181.81n, nil, 1.3, 0, nil)
```

Calculates the EVM value when  $g_{autoLevelDetect}$  is set to nil. In this case, values are specified for n voltage and n offset.

### Predefined and Waveform (Calculator) Functions

# eyeDiagram

```
eyeDiagram(
    o_waveform
    n_start
    n_stop
    n_period
    [ ?advOptions t_advOptions ]
    [ ?intensityPlot g_intensityPlot ]
    )
    => o_waveform / nil
```

## **Description**

Returns an eye-diagram plot of the input waveform signal. It returns the waveform object of the eye-diagram plot. Using an advanced option, the function also calculates the maximum vertical and horizontal opening of the eye formed when the input waveform is folded by the specified period to form the eye.

## **Arguments**

o_waveform	Input waveform signal.
n_start	The X-axis start value from where the eye-diagram plot is to begin.
n_stop	The X-axis stop value where the eye-diagram plot is to terminate.
n_period	The period after which the waveform is to be folded to form the eye.
?advOptions t_advOptions	Specifies whether the vertical (Max Vertical Opening) or horizontal opening (Max Horizontal Opening) of the eye is to be calculated.
	Valid values: vertical, horizontal
	Default value: nil
?intensityPlot g_intensityPlot	Boolean used to specify whether to generate a high intensity eye diagram plot.

**Note:** If  $t\_advOptions$  is specified, the function approximates vertical eye height and horizontal eye width to assume the symmetry of the eye. The function returns the most optimum results for single-eye scenarios.

## Predefined and Waveform (Calculator) Functions

#### **Value Returned**

o\_waveform Returns a waveform object representing the eye-

diagram plot of the input waveform

nil Returns nil and an error message otherwise

## **Example**

```
eyeDiagram( v("/out" ) On 500n 12.5n )
```

Returns a waveform that represents an eye-diagram plot.

```
eyeDiagram( v("/out" ) On 500n 12.5n ?advOptions "vertical" )
```

Calculates the maximum vertical opening of the eye that is formed when the input waveform is folded after 12.5n

```
eyeDiagram( v("/out" ) On 500n 12.5n ?advOptions "horizontal" )
```

Calculates the maximum horizontal opening of the eye that is formed when the input waveform is folded after 12.5n

## Predefined and Waveform (Calculator) Functions

# eyeHeightAtXY

```
eyeHeightAtXY(
    o_eyeDiagram
    f_x
    f_y
)
=> f eyeHeight / nil
```

## Description

Calculates the eye height at the specified point (x,y) inside the eye diagram. Eye height is the difference of two intercepts made with the innermost traces of the eye in the y-axis direction.

**Note:** The specified point (x,y) must lie within the open eye whose height you want to calculate.

For more information about how the eye height is calculated, see the <u>eyeHeightAtXY</u> calculator function.

## **Arguments**

o_eyeDiagram	The eye diagram waveform that is used to calculate the eye height.
f_x	The x-axis value that is used to calculate the eye height.
f_y	The y-axis value that is used to calculate the eye height.

#### **Values Returned**

f_eyeHeight	Returns the eye height at the specified point $(x,y)$ inside the eye diagram.
nil	Returns nil if there is an error.

### **Example**

The following example calculates the eye height at the point, x = 70p and y = 2.2:

```
eyeHeightAtXY(eyeDiagram(v("/example_1" ?result "tran") 560p 5.000n 140p ?triggerPeriod 7e-11 ?autoCenter t) \overline{7}0p 2.2 ) => 1.494594
```

## Predefined and Waveform (Calculator) Functions

# eyeWidthAtXY

```
eyeWidthAtXY(
    o_eyeDiagram
    f_x
    f_y
)
    => f eyeWidth / nil
```

## Description

Calculates the eye width at the specified point (x,y) inside the eye diagram. Eye width is the difference of two intercepts made with the innermost traces of the eye in the x-axis direction.

**Note:** The specified point (x,y) must lie within the open eye whose width you want to calculate.

For more information about how the eye width is calculated, see the <u>eyeWidthAtXY</u> calculator function.

## **Arguments**

o_eyeDiagram	The eye diagram waveform that is used to calculate the eye width.
f_x	The x-axis value that is used to calculate the eye width.
f_y	The y-axis value that is used to calculate the eye width.

## Values Returned

f_eyeWidth	Returns the eye width at the specified point $(x,y)$ inside the eye diagram.
nil	Returns nil if there is an error.

#### **Example**

The following example calculates the eye height at the point, x = 70p and y = 2.2:

```
eyeWidthAtXY(eyeDiagram(v("/example_1" ?result "tran") 560p 5.000n 140p
?triggerPeriod 7e-11 ?autoCenter t) 70p 2.2 )
=> 2.388933e-11
```

## Predefined and Waveform (Calculator) Functions

# eyeAperture

```
eyeAperture(
    o_waveform
    f_vref
    f_acHeight
    f_dcHeight
    g_plotBox
    ?optimize g_optimize
)
=> o waveform / aperture width / nil
```

# **Description**

Returns the aperture of the input eye diagram signal.

#### Predefined and Waveform (Calculator) Functions

### **Arguments**

o waveform

_	aperture is to be calculated.
f_vref	Reference voltage.
f_acHeight	AC height, which specifies the height of the left side of the aperture window.
f_dcHeight	DC height, which specifies the height of the right side of

the aperture window.

 $g_plotBox$  Specifies whether to display the aperture in the eye

diagram or calculate the eye width.

Valid values: t or nil. When this argument is set to t, the eye aperture is displayed in the output plot. When set to

Input signal, which is a eye diagram waveform for which

nil, the eye aperture width is returned.

?optimize  $g_{optimize}$  Specifies whether to calculate the reference voltage that

can be used to achieve the maximum eye aperture width.

Valid values: t or nil.

#### **Values Returned**

o_waveform	Returns the output waveform with eye aperture plotted.
aperture_width	Returns the aperture width.
nil	Returns nil and an error message otherwise.

## **Example**

This example calculates the eye aperture on the two eye diagram signals, signal1 and signal2, with following values:

- $\blacksquare$  vref=0.75
- acHeight=0.6
- $\blacksquare$  dcHeight=0.5
- plotbox=t (to plot the eye aperture)

### Predefined and Waveform (Calculator) Functions

# eyeMeasurement

```
eyeMeasurement(
    o waveform
    [ ?sample n_sample ]
    [ ?auto q auto ]
    [ ?horizThreshold n horizThreshold ]
    [ ?sample n sample ]
    [ ?xTypePercent0 g xTypePercent0 ]
    [ ?startx0 n_startx0 ]
    [ ?starty0 n_starty0 ]
    [ ?yTypePercent0 g_yTypePercent0 ]
    [ ?endx0 n endx0 ]
    [ ?endy0 n endy0 ]
    [ ?xTypePercent1 g_xTypePercent1 ]
    [ ?startx1 n_startx1 ]
    [ ?starty1 n_starty1 ]
    [ ?yTypePercent1 g yTypePercent1 ]
    [ ?endx1 n_endx1 ]
    [ ?endy1 n endy1 ]
    [ ?noOfBins n_noOfBins ]
    [ ?measure t measure ]
    => o waveform / nil
```

# Description

Evaluates the measurements for the eye diagram plot.

#### Predefined and Waveform (Calculator) Functions

#### **Arguments**

o waveform

?sample n sample

?auto g\_auto

?horizThreshold
n\_horizThreshold

?sample n sample

?xTypePercent0
q xTypePercent0

?startx0 n startx0

The eye diagram waveform.

The time interval after which the signals are divided in the eye diagram plot. If this field is left blank, the data within the level 1 and level 0 regions are used to analyze the amplitude variation of the signal. This means there is some sensitivity to the actual spacing between the data points in the signal, which is caused by the variable time steps in the simulator. If the points are clustered in the curve portion, the distribution can be skewed. To perform the analysis, the sampling interval you specify in this field is divided into even time points.

When this argument is set to true, then the following agument values are computed automatically:

horizThreshold, startx0, startx1, xTypePercent0, xTypePercent1, endx0, endx1, yTypePercent0, yTypePercent1, starty0, endy0, starty1, and endy1.

The Y-axis level (for example voltage) that represents the switching threshold of the signal, typically half the signal range. This is used to compute statistical information about the threshold.

The time interval after which the signals are divided in the eye diagram plot. If this field is left blank, the data within the level 1 and level 0 regions are used to analyze the amplitude variation of the signal. This means there is some sensitivity to the actual spacing between the data points in the signal, which is caused by the variable time steps in the simulator. If the points are clustered in the curve portion, the distribution can be skewed. To perform the analysis, the sampling interval you specify in this field is divided into even time points.

Level0 X-range specified whether specified in "%". If the value is t, it signifies the "%" value and if the value is nil, it signifies the absolute value.

Default value is t.

Level0 X-range start value. Default value is 40.

# Predefined and Waveform (Calculator) Functions

?starty0 n_starty0	Level0 Y-range start value. Default value is 0.
<pre>?yTypePercent0 g_yTypePercent0</pre>	Level0 Y-range specified whether specified in "%". If the value is $t$ , it signifies the "%" value and if the value is nil, it signifies the absolute value. Default value is $t$ .
?endx0 n_endx0	Level0 X-range end value. Default value is 60.
?endy0 n_endy0	Level0 Y-range end value. Default value is 50.
?xTypePercent1 g_xTypePercent1	Level1 X-range specified whether specified in "%". If the value is $t$ , it signifies the "%" value and if the value is nil, it signifies the absolute value. Default value is $t$ .
?startx1 n_startx1	Level1 X-range start value. Default value is 40.
?starty1 <i>n_starty1</i>	Level1 Y-range start value. Default value is 50.
<pre>?yTypePercent1 g_yTypePercent1</pre>	Level1 Y-range specified whether specified in "%". If the value is $t$ , it signifies the "%" value and if the value is nil, it signifies the absolute value. Default value is $t$ .
?endx1 n_endx1	Level1 X-range end value. Default value is 60.
?endy1 <i>n_endy1</i>	Level1 Y-range end value. Default value is 100.
?noOfBins n_noOfBins	Number of signal bins to be displayed in the eye diagram plot. These signals bins are used to form the horizontal (threshold crossing times) and vertical (amplitude variation) histograms.

### Predefined and Waveform (Calculator) Functions

?measure t measure

Computes one of the measurement values described below:

- level0Mean (Level0 Mean)—Mean of the Y-values within the level0 region.
- level0Stddev (Level0 Stddev)—Standard deviation of the Y-values within the level0 region.
- level1Mean (Level1 Mean)—Mean of the Y-values within the level1 region.
- level1Stddev (Level1 Stddev)—Standard deviation of the Y-values within the level1 region.
- amplitude (Eye amplitude)—Mean to mean amplitude of the eye, computed as: Meanlevel1 - Meanlevel0.
- height (Eye height)—Vertical opening of the eye, computed as:

```
(Meanlevel1 - 3*level1) -
(Meanlevel0 - 3*level0).
```

signalToNoise (Eye signalToNoise)—Signal
to noise ratio of the eye, computed as:
 (Meanlevel1 - Meanlevel0) / (level1
+ level0).

- thresholdCrossingStddev (Threshold crossing stddev)—Threshold crossing standard deviation is computed only when there is a single transition region in the eye diagram because it is analyzed over the entire period.
- thresholdCrossingAverage (Threshold crossing average)—This is computed over the entire period.
- width (Eye width)—Represents the opening of the eye in the X direction. It is computed as:

```
(Meantransition2 -
3*std(transition2)) -
(Meantransition1 -
3*std(transition1)).
```

### Predefined and Waveform (Calculator) Functions

- riseTime (Eye Rise Time)—Two thresholds taken at the 20% and 80% points between the level0 mean and level1 mean. At each of these two thresholds, a horizontal histogram is computed, which is an analysis of the crossing points of these two thresholds, and the resulting rise time is the difference in the mean crossing point at each of these two thresholds.
- fallTime (Eye Fall Time)—Signal measured between the percent high and percent low of the difference between the initial and final value.
- randJitterLeft—Random jitter calculated from the crossing histogram of the left crossing area
- randJitterRight—Random jitter calculated from the crossing histogram of the right crossing area
- determJitter—Average deterministic jitter of the crossing areas

#### Value Returned

o waveform

nil

Returns the computed scalar value or a waveform for the specific measure that was passed.

Returns nil and an error message otherwise.

## Example

The following function computes the threshold crossing average for the eye diagram for signal /net1 from 0.0 to 2e-08 with a period of 5e-10:

```
wave_eyeDiagram=eyeDiagram(vtime('tran "/net1") 0.0 2e-08 5e-10 ?intensityPlot t)
eyeMeasurement(wave_eyeDiagram ?horizThreshold 0 ?startx0 40 ?starty0 0
?xTypePercent0 "t" ?endx0 60 ?endy0 50 ?yTypePercent0 "t" ?startx1 40 ?starty1 50
?xTypePercent1 "t" ?endx1 60 ?endy1 100 ?yTypePercent1 "t" ?noOfBins 10 ?measure
"thresholdCrossingAverage")
```

### The following function is used to get the eye open width:

```
wave eyeDiagram=eyeDiagram(vtime('tran "/net1") 0.0 2e-08 5e-10 ?intensityPlot t)
```

## Predefined and Waveform (Calculator) Functions

eyeMeasurement(wave\_eyeDiagram ?horizThreshold 0 ?startx0 40 ?starty0 0 ?xTypePercent0 "t" ?endx0 60 ?endy0 50 ?yTypePercent0 "t" ?startx1 40 ?starty1 50 ?xTypePercent1 "t" ?endx1 60 ?endy1 100 ?yTypePercent1 "t" ?noOfBins 10 ?measure "width")

## Predefined and Waveform (Calculator) Functions

# edgeTriggeredEyeDiagram

```
edgeTriggeredEyeDiagram(
    o_waveform
    n_start
    n_stop
    o_triggerWave
    n_threshold
    s_edgeType
    n_triggerOffset
    [?intensityPlot g_intensityPlot]
)
    => o waveform / nil
```

## **Desription**

Returns a signal triggered at the beginning of the eye diagram instead of a fixed period.

### Predefined and Waveform (Calculator) Functions

### **Arguments**

0	waveform	The eye diagram waveform.

n start The X-axis start value from where the eye diagram plot

is to begin.

n stop The X-axis stop value where the eye diagram plot is to

terminate.

o\_triggerWave The waveform that is used for triggering the eye

diagram.

n threshold The Y-axis value of trigger wave at which the

corresponding cross points of the trigger wave are

calculated.

s edgeType Type of the crossing.

Valid values: rising, falling, either.

n triggerOffset The value by which the trigger wave should be l-shifted

to align with the input waveform signal.

?intensityPlot g\_intensityPlot

Controls the intensity based plotting of the eye

diagram.

#### Value Returned

o waveform Returns the computed scalar value or a waveform for

the specific measure that was passed.

nil Returns nil and an error message otherwise.

#### **Examples**

In the following example VT("/out") is an input waveform for which eye diagram is to be determined from 0n to 10n. The period to wrap or fold the eye diagram is determined by the cross points of the trigger waveform VT("/clk") at the given threshold.

```
edgeTriggeredEyeDiagram(VT("/out") On 10n VT("/clk") 2.5 "either" On)
```

The above function returns a waveform with the relevant edge Trigger eye diagram attributes set so that when plotted the edge trigger eye diagram is displayed.

### Predefined and Waveform (Calculator) Functions

The following example shows that an offset of ln signifies that VT("/clk") is be l-shifted by ln, lshift(VT("/clk") ln), before determining the cross points. Also, intensity-based plotting is turned on.

edgeTriggeredEyeDiagram(VT("/out") 0n 10n VT("/clk") 2.5 "rising" 1n
?intensityPlot t)

## Predefined and Waveform (Calculator) Functions

# flip

```
flip(
    o_waveform
)
    => o waveform / nil
```

## **Description**

Returns a waveform with the X vector values negated.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be
	displayed as a series of points on a grid. (A waveform object
	identifier looks like this: srrWaye : XXXXX)

#### Value Returned

o_waveform	Returns a waveform object representing the input waveform mirrored about its Y axis. Returns a family of waveforms if the input argument is a family of waveforms.
nil	Returns nil and an error message otherwise.

## **Example**

```
plot( flip( v("/net4") ) )
```

Plots the waveform for the voltage of "/net4" with the X vector values negated.

## Predefined and Waveform (Calculator) Functions

## fourEval

```
fourEval(
    o_waveform
    n_from
    n_to
    n_by
    [ ?baseBand g_baseBand ]
    )
    => o_waveform / nil
```

## **Description**

Evaluates the Fourier series represented by an expression.

This function is an inverse Fourier transformation and thus the inverse of the <u>dft</u> command. The fourEval function transforms the expression from the frequency domain to the time domain.

## Predefined and Waveform (Calculator) Functions

### **Arguments**

o waveform Waveform object representing simulation results that

can be displayed as a series of points on a grid. (A

waveform object identifier looks like this:

srrWave: XXXXX.)

n from Starting point on the X axis at which to start the

evaluation.

n\_to Increment.

n by Ending point on the X axis.

?baseBand g baseBand Accepts boolean values t or nil. The default value is

nil.

When set to t, the function evaluates the baseband version of the inverse of the dft function by converting the unsymmetrical spectrum to a symmetrical one.

#### Value Returned

o waveform Returns a waveform object representing the inverse

Fourier transformation of the input waveform. Returns a family of waveforms if the input argument is a family of waveforms. Returns the baseband version of the inverse of the  $\mathtt{dft}$  function if  $\mathtt{baseBand}$  is set to  $\mathtt{t}$ .

nil Returns nil and an error message otherwise.

## Predefined and Waveform (Calculator) Functions

## Example

```
plot( fourEval( v( "/net3" ) 1k 10k 10 )
```

Plots the waveform representing the inverse Fourier transformation of the voltage of "/net3" from 1k to 10k.

### Predefined and Waveform (Calculator) Functions

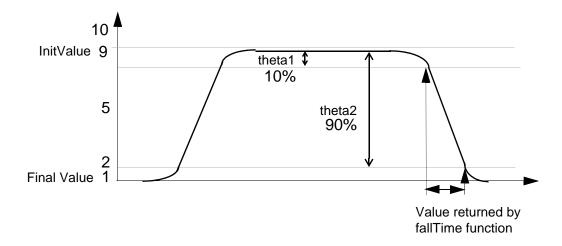
### fallTime

```
fallTime(
    o_waveform
    n_initVal
    g_initType
    n_finalVal
    g_finalType
    n_theta1
    n_theta2
    [ g_multiple [ s_Xname ][ g_histoDisplay ][ x_noOfHistoBins ] ]
    )
    => o waveform / n value / nil
```

## **Description**

Returns the fall time measured between theta1 (percent high) to theta2 (percent low) of the difference between the initial value and the final value.

The fallTime function can also be used to compute the rise time if <code>initVal</code> is lower than <code>finalVal</code>.



## Predefined and Waveform (Calculator) Functions

# **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_initVal	Initial value at which to start the computation.
g_initType	Specifies how n_initVal functions.
	Valid values: a non-nil value specifies that the initial value is taken to be the value of the waveform, interpolated at $n\_initVal$ , and the waveform is clipped from below as follows:
	<pre>o_waveform = clip( o_waveform g_initVal nil )</pre>
	where nil specifies that $n\_initVal$ is defined by the X value entered. (The command gets the Y value for the specified X value and uses that value for $n\_initVal$ .)
n_finalVal	Final value at which to end the computation.
$g\_finalType$	Specifies how the $n_finalVal$ argument functions.
	Valid values: a non-nil value specifies that the final value is taken to be the value of the waveform, interpolated at $n\_finalVal$ , and the waveform is clipped from above, as follows:
	<pre>o_waveform = clip(o_waveform nil n_finalVal)</pre>
	where nil specifies that the $n\_finalVal$ argument is defined by the X value entered. (The command gets the Y value for the specified X value and uses that value for $n\_finalVal$ .)
n_theta1	Percent high.
n_theta2	Percent low.
g_multiple	An optional boolean argument that takes the value $\mathtt{nil}$ by default. If set to $\mathtt{t}$ , the function returns multiple occurrences of the fallTime event.
s_xName	An optional argument that is used only when $g_{multiple}$ is set to t. It takes the value time by default. It controls the contents of the x vector of the waveform object returned by the function.
	Valid values: 'time, 'cycle

### Predefined and Waveform (Calculator) Functions

 $g_histoDisplay$  When set to t, returns a waveform that represents the

statistical distribution of the fallTime data in the form of a histogram. The height of the bars (bins) in the histogram

represents the frequency of the occurrence of values within the

range of fallTime data.

Valid values: t nil
Default value: nil

x noOfHistoBins Denotes the number of bins represented in the histogram

representation. Valid values: Any positive integer Default value:

nil

**Note:**  $g_histoDisplay$  and  $x_noOfHistoBins$  are added for backward compatibility only. It will be deprecated in future releases. Use the histo function for plotting the histogram of the resulting function.

#### Value Returned

o_waveform	Returns a waveform representing the fall time for a family of
------------	---

waveforms if the input argument is a family of waveforms or if

g multiple is set to t.

n value Returns a value for the fall time if the input is a single waveform.

nil Returns nil and an error message otherwise.

### **Example**

```
fallTime( v( "/net8" ) 9 nil 1 nil 10 90 )
```

Computes the fall time for the waveform representing the voltage of "/net8" from 9 to 1.

## Predefined and Waveform (Calculator) Functions

## freq

```
freq(
    o_waveform
    t_crossType
    [?threshold n_threshold]
    [?mode t_mode]
    [?xName xName]
    [?histoDisplay g_histoDisplay]
    [?noOfHistoBins x_noOfHistoBins]
)
    => o_outputWave / nil
```

## **Description**

Computes the frequency of the input waveform(s) as a function of time or cycle.

## Predefined and Waveform (Calculator) Functions

Waveform, expression, or a family of waveforms.

statistical distribution of the riseTime data in the form of a histogram. The height of the bars (bins) in the

occurrence of values within the range of riseTime

histogram represents the frequency of the

### **Arguments**

o waveform

g histoDisplay

t_crossType	The points at which the curves of the waveform intersect with the threshold. While intersecting, the curve may be either rising or falling. For the freq function, you may specify the frequency to be calculated against either the rising points or the falling points by setting crossType to rising or falling, respectively. The default crossType is rising.
?threshold n_threshold	The threshold value against which the frequency is to be calculated. This needs to be specified only when the mode selected is user.
?mode t_mode	The mode for calculating the threshold. This is $\mathtt{auto}$ , by default, in which case $\mathtt{n\_threshold}$ is calculated internally. It can alternatively be set to $\mathtt{user}$ , in which case, an $\mathtt{n\_threshold}$ value needs to be provided. Default Value: $\mathtt{auto}$
?xName t_xName	The X-axis of the output waveform. The default value is time but cycle is also a valid value.
	Default Value: time
?histoDisplay	When set to t, returns a waveform that represents the

Valid values: t nil
Default value: nil

?noOfHistoBins Denotes the number of bins represented in the  $x_noOfHistoBins$  bistogram representation.

data.

Valid values: Any positive integer

Default value: 1

**Note:**  $g_histoDisplay$  and  $x_noOfHistoBins$  are added for backward compatibility only. It will be deprecated in future releases. Use the histo function for plotting the histogram of the resulting function.

## Predefined and Waveform (Calculator) Functions

### **Value Returned**

o\_outputWave Returns the frequency as a function of time or cycle.

nil Returns nil if the frequency cannot be calculated.

## **Example**

```
freq( wave1 "rising" ?mode "user" ?threshold 18.5 ?xName "cycle" )
=> srrWave: 170938688
```

Returns the frequency for wave1 with the threshold at 18.5 against cycle on the x-axis.

## Predefined and Waveform (Calculator) Functions

# freq\_jitter

```
freq_jitter(
    o_waveform
    t_crossType
    [?mode t_mode]
    [?threshold n_threshold]
    [?binSize n_binSize]
    [?xName t_xName]
    [?outputType t_outputType]
)
    => o_waveform / f_val / nil
```

## **Description**

Calculates the frequency jitter.

### Predefined and Waveform (Calculator) Functions

### **Arguments**

o waveform, expression, or a family of waveforms.

t crossType The points at which the curves of the waveform intersect

with the threshold. While intersecting, the curve may be

either rising or falling.

Valid values: rising and falling.

Default value: rising

?mode t mode The mode for calculating the threshold.

Valid values: auto and user.

If set to user, an n threshold value needs to be

provided.

If set to auto, n threshold is calculated internally.

Default value: auto

?threshold n threshold The threshold value against which the frequency is to be

calculated. It needs to be specified only when the mode

selected is user.

?binSize n binSize The width of the moving average window. The deviation of

value at the particular point from the average of this

window is the litter.

Default value: 0

?xName t xName The X-axis of the output waveform. Valid values: time

and cycle. Default value: time

?outputType Type of output.

t\_outputType Valid values: sd and plot.

If set to sd, the output is a standard deviation jitter.

If set to plot, the output is a waveform.

Default value: plot

#### Value Returned

o waveform Returns the frequency jitter values as a function of time or

cycle when the output Type is set to plot.

## Predefined and Waveform (Calculator) Functions

f val Returns the standard deviation value when the

output Type is set to sd.

nil Returns nil otherwise.

## **Example**

```
freq_jitter( wavel "rising" ?mode "user" ?threshold 1 ?binSize 2 ?xName "cycle"
?outputType "sd" )
=> 0.1338585
```

Returns the standard deviation for the frequency jitter of wave1 with the threshold of 1 against the cycle on the x-axis.

## Predefined and Waveform (Calculator) Functions

## frequency

```
frequency(
    o_waveform
)
=> o_waveform / n_value / nil
```

## **Description**

Computes the reciprocal of the average time between two successive midpoint crossings of the rising waveform.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be
	displayed as a series of points on a grid. (A waveform object
	identifier looks like this: srrWave: XXXXX.)

#### Value Returned

o_waveform	Returns a waveform representing the frequency of a family of waveforms if the input argument is a family of waveforms.
n_value	Returns a number representing the frequency of the specified waveform.
nil	Returns nil and an error message otherwise.

## **Example**

```
frequency( v( "/net12" ) )
```

Returns the frequency of "/net12".

## Predefined and Waveform (Calculator) Functions

### ga

```
ga(
    o_s11
    o_s12
    o_s21
    o_s22
    [?gs n_gs]
)
    => o_waveform / nil
```

## **Description**

Returns the available gain in terms of the supplied parameters and the optional source reflection coefficient (Gs).

## **Arguments**

o_s11	Waveform object representing s11.
o_s12	Waveform object representing s12.
o_s21	Waveform object representing s21.
o_s22	Waveform object representing s22.
n_gs	Source reflection coefficient.
	Default value: 0

#### Value Returned

o_waveform	Waveform object representing the available gain.
nil	Returns nil and an error message otherwise.

```
s11 = sp(1 1)
s12 = sp(1 2)
s21 = sp(2 1)
s22 = sp(2 2)
plot(ga(s11 s12 s21 s22))
```

## Predefined and Waveform (Calculator) Functions

## gac

```
gac(
    o_s11
    o_s12
    o_s21
    o_s22
    g_level
    g_frequency
)
    => o waveform / nil
```

## **Description**

Computes the available gain circles.

The g data type on  $g\_level$  and  $g\_frequency$  allows either the level or the frequency to be swept while the other remains fixed.

### Predefined and Waveform (Calculator) Functions

Waveform object representing s11.

### **Arguments**

o s11

o_s12	Waveform object representing s12.
o_s21	Waveform object representing s21.
o_s22	Waveform object representing s22.
g_level	Level in dB. It can be specified as a scalar or a vector. If it is specified as a vector, the level is swept. The $linRg$ function can be called to generate a linear range. For example, $linRg(-30\ 30\ 5)$ is the same as $list(-30\ -25\ -20\ -15\ -10\ -5\ 0\ 5\ 10\ 15\ 20\ 25\ 30)$ and the $g\_level$ argument can be specified as either of the above. In that case, an available gain circle is calculated at each one of the 13 levels.
g_frequency	Frequency, which can be specified as a scalar or a linear range. If it is specified as a linear range, the frequency is swept. The linear range is specified as a list with three values: the start of

the range, the end of the range, and the increment. For example, <code>list(100M 1G 100M)</code> specifies a linear range with the following values:

{ 100M, 200M, 300M, 400M, 500M, 600M, 700M, 800M, 900M, 1G }

In that case, an available gain circle is calculated at each one of the 10 frequencies.

#### Value Returned

o\_waveform Waveform object representing the available gain circles.

nil Returns nil and an error message otherwise.

```
s11 = sp(1 1 ?result "sp")
s12 = sp(1 2 ?result "sp")
s21 = sp(2 1 ?result "sp")
s22 = sp(2 2 ?result "sp")
plot(gac(s11 s12 s21 s22 linRg(-30 30 5) 900M))
```

### Predefined and Waveform (Calculator) Functions

## gainBwProd

```
gainBwProd(
    o_waveform
)
=> o waveform / n value / nil
```

## **Description**

Calculates the gain-bandwidth product of a waveform representing the frequency response of interest over a sufficiently large frequency range.

Returns the product of the zero-frequency-gain and 3dB-gain-frequency.

.

gainBwProd (gain) = 
$$A_0 * f2$$

The gain-bandwidth product is calculated as the product of the DC gain  $A_o$  and the critical frequency f2. The critical frequency f2 is the smallest frequency for which the gain equals  $1/\sqrt{2}$  times the DC gain  $A_o$ .

### Predefined and Waveform (Calculator) Functions

### **Arguments**

o\_waveform Waveform object representing simulation results that can be

displayed as a series of points on a grid. (A waveform object

identifier looks like this: srrWave: XXXXX.)

#### **Value Returned**

o waveform Returns a waveform representing the gain-bandwidth product

for a family of waveforms if the input argument is a family of

waveforms.

n value Returns a value for the gain-bandwidth product for the specified

waveform.

nil Returns nil and an error message otherwise.

### **Example**

```
gainBwProd( v( "/OUT" ) )
```

Returns the gain-bandwidth product for the waveform representing the voltage of the "/OUT" net.

### Predefined and Waveform (Calculator) Functions

## gainMargin

```
gainMargin(
    o_waveform
    [ g_stable ]
    )
    => o waveform / n value / nil
```

### **Description**

Computes the gain margin of the loop gain of an amplifier.

The first argument is a waveform representing the loop gain of interest over a sufficiently large frequency range. This command returns the dB value of the waveform when its phase crosses negative pi.

```
gainMargin( gain ) = 20 * log10( value( gain f0 ) )
```

The gain margin is calculated as the magnitude of the gain in dB at f0. The frequency f0 is the lowest frequency in which the phase of the gain provided is -180 degrees. For stability, the gain margin will be negative when g\_stable is set to nil. If g\_stable value is set to t, then a stable design will have a positive value.

### **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
g_stable	Boolean optional value that takes the value nil by default.

#### Value Returned

o_waveform	Returns a waveform representing the gain margin for a family of waveforms if the input argument is a family of waveforms.
n_value	Returns the value for the gain margin of the specified waveform.
nil	Returns nil and an error message otherwise.

```
gainMargin( v( "/OUT" ) ) = -9.234 gainMargin( v( "/OUT" ) nil ) = -9.234
```

## Predefined and Waveform (Calculator) Functions

gainMargin(v("/OUT")t) = 9.234

### Predefined and Waveform (Calculator) Functions

## gmax

```
gmax(
    o_s11
    o_s12
    o_s21
    o_s22
)
    => o_waveform / nil
```

## **Description**

Returns the maximum power gain in terms of the supplied parameters.

## **Arguments**

o_s11	Waveform object representing s11.
0_s12	Waveform object representing s12.
0_s21	Waveform object representing s21.
o_s22	Waveform object representing s22.

#### Value Returned

```
o_waveformnilReturns nil and an error message otherwise.
```

```
s11 = sp(1 1)
s12 = sp(1 2)
s21 = sp(2 1)
s22 = sp(2 2)
plot(gmax(s11 s12 s21 s22))
```

### Predefined and Waveform (Calculator) Functions

## gmin

```
gmin(
    o_Gopt
    o_Bopt
    f_zref
)
    => o_gminWave / nil
```

## **Description**

Returns the optimum noise reflection coefficient in terms of  $o\_Gopt$ ,  $o\_Bopt$ , and  $f\_zref$ .

gmin is returned as follows:

## **Arguments**

o_Gopt	Waveform object representing the optimum source conductance.
o_Bopt	Waveform object representing the optimum source susceptance.
f_zref	Reference impedance.

#### Value Returned

o_gminWave	Waveform object representing the optimum noise reflection coefficient.
nil	Returns nil and an error message otherwise.

```
Gopt = getData("Gopt")
Bopt = getData("Bopt")
Zref = zref(1 ?result "sp")
plot(gmin(Gopt Bopt Zref))
```

## Predefined and Waveform (Calculator) Functions

## gmsg

```
gmsg(
    o_s11
    o_s12
    o_s21
    o_s22
)
    => o_waveform / nil
```

## **Description**

Returns the maximum stable power gain in terms of the supplied parameters.

## **Arguments**

o_s11	Waveform object representing s11.
o_s12	Waveform object representing s12.
o_s21	Waveform object representing s21.
0_s22	Waveform object representing s22.

#### Value Returned

o_waveform	Waveform object representing the maximum stable power gain.
nil	Returns nil and an error message otherwise.

```
s11 = sp(1 1)
s12 = sp(1 2)
s21 = sp(2 1)
s22 = sp(2 2)
plot(gmsg(s11 s12 s21 s22))
```

## Predefined and Waveform (Calculator) Functions

## gmux

## **Description**

Returns the maximum unilateral power gain in terms of the supplied parameters.

## **Arguments**

o_s11	Waveform object representing s11.
o_s12	Waveform object representing s12.
o_s21	Waveform object representing s21.
o_s22	Waveform object representing s22.

#### Value Returned

o_waveform	Waveform object representing the maximum unilateral power gain.
nil	Returns nil and an error message otherwise.

```
s11 = sp(1 1)

s12 = sp(1 2)

s21 = sp(2 1)

s22 = sp(2 2)

plot(gmux(s11 s12 s21 s22))
```

### Predefined and Waveform (Calculator) Functions

## gp

```
gp(
    o_s11
    o_s12
    o_s21
    o_s22
    [?gl n_gl]
)
    => o_waveform / nil
```

## **Description**

Computes the power gain in terms of the S-parameters.

### **Arguments**

o_s11	Waveform object representing s11.
o_s12	Waveform object representing s12.
o_s21	Waveform object representing s21.
o_s22	Waveform object representing s22
n gl	Load reflection coefficient. Default value: 0

#### **Value Returned**

o_waveform	Waveform object representing the power gain.
nil	Returns nil and an error message otherwise.

## **Example**

```
s11 = sp(1 1)

s12 = sp(1 2)

s21 = sp(2 1)

s22 = sp(2 2)

plot(gp(s11 s12 s21 s22))
```

**Note:** gl is an imaginary number which should be input in the following format:

```
gp( s11 s12 s21 s22 ?gl complex(<realPart> <imagPart>))
```

## Predefined and Waveform (Calculator) Functions

### gpc

```
gpc(
    o_s11
    o_s12
    o_s21
    o_s22
    g_level
    g_frequency
)
    => o waveform / nil
```

## **Description**

Computes the power gain circles.

The g datatype on  $g\_level$  and  $g\_frequency$  allows either the level or the frequency to be swept while the other remains fixed.

### Predefined and Waveform (Calculator) Functions

### **Arguments**

 $o\_s11$  Waveform object representing s11.  $o\_s12$  Waveform object representing s12.  $o\_s21$  Waveform object representing s21.  $o\_s22$  Waveform object representing s22.  $o\_s22$  Level in dB. It can be specified as a second content of the specified con

Level in dB. It can be specified as a scalar or a vector. If it is specified as a vector, the level is swept. The linRg function can be called to generate a linear range. For example, linRg ( -30 30 5 ) is the same as list (-30 -25 -20 -15 -10 -5 0 5 10 15 20 25 30) and the  $g\_level$  argument can be specified as either. In that case, a power gain circle is calculated

at each one of the 13 levels.

g\_frequency The frequency. It can be specified as a scalar or a linear range.

If it is specified as a linear range, the frequency is swept. The linear range is specified as a list with three values: the start of the range, the end of the range, and the increment. For

example, list (100M 1G 100M) specifies a linear range with

the following values:

{ 100M, 200M, 300M, 400M, 500M, 600M, 700M, 800M, 900M,

1G }

In that case, a power gain circle is calculated at each one of the

10 frequencies.

#### Value Returned

o waveform bject representing the power gain circles.

nil Returns nil and an error message otherwise.

### Predefined and Waveform (Calculator) Functions

## groupDelay

```
groupDelay(
    o_waveform
)
=> o_waveform / nil
```

### **Description**

Computes the group delay of a waveform.

This command returns the derivative of the phase of  $o\_waveform$  / 2pi. Group delay is defined as the derivative of the phase with respect to frequency. Group delay is expressed in seconds.

It is calculated using the vp function as shown below:

Group Delay = 
$$\frac{d\phi}{d\omega} = \frac{d}{df} \left[ \frac{phase(/\text{net}X)}{360} \right]$$

### **Arguments**

o_waveform	Waveform object representing simulation results that can be
	displayed as a series of points on a grid. (A waveform object
	identifier looks like this: srrWave: XXXXX.)

#### Value Returned

o_waveform	Returns a waveform representing the group delay of the specified waveform.
nil	Returns nil and an error message otherwise.

## **Example**

```
plot( groupDelay( v( "/net3" ) ) )
```

Plots the waveform representing the group delay of the voltage of "/net3".

### Predefined and Waveform (Calculator) Functions

## gt

```
gt(
    o_s11
    o_s12
    o_s21
    o_s22
    [ ?gs n_gs ]
    [ ?gl n_gl ]
    )
    => o waveform / nil
```

## **Description**

Returns the transducer gain in terms of the supplied parameters and the optional source reflection coefficient (Gs) and the input reflection coefficient (Gl).

## **Arguments**

o_ <i>s</i> 11	Waveform object representing s11.
o_s12	Waveform object representing s12.
o_s21	Waveform object representing s21.
o_s22	Waveform object representing s22.
n_gs	Source reflection coefficient. Default value: 0
n_gl	Input reflection coefficient. Default value: 0

#### Value Returned

o_waveform	Waveform object representing the transducer gain.
nil	Returns nil and displays a message if there is an error.

```
s11 = sp(1 1)
s12 = sp(1 2)
s21 = sp(2 1)
s22 = sp(2 2)
plot(gt(s11 s12 s21 s22))
```

# Predefined and Waveform (Calculator) Functions

Note: gl is an imaginary number which should be input in the following format:

gt( s11 s12 s21 s22 ?gl complex(<realPart> <imagPart>))

## Predefined and Waveform (Calculator) Functions

## harmonic

```
harmonic(
    o_waveform
    h_index
)
=> o_waveform / g_value / nil
```

# **Description**

Returns the waveform for a given harmonic index.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
h_index	The index number that designates the harmonic information to be returned. For the 'pss, 'pac, and 'pxf analyses, the index is an integer number. For the 'pdisto analysis, the index is a list of integers that correspond with the frequency names listed in the funds analysis parameter in the netlist. If more than one $h\_index$ is desired at one time, a list can be specified.

### Value Returned

o_waveform	Returns a waveform (when a single $h\_index$ is specified) or family of waveforms (when more than one $h\_index$ is specified) if the input argument is a family of waveforms.
g_value	Returns the harmonic value if the input is a single waveform with the X values being harmonics
nil	Returns nil and displays a message if there is an error.

## **Example**

### For each of the following commands:

```
harmonic(v("/net49" ?result "pss-fd.pss") 1)
harmonic(v("/Pif" ?result "pdisto-fi.pdisto") list(1 -1))
```

Each result is a complex number.

### Predefined and Waveform (Calculator) Functions

### For each of the following commands:

```
harmonic(v("/net54" ?result "pac-pac") 1)
harmonic(v("/net51" ?result "sweeppss_pss_fd-sweep") list(8))
harmonic(v("/Pif" ?result "sweeppss_pac-sweep") -8)
harmonic(v("/net36" ?result "sweeppdisto pdisto fi-sweep") '(1 -1))
```

#### Each result is a waveform.

### For each of the following commands:

### Each result is a family of waveforms.

## Neither of the following commands should be entered:

```
harmonic(v("/net49" ?result "pss-fd.pss") list(0 1))
harmonic(v("/Pif" ?result "pdisto-fi.pdisto") '((1 -1) (-1 2)))
```

Each resulting waveform is not in a useful format.

## Predefined and Waveform (Calculator) Functions

# harmonicFreqList

```
harmonicFreqList(
    [ ?resultsDir t_resultsDir ]
    [ ?result S_resultName ]
)
    => n list / nil
```

# **Description**

Returns a list of lists, with each sublist containing a harmonic index and the minimum and maximum frequency values that the particular harmonic ranges between.

If both of these frequency values are the same, just one frequency value is returned.

## **Arguments**

t_resultsDir	Directory containing the PSF files (results). If you supply this argument, you must also supply the resultName argument.
S_resultName	Results from an analysis.

### Value Returned

n_list	Returns a list of lists. For the 'pss, 'pac, and 'pxf analyses, the first element of each sublist is an integer number. For the 'pdisto analysis, the first element of each sublist is a list of integers that correspond with the frequency names listed in the funds analysis parameter in the netlist. For all sublists, the remaining entries are the minimum and maximum frequency values that the particular harmonic ranges between. If both of these frequency values are the same, just one frequency value is returned.
nil	Returns nil if no harmonics are found in the data.

### **Example**

### For each of the following commands:

```
harmonicFreqList( ?result "pss-fd.pss" )
harmonicFreqList( ?result "pac-pac" )
harmonicFreqList( ?result "sweeppss pss fd-sweep" )
```

### Predefined and Waveform (Calculator) Functions

```
harmonicFreqList( ?result "sweeppss pac-sweep" )
```

Each result is a list of integers.

### For each of the following commands:

```
harmonicFreqList( ?result "pdisto-fi.pdisto" )
harmonicFreqList( ?result "sweeppdisto pdisto fi-sweep" )
```

Each result is a list of lists, with each sublist containing a combination of integer numbers that correspond with the frequency names listed in the funds analysis parameter in the netlist. These names can also be extracted from the PSF data by using the resultParam function to find the 'largefundname and 'moderatefundnames values. For example:

```
strcat(resultParam( 'largefundname ?result "pdisto-fi.pdisto" ) " "
resultParam( 'moderatefundnames ?result "pdisto-fi.pdisto" ))
```

Returns a string representing the order of the frequency names.

### Predefined and Waveform (Calculator) Functions

## harmonicList

```
harmonicList(
    [ ?resultsDir t_resultsDir ]
    [ ?result S_resultName ]
    )
    => n list
```

## **Description**

Returns the list of harmonic indices available in the resultName or current result data.

### **Arguments**

t_resultsDir	Directory containing the PSF files (results). If you supply this argument, you must also supply the resultName argument.
$S\_resultName$	Results from an analysis.

### Value Returned

n_list	Returns a list of harmonic indices. For the 'pss, 'pac, and 'pxf analyses, the index is an integer number. For the 'pdisto analysis, the index is a list of integers that correspond with the frequency names listed in the 'funds analysis parameter in the netlist.
nil	Returns nil if no harmonics are found in the data.

### **Example**

### For each of the following commands:

```
harmonicList( ?result "pss-fd.pss" )
harmonicList( ?result "pac-pac" )
harmonicList( ?result "sweeppss_pss_fd-sweep" )
harmonicList( ?result "sweeppss pac-sweep" )
```

Each result is a list of integers.

## For each of the following commands:

```
harmonicList( ?result "pdisto-fi.pdisto" )
harmonicList( ?result "sweeppdisto pdisto fi-sweep" )
```

## Predefined and Waveform (Calculator) Functions

Each result is a list of lists, with each sublist containing a combination of integer numbers that correspond with the frequency names listed in the 'funds analysis parameter in the netlist. These names can also be extracted from the PSF data by using the 'resultParam function to find the 'largefundname and 'moderatefundnames values. For example:

```
strcat(resultParam( 'largefundname ?result "pdisto-fi.pdisto" ) " "
resultParam( 'moderatefundnames ?result "pdisto-fi.pdisto" ))
```

Returns a string representing the order of the frequency names.

## Predefined and Waveform (Calculator) Functions

### histo

```
histo(
    o_waveform
    x_bins
    n_min
    n_max
)
=> o histoWaveform / nil
```

### **Description**

Returns a waveform that represents the statistical distribution of input data in the form of a histogram. The height of the bars (or bins) in the histogram represents the frequency of the occurrence of values within a specific period. Using the histo function, the range for capturing these frequencies can be specified through the  $n\_min$  and  $n\_max$  values.

# **Arguments**

o_waveform	Input waveform.
$x\_bins$	Number of bins to represent the input data.
n_min	The first value on the horizontal axis of the histogram. By default, it assumes the minimum value of the input waveform.
n_max	The last value on the horizontal axis of the histogram. By default, it assumes the maximum value of the input waveform.

### Value Returned

o_histoWaveform	Returns a waveform representing the statistical distribution of the input waveform o_waveform.
nil	Returns nil in case of an error.

## **Example**

```
histo( VT("/vin") 3 1.5 3.5)
=> out_wave
plot( out wave )
```

Plots the output waveform  $out\_wave$  as a histogram, which represents the statistical distribution of the input waveform VT("/vin").

# Predefined and Waveform (Calculator) Functions

# histogram2D

```
histogram2D(
    o_waveform
    x_nbins
    s_type
    g_setAnnotation
    g_setDensityEstimator)
=> o waveform / nil
```

# **Description**

Returns a waveform that represents the statistical distribution of input data in the form of a histogram. The height of the bars (or bins) in the histogram represents the frequency of the occurrence of values within a specific period.

## Predefined and Waveform (Calculator) Functions

### **Arguments**

o waveform Input waveform.

x nbins Number of bins (bars) to be plotted in the resulting histogram

plot.

Valid values: 1 to 50.

Default value:10.

s type Type of histogram to be plotted.

Valid values: Standard, Cumulative line, and

Cumulative box.

Default value: Standard.

g setAnnotation Boolean specifying whether to display the standard deviation

lines in the resulting histogram plot.

Valid values: t or nil

Default value: nil

g\_setDensityEstim

ator

Boolean specifying whether the resulting histogram plot display

a curve that estimates the distribution concentration.

Valid values: t or nil

Default value: nil

#### Value Returned

o waveform Returns a waveform representing the statistical distribution of

the input waveform o waveform.

nil Returns nil in case of an error.

### Example

histogram2D(i("/V2/PLUS" ?result "tran") 10 "standard" t t )

Plots the output waveform out\_wave as a histogram, which represents the statistical distribution of the input waveform /V2/PLUS.

## Predefined and Waveform (Calculator) Functions

# iinteg

```
iinteg(
    o_waveform
)
=> o waveform / nil
```

# **Description**

Computes the indefinite integral of a waveform with respect to the X-axis variable.

# **Arguments**

o_waveform	Waveform object representing simulation results that can be
	displayed as a series of points on a grid. (A waveform object
	identifier looks like this: srrWave: XXXXX.)

## **Value Returned**

o_waveform	Returns a waveform representing the indefinite integral of the input waveform.
nil	Returns nil and an error message otherwise.

# Example

```
plot( iinteg( v( "/net8" )))
```

Computes the indefinite integral of the waveform representing the voltage of "/net8".

### Predefined and Waveform (Calculator) Functions

# imag

```
imag(
     { o_waveform | n_input }
)
=> o waveformImag / n numberImag / nil
```

### **Description**

Returns the imaginary part of a waveform representing a complex number or returns the imaginary part of a complex number.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_input	Complex number.

### Value Returned

o_waveformImag	Returns a waveform when the input argument is a waveform.
n_numberImag	Returns a number when the input argument is a number.
nil	Returns nil and an error message otherwise.

### **Example**

```
imag( v( "/net8" ) )
```

Returns a waveform representing the imaginary part of the voltage of "/net8". You also can use the vim alias to perform the same command, as in

```
vim( "net8" ).
x=complex( -1 -2 ) => complex(-1, -2)
imag( x ) => -2.0
```

Creates a variable  ${\bf x}$  representing a complex number, and returns the real portion of that complex number.

## Predefined and Waveform (Calculator) Functions

# inl

```
inl(
    o_dacSignal
    o_sample | o_pointList | n_interval
    [ ?mode t_mode ]
    [ ?threshold n_threshold ]
    [ ?crossType t_crossType ]
    [ ?delay f_delay ]
    [ ?units x_units ]
    [ ?nbsamples n_nbsamples ]
    )
    => n_inl / nil
```

# **Description**

Computes the integral non-linearity of a transient simple or parametric waveform.

# Predefined and Waveform (Calculator) Functions

# **Arguments**

o_dacSignal	Waveform for which the integral non-linearity is to be calculated.
o_sample	Waveform used to obtain the points for sampling the dacSignal. These are the points at which the waveform crosses the threshold while either rising or falling (defined by the crossType argument) with the delay added to them.
n_pointList	List of domain values at which the sample points are obtained from the dacSignal.
n_interval	The sampling interval.
?mode t_mode	The mode for calculating the threshold.
	Valid values: auto and user.
	Default value: auto.
	If set to user, an n_threshold value needs to be provided.
	If set to auto, n_threshold is calculated internally.
?threshold n_threshold	The threshold value against which the integral non-linearity is to be calculated. It needs to be specified only when the mode is selected as user.
?crossType t_crossType	The points at which the curves of the waveform intersect with the threshold. While intersecting, the curve may be either rising or falling.
	Valid values: rising and falling, respectively.
	Default crossType is rising.
?delay f_delay	The delay time after which the sampling begins.
	Valid values: Any valid time value.
	Default value: 0.
?unit x_units	Unit for expressing the output waveform.
	Valid values: abs (absolute) and 1sb (multiples of least significant bit).
	Default value: abs.

?nbsamples  $n_nbsamples$ 

## Predefined and Waveform (Calculator) Functions

Number of samples used for calculating the non-linearity. If not specified, the samples are taken against the entire data window.

**Note:** For each of the three ways in which the sample points can be specified, only a few of the other optional arguments are meaningful, as indicated below:

- For o\_sample, the arguments  $t_mode, n_threshold, t_crossType, f_delay,$  and  $x_units$  are meaningful.
- For  $n_{pointList}$ , the arguments  $x_{units}$  are meaningful.
- For n interval, the arguments x units, and n nbsamples are meaningful.

#### Value Returned

$n\_inl$	Returns the integral non-linearity waveform.
nil	Returns nil and an error message otherwise.

## **Example**

```
inl( wave1 wave2 ?crossType "rising" ?delay 0.4 )
=> srrWave:175051544
```

Returns the integral non-linearity for wave1 by taking the points at which wave2 crosses the internally calculated threshold while rising as the sample points and adding a delay of 0.4 to them.

### Predefined and Waveform (Calculator) Functions

# integ

```
integ(
    o_waveform
    [ n_intial_limit , n_final_limit ]
)
    => o waveform / n value / nil
```

# **Description**

Computes the definite integral of the waveform with respect to a range specified on the X-axis of the waveform. The result is the value of the area under the curve over the range specified on the X-axis.

You should specify either both the limits or neither. In case you do specify the limits, they become the end points of the range on the X-axis for definite integration. If you do not specify the limits, then the range for definite integration is the entire range of the sweep on the X-axis.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
initial_limit_n	Initial limit for definite integration.
final_limit_n	Final limit for definite integration.

### Value Returned

o_waveform	Returns a waveform representing the definite integral for a family of waveforms if the input argument is a family of waveforms.
n_value	Returns a numerical value representing the definite integral of the input waveform if the input argument is a single waveform.
nil	Returns nil and an error message otherwise.

# **Example**

```
integ( v( "/out" ) )
```

# Predefined and Waveform (Calculator) Functions

Returns the definite integral of the waveform representing the voltage of "/out" over its entire range.

```
integ( VT( "/out" ),12.5n,18n)
```

Returns the definite integral of the waveform representing the voltage of "/out" within a specified range.

# Predefined and Waveform (Calculator) Functions

### intersect

```
intersect(
    o_waveform1
    o_waveform2
)
    => o wave / nil
```

# **Description**

Returns a waveform containing the points of intersection for two waveforms passed as arguments.

# **Arguments**

o_waveform1	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
o_waveform2	Additional waveform object.

### Value Returned

o_wave	Returns a waveform containing the points of intersection for the two waveforms passed as arguments.
nil	Returns nil if the two waveforms are disjoint or overlap each other, and an error message, if the arguments to the function are not correct.

# **Example**

```
intersect( VT("/inp1") VT("/inp2") )
```

### Predefined and Waveform (Calculator) Functions

# ipn

```
ipn(
    o_spurious
    o_reference
    [ f_ordspur f_ordref f_epspur f_epref g_psweep s_measure ]
    )
    => o_waveform / f_number / nil
```

# Description

Performs an intermodulation *n*th-order intercept measurement.

The data for this measurement can be either a single input power value or a parametric input power sweep.

From each of the spurious and reference power waveforms (or points), the <code>ipn</code> function extrapolates a line of constant slope (dB/dB) according to the specified order and input power level. These lines represent constant small-signal power gain (ideal gain). The <code>ipn</code> function calculates the intersection of these two lines and returns the value of either the X coordinate (input referred) or Y coordinate.

# Predefined and Waveform (Calculator) Functions

# **Arguments**

o_spurious	Waveform or number representing the spurious output power (in dBm).
o_reference	Waveform or number representing the reference output power (in dBm).
f_ordspur	Order or slope of the spurious constant-slope power line.
	Default value: 3
f_ordref	Order or slope of the reference constant-slope power line.
	Default value: 1
f_epspur	Value (in dBm) used to indicate the point where the spurious constant-slope power line begins. If $g\_psweep$ is t, this value is the input power value of the point on the $o\_spurious$ waveform, otherwise this value is paired with the $o\_spurious$ value to define the point. This point should be in the linear region of operation. (If $g\_psweep$ is t, $f\_spspur$ defaults to the X coordinate of the first point of the $o\_spurious$ wave; if $s\_measure$ is 'input, a number must be specified.)
f_epref	Value (in dBm) used to indicate the point where the reference constant-slope power line begins. If $g_psweep$ is t, this value is the input power value of the point on the $o_reference$ waveform, otherwise this value is paired with the $o_reference$ value to define the point. This point should be in the linear region of operation. (If $g_psweep$ is t, $f_epref$ defaults to the X coordinate of the first point of the $o_reference$ wave; if $s_measure$ is 'input, a number must be specified.)
g_psweep	Boolean indicating that the input power to the circuit was a parametric sweep. The power sweep must both be in dBm and be performed at the lowest parametric level.
	Default value: t
s_measure	Name indicating if measurement is to be input referred ('input) or output referred ('output).
	Default value: 'input

## Predefined and Waveform (Calculator) Functions

#### Value Returned

o_waveform	Depending on setting of $g_psweep$ and the dimension of the input waveforms, returns either a waveform or a family of waveforms.
f_number	If $o\_spurious$ and $o\_reference$ are numbers or they are waveforms when $g\_psweep$ is t, returns a number.
nil	Returns nil and an error message otherwise.

### **Example**

```
spurWave = db20(harmonic(wave signalHarmonic))
refWave = db20(harmonic(wave referenceHarmonic))
xloc = ipn( spurWave refWave 3.0 1.0 -25 -25 )
yloc = ipn( spurWave refWave 3.0 1.0 -25 -25 t "Output")
```

Computes the IP3 point for the given wave.

### Each of the following examples returns an ip3 measurement.

```
ipn(dB20(harmonic(v("/Pif" ?result "pss fd") 9))
    dB20(harmonic(v("/Pif" ?result "pss fd") 8)))
ipn(dbm(harmonic(spectralPower(v("/Pif" ?result "pss fd")/50.0
    v("/Pif" ?result "pss fd")) 9))
    dbm(harmonic(spectralPower(v("/Pif" ?result "pss fd")/50.0
    v("/Pif" ?result "pss fd")) 8)))
ipn(dbm(harmonic(spectralPower(v("/Pif" ?result "pss fd")
    /resultParam("rif:r" ?result "pss td")
    v("/Pif" ?result "pss fd")) 9))
    \label{thm:dbm:equation:dbm:equation} $$\operatorname{dbm}(\operatorname{harmonic}(\operatorname{spectral}\overline{P}\operatorname{ower}(\operatorname{v}("/\operatorname{Pif"}\ \operatorname{?result}\ "\operatorname{pss}\ \operatorname{fd"}))$
    /resultParam("rif:r" ?result "pss td")
    v("/Pif" ?result "pss fd")) 8)))
ipn(dbm(harmonic(spectralPower(i("/rif/PLUS" ?result "pss fd")
    v("/Pif" ?result "pss fd")) 9))
    dbm(harmonic(spectralPower(i("/rif/PLUS" ?result "pss fd")
    v("/Pif" ?result "pss fd")) 8))
    3. 1. -25 -25 t "Output")
ipn(dbm(harmonic(spectralPower(v("/Pif" ?result "pac")
    /resultParam("rif:r" ?result "pss td")
    v("/Pif" ?result "pac")) -21))
    dbm(harmonic(spectralPower(v("/Pif" ?result "pac")
    /resultParam("rif:r" ?result "pss td")
    v("/Pif" ?result "pac")) -25)))
```

### Predefined and Waveform (Calculator) Functions

# ipnVRI

```
ipnVRI(
    o_vport
    x_harmspur
    x_harmref
    [?iport o_iport]
    [?rport f_rport]
    [?ordspur f_ordspur] | [?epoint f_epoint]
    [?psweep g_psweep]
    [?epref f_epref]
    [?ordref f_ordref]
    [?measure s_measure]
)
    => o waveform / f number / nil
```

## **Description**

Performs an intermodulation *n*th-order intercept point measurement.

Use this function to simplify the declaration of an ipn measurement. This function extracts the spurious and reference harmonics from the input waveform(s), and uses  ${\tt dBm} ({\tt spectralPower} ((i \ {\tt or} \ {\tt v/r}), {\tt v})) \ to \ {\tt calculate} \ the \ respective \ powers. The function passes these power curves or numbers and the remaining arguments to the ipn function to complete the measurement.$ 

From each of the spurious and reference power waveforms (or points), the <code>ipn</code> function extrapolates a line of constant slope (dB/dB) according to the specified order and input power level. These lines represent constant small-signal power gain (ideal gain). The <code>ipn</code> function calculates the intersection of these two lines and returns the value of either the X coordinate (input referred) or the Y coordinate.

### **Arguments**

o_vport	Voltage across the output port. This argument must be a family of spectrum waveforms (1 point per harmonic), with the option of containing a parametric input power sweep (in dBm).
x_harmspur	Harmonic number of the spurious voltage contained in <code>o_vport</code> . When <code>o_iport</code> is specified, also applies to a current waveform contained in <code>o_iport</code> .
x_harmref	Harmonic index of the reference voltage contained in o_vport. When o_iport is specified, also applies to a current waveform contained in o_iport.

### Predefined and Waveform (Calculator) Functions

?iport o\_iport Current into the output port. This argument must be a family of

spectrum waveforms (1 point per harmonic), with the option of containing a parametric input power sweep (in dBm). When specified, power is calculated using voltage and current.

?rport f rport Resistance into the output port. When specified and o iport

is nil, the output power is calculated using voltage and

resistance.

Default value: 50

?ordspur Order or slope of the spurious constant-slope power line.

*f\_ordspur* Default value: 3

?epoint f\_epoint Value (in dBm) used to indicate the point where the spurious

constant-slope power line begins. If  $g_psweep$  is t, this value is the input power value of the point on the  $o_spurious$  waveform, otherwise this value is paired with the  $o_spurious$  value to define the point. This point should be in the linear

region of operation.

Default value: If  $g_psweep$  is t, the lowest input power value;

if  $s_{measure}$  is 'input, a number must be specified.

?psweep g psweep Boolean indicating that the input power to the circuit was a

parametric sweep. The power sweep must be in dBm and must

be performed at the lowest parametric level.

Default value: t

?epref f\_epref Value (in dBm) used to indicate the point where the reference

constant-slope power line begins. If  $g_psweep$  is t, this value is the input power value of the point on the  $o_reference$ 

waveform, otherwise this value is paired with the

o reference value to define the point. This point should be

in the linear region of operation.

Default value: If  $f_{epoint}$  is not nil,  $f_{epoint}$ ; else if  $g_{epoint}$  is t, the X coordinate of the first point of the

o reference wave; else if s measure is 'input, a

number must be specified.

?ordref f ordref Order or slope of the reference constant-slope power line.

Default value: 1

## Predefined and Waveform (Calculator) Functions

?measure Symbol indicating if measurement is to be input referred s measure

('input) or output referred ('output).

Default value: 'input

### Value Returned

Depending on the setting of q psweep and the dimension of o waveform

input waveform(s), the ipnVRI function returns either a

waveform or a family of waveforms.

Depending on the setting of g psweep and the dimension of f number

input waveform(s), the ipnVRI function returns a number.

nil Returns nil and an error message otherwise.

### **Example**

### Each of following examples returns an ip3 measurement:

```
ipnVRI(v("/Pif" ?result "pss fd") 9 8)
ipnVRI(v("/Pif" ?result "pss_fd") 9 8
    ?iport i("/rif/PLUS" ?result "pss_fd") ?epoint -25
   ?measure "Output")
ipnVRI(v("/Pif" ?result "pac") -21 -25
    ?rport resultParam("rif:r" ?result "pss td"))
```

### Predefined and Waveform (Calculator) Functions

# **ipnVRICurves**

```
ipnVRICurves(
    o_vport
    x_harmspur
    x_harmref
    [?iport o_iport]
    [?rport f_rport]
    [?ordspur f_ordspur]
    [?epoint f_epoint]
    [?psweep g_psweep]
    [?epref f_epref]
    [?ordref f_ordref]
)
    => o waveform / nil
```

### Description

Constructs the waveforms associated with an ipn measurement.

Use this function to simplify the creation of waves associated with an ipn measurement. This function extracts the spurious and reference harmonics from the input waveform(s), and uses dBm(spectralPower((i or v/r), v)) to calculate the respective powers.

From each of the spurious and reference power waveforms (or points), the <code>ipnVRICurves</code> function extrapolates a line of constant slope (dB/dB) according to the specified order and input power level. These lines represent constant small-signal power gain (ideal gain). The function returns these lines and power waveforms (when present) as a family of waveforms.

This function only creates waveforms and does not perform an ipn measurement or include labels with the waveforms. Use the ipn or ipnVRI function for making measurements.

# Predefined and Waveform (Calculator) Functions

# **Arguments**

o_vport	Voltage across the output port. This argument must be a family of spectrum waveforms (1 point per harmonic), with the option of containing a parametric input power sweep (in dBm).
x_harmspur	Harmonic index of the spurious voltage contained in $o\_vport$ . When $o\_iport$ is specified, also applies to a current waveform contained in $o\_iport$ .
x_harmref	Harmonic index of the reference voltage contained in <code>o_vport</code> . When <code>o_iport</code> is specified, also applies to a current waveform contained in <code>o_iport</code> .
?iport o_iport	Current into the output port. This argument must be a family of spectrum waveforms (1 point per harmonic), with the option of containing a parametric input power sweep (in dBm). When specified, power is calculated using voltage and current.
?rport f_rport	Resistance into the output port. When specified and $o\_iport$ is nil, the output power is calculated using voltage and resistance.
	Default value: 50
?ordspur	Order or slope of the spurious constant-slope power line.
f_ordspur	Default value: 3
?epoint f_epoint	Value (in dBm) used to indicate the point where the spurious constant-slope power line begins. If <code>g_psweep</code> is t, this value is the input power value of the point on the <code>o_spurious</code> waveform, otherwise this value is paired with the <code>o_spurious</code> value to define the point. This point should be in the linear region of operation.
	Default value: If $g_psweep$ is t, the X coordinate of the first point of the $o_spurious$ wave; otherwise a number must be specified.
?psweep <i>g_psweep</i>	Boolean indicating that the input power to the circuit was a parametric sweep. The power sweep must be in dBm and must be performed at the lowest parametric level.
	Default value: t

### Predefined and Waveform (Calculator) Functions

?epref  $f_{epref}$  Value (in dBm) used to indicate the point where the reference

constant-slope power line begins. If  $g_psweep$  is t, this value is the input power value of the point on the o reference

waveform, otherwise this value is paired with the

o\_reference value to define the point. This point should be

in the linear region of operation.

Default value: If  $f_{epoint}$  is not nil,  $f_{epoint}$ ; else if  $g_{psweep}$  is t, the X coordinate of the first point of the oreference wave; else a number must be specified.

?ordref f ordref Order or slope of the reference constant-slope power line.

Default value: 1

#### Value Returned

o waveform A family of waveforms that contains the spurious and reference

tangent lines, and when g psweep is t, contains the spurious

and reference waveforms.

nil Returns nil and an error message otherwise.

#### **Example**

Each of following examples returns curves related to an ip3 measurement:

# Predefined and Waveform (Calculator) Functions

## kf

```
kf(
     o_s11
     o_s12
     o_s21
     o_s22
)
     => o_waveform / nil
```

# **Description**

Returns the stability factor in terms of the supplied parameters.

# **Arguments**

o_s11	Waveform object representing s11.
0_s12	Waveform object representing s12.
0_s21	Waveform object representing s21.
o_s22	Waveform object representing s22.

### Value Returned

o_waveform	Waveform object representing the stability factor.
nil	Returns nil if there is an error.

# **Example**

```
s11 = sp(1 1)

s12 = sp(1 2)

s21 = sp(2 1)

s22 = sp(2 2)

plot(kf(s11 s12 s21 s22))
```

## Predefined and Waveform (Calculator) Functions

### In

## **Description**

Gets the base-e (natural) logarithm of a waveform or number.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_number	Number.

### Value Returned

o_waveform	Returns a waveform object representing the base-e (natural) logarithm of the input waveform if the input argument is a waveform object, or returns a family of waveforms if the input argument is a family of waveforms
f_number	Returns a number if the input argument is a number.
nil	Returns nil and an error message otherwise.

### **Example**

```
ln( v( "/net9" ) )
```

Gets a waveform that is calculated as the natural logarithm of the input waveform.

```
ln(ymax(v("/net9")))
```

Gets a waveform that is calculated as the natural logarithm of the following: ymax (v ( " / net 9 " ) ) .

```
ln(100) => 4.60517
```

Gets the natural logarithm of 100.

### Predefined and Waveform (Calculator) Functions

# log10

## **Description**

Gets the base-10 logarithm of a waveform or a number.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n number	Number.

### Value Returned

o_waveform	Returns a waveform object if the input argument is a waveform object or returns a family of waveforms if the input argument is a family of waveforms.
n_number	Returns a number that is calculated as the base-10 logarithm of the input number.
nil	Returns nil and an error message otherwise.

# **Example**

```
log10(v("/net9"))
```

Gets a waveform that is calculated as the base-10 logarithm of the input waveform.

```
log10( ymax( v( "/net9" ) ) )
```

Gets a waveform representing the base-10 logarithm of ymax (v("/net9")).

```
log10 ( 100 ) => 2.0
```

Gets the base-10 logarithm of 100, or 2.

# Predefined and Waveform (Calculator) Functions

# Isb

```
lsb(
    o_s11
    o_s12
    o_s21
    o_s22
    g_frequency
)
    => o_waveform / nil
```

# **Description**

Computes the load stability circles.

## Predefined and Waveform (Calculator) Functions

### **Arguments**

o_s11	Waveform object representing s11.
o_s12	Waveform object representing s12.
o_s21	Waveform object representing s21.
o_s22	Waveform object representing s22.

g\_frequency Frequency. It can be specified as a scalar or a linear range. If it

is specified as a linear range, the frequency is swept. The linear range is specified as a list with three values: the start of the range, the end of the range, and the increment. For example, list (100M 1G 100M) specifies a linear range with the

following values:

```
{ 100M, 200M, 300M, 400M, 500M, 600M, 700M, 800M, 900M, 1G }
```

In that case, a load stability circle is calculated at each one of

the 10 frequencies

### Value Returned

o\_waveform Waveform object representing the load stability circles.

nil Returns nil and an error message otherwise.

### **Example**

plot(lsb(s11 s12 s21 s22 list(800M 1G 100M)))

# Predefined and Waveform (Calculator) Functions

## **Ishift**

```
lshift(
    o_waveform
    n_delta
)
=> o waveform / nil
```

# **Description**

Shifts the waveform to the left by the delta value.

This command is the inverse of the <u>rshift</u> command.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_delta	Value by which the waveform is to be shifted.

### Value Returned

o_waveform	Returns a waveform object representing the input waveform shifted to the left. Returns a family of waveforms if the input argument is a family of waveforms.
nil	Returns nil and an error message otherwise.

### Example

```
plot( lshift( v( "/net8" ) 30u ) )
```

Shifts the waveform representing the voltage of "/net8" to the left by 30u and plots the resulting waveform.

## Predefined and Waveform (Calculator) Functions

## mag

```
mag(
     { o_waveform | n_number }
)
=> o waveform / n number / nil
```

### **Description**

Gets the magnitude of a waveform or number.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_number	Number.

### Value Returned

o_waveform	Returns a waveform object if the input argument is a waveform object or returns a family of waveforms if the input argument is a family of waveforms.
n_number	Returns a number if the input argument is a number.
nil	Returns nil and an error message otherwise.

### **Example**

```
mag( v( "5" ) )
```

Gets the magnitude of the waveform representing the voltage at net 5. You can also use the vm alias to perform the same command, as in vm ( "5" ).

```
mag( i( "VFB" ) )
```

Gets the magnitude of the waveform representing current through the VFB component. You can also use the im alias to perform the same command, as in im ( "VFB" ).

```
mag(-10) => 10
```

Returns the magnitude of -10.

# Predefined and Waveform (Calculator) Functions

# nc

```
nc(
    o_Fmin
    o_Gmin
    o_rn
    g_level
    g_frequency
)
    => o_waveform / nil
```

# **Description**

Computes the noise circles.

### Predefined and Waveform (Calculator) Functions

## **Arguments**

Fmin
 Waveform object representing the minimum noise factor.

O Gmin Waveform object representing the optimum noise reflection.

o\_rn Waveform object representing the normalized equivalent noise

resistance.

g level Level in dB. It can be specified as a scalar or a vector. The level

is swept, if it is specified as a vector. The linRg function can be called to generate a linear range. For example, linRg(-30 30 5) is the same as list(-30 -25 -20 -15 -10 -5 0 5 10 15 20 25 30) and the  $g_level$  argument can be specified as either of the above. In that case, a noise circle is

calculated at each one of the 13 levels.

g\_frequency Frequency. It can be specified as a scalar or a linear range. The

frequency is swept if it is specified as a linear range. The linear range is specified as a list with three values: the start of the range, the end of the range, and the increment. For example, list (100M 1G 100M) specifies a linear range with the

following values:

{100M, 200M, 300M, 400M, 500M, 600M, 700M, 800M, 900M,

1G }

In that case, a noise circle is calculated at each one of the 10

frequencies.

### Value Returned

o waveform Waveform object representing the noise circles.

nil Returns nil and an error message otherwise.

### **Example**

```
Gopt = getData("Gopt")
Bopt = getData("Bopt")
Zref = zref(1 ?result "sp")
Gmin = gmin(Gopt Bopt Zref)
Fmin = getData("Fmin")
rn = getData("NNR")
NC = nc(Fmin Gmin rn 10 list(100M 1G 100M))
```

## Predefined and Waveform (Calculator) Functions

displayMode("smith")
smithType("impedance")
plot(NC)

## Predefined and Waveform (Calculator) Functions

## normalQQ

```
normalQQ(
    o_waveform
)
=> o waveform / nil
```

## **Description**

Returns a quantile-quantile plot of the sample quantiles versus theoretical quantiles from a normal distribution. If the distribution is normal, the plot is close to a linear waveform.

## **Argument**

o_waveform	Waveform object representing simulation results that can be
	displayed as a series of points on a grid. (A waveform object
	identifier looks like this: srrWave: XXXXX.)

#### **Values Returned**

o_waveform	Returns the waveform representing the normal quantile plot.
nil	Returns nil and an error message otherwise.

## **Example**

```
normalQQ(v("net10" ?result "tran"))
```

Returns the quantile plot for the v("net10" ?result "tran") signal.

# Predefined and Waveform (Calculator) Functions

### overshoot

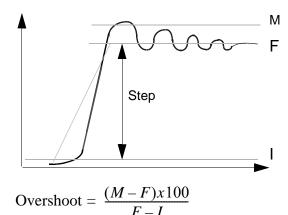
```
overshoot(
    o waveform
    n_initVal
    g_initType
    n finalVal
    g finalType
     [ g multiple [ s_Xname ] ]
     [ g_histoDisplay ]
     [ x noOfHistoBins ]
    => o waveform / n value /nil
```

## **Description**

Computes the percentage by which an expression overshoots a step going from the initial value to the final value you enter.

This command returns the overshoot of o waveform as a percentage of the difference between the initial value and the final value.

In the equation below, M represents Maximum Value of the peak wave, F represents Final Value of the settled wave, and I represents Initial Value of the wave.



## Predefined and Waveform (Calculator) Functions

## Arguments

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_initVal	Initial X value at which to start the computation.
g_initType	Specifies how initVal functions.
	Valid values: a non-nil value specifies that the initial value is taken to be the value of the waveform, interpolated at <pre>initVal</pre> , and the waveform is clipped from below, as follows:
	<pre>o_waveform = clip( o_waveform initVal nil )</pre>
	nil specifies that $initVal$ is defined by the X value entered. (The command gets the Y value for the specified X value and uses that value for $initVal$ .)
n_finalVal	Final value at which to end the computation.
g_finalType	Specifies how final Val functions.
	Valid values: a non-nil value specifies that the final value is taken to be the value of the waveform, interpolated at final Val, and the waveform is clipped from above, as follows:
	<pre>o_waveform = clip( o_waveform nil finalVal)</pre>
	nil specifies that finalVal is defined by the X value entered. (The command gets the Y value for the specified X value and uses that value for finalVal.)
g_multiple	An optional boolean argument that takes the value $\tt nil$ by default. If set to $\tt t$ , the function returns multiple occurrences of the overshoot event.
s_xName	An optional argument that is used only when $g_{multiple}$ is set to t. It takes the value time by default. It controls the contents of the x vector of the waveform object returned by the function.
	Valid values: 'time, 'cycle

## Predefined and Waveform (Calculator) Functions

g histoDisplay When set to t, returns a waveform that represents the statistical

distribution of the riseTime data in the form of a histogram. The

height of the bars (bins) in the histogram represents the frequency of the occurrence of values within the range of

riseTime data.

Valid values: t nil

Default value: nil

x noOfHistoBins Denotes the number of bins represented in the histogram

representation.

Valid values: Any positive integer

Default value: nil

**Note:**  $g_histoDisplay$  and  $x_noOfHistoBins$  are added for backward compatibility only. It will be deprecated in future releases. Use the histo function for plotting the histogram of the resulting function.

#### Value Returned

o_waveform	Returns a waveform (or family of waveforms) representing the amount of overshoot in comparison to the whole signal if the input argument is a family of waveforms or if $g_{multiple}$ is set to t.

 $n\_value$  Returns a value for the amount of overshoot in comparison to

the whole signal if the input is a single waveform.

nil Returns nil and an error message otherwise.

## **Example**

```
overshoot( v( "/net8" ) 7n t 3.99u t )
```

Returns the value of the overshoot for the waveform representing the voltage of "/net8".

```
overshoot(VT("/out") 0.5 nil 4.95 nil 5 t 'time)
```

Returns multiple occurrences of overshoot specified against time-points at which each overshoot event occurs.

```
overshoot(VT("/out") 0.5 nil 4.95 nil 5 t 'cycle)
```

Returns multiple occurrences of overshoot specified against cycle numbers, where a cycle number refers to the n'th occurrence of the overshoot event in the input waveform.

## Predefined and Waveform (Calculator) Functions

## pavg

```
pavg(
    o_waveform
    n_from
    n_to
    [ n_period [ n_sfactor ] ]
)
    => o_waveform / nil
```

## **Description**

Computes the periodic average of a family of signals for each time point.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like srrWave:XXXXX.).
n_from	Starting numeric value for the range on the X-axis.
n_to	Ending numeric value for the range on the X-axis.
n_period	Numeric value for the period of the input waveform.
n_sfactor	Sampling factor. This can be increased in order to increase the accuracy of the output.
	Default value: 1

## **Values Returned**

o_waveform	Returns a waveform representing the periodic average of a family of signals.
nil	Returns nil and an error message otherwise.

## **Example**

```
pavg( v("/net8") ?from 1n ?to 20n ?period 2n ?sfactor 1)
```

Returns the value of the periodic average for the family of waveforms representing the voltage of "/net8".

## Predefined and Waveform (Calculator) Functions

## peak

## **Description**

Detects the peaks in the input waveform and returns the X and Y coordinates of these peak points in the form of a waveform.

**Note:** The function will not work for waveforms that comprise of complex numbers.

## **Arguments**

o_waveform	Input waveform.
?from f_from	The initial point on the given waveform to start determining the peaks. By default, the first point of the waveform is the starting point.
?to <i>f_to</i>	The final point on the given waveform up to which the peaks are to be determined. By default, the last point of the waveform is the end point.
?xtol f_xtol	The distance on the X axis within which all peaks are to be filtered.Default: 0.0
?ytol f_ytol	The distance on the Y axis within which all peaks are to be filtered. Default: 0.0
?withLast $g_withLast$	Determines whether to include the last point in the peak calculation or not. When this argument is set to $t$ , the last point is included in the plot if it is a higher than the previous point. When set to $nil$ , the plot stops at the last peak.

**Note:** If both  $f_xtol$  and  $f_ytol$  are specified, the filtering mechanism will operate as follows:

## Predefined and Waveform (Calculator) Functions

- The maximum peak is selected first.
- All adjacent peaks in the neighborhood of both *f\_xtol* in the X-axis direction and *f ytol* in the Y-axis direction are then filtered.
- Next, all the peaks in the rectangular window thus formed are filtered based on both f xtol and f ytol.

If only one of  $f_xtol$  or  $f_ytol$  is specified, the peaks are filtered only in either the X-axis direction or the Y-axis direction, respectively.

#### Value Returned

o_waveform	Returns a waveform whose X and Y coordinates of the peaks are determined from the input waveform and the peaks are filtered based on the $f_xtol$ and $f_ytol$ criteria.
nil	Returns nil and an error message otherwise.

## **Example**

```
peak( vt("/out") ?from 1n ?to 20n ?xtol 2n ?ytol 0.5)
```

Out of all the peaks in the region starting from 1n to 20n, the function returns a waveform comprising of some of these peaks that satisfy the criteria of x-t01 (2n) and yt01 (0.5).

## Predefined and Waveform (Calculator) Functions

## peakToPeak

```
peakToPeak(
    o_waveform
    [ ?overall type_overall ]
    )
    => o waveform / n value / nil
```

## **Description**

Returns the difference between the maximum and minimum values of a waveform.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
?overall overall	

### Value Returned

o_waveform	Returns a waveform or a family of waveforms if the input argument is a family of waveforms.
n_value	Returns the difference between the maximum and minimum values of a waveform if the input argument is a single waveform.
nil	Returns nil and an error message otherwise.

## **Example**

```
peakToPeak( v( "/net2" ) )
```

Returns the difference between the maximum and minimum values of the waveform representing the voltage of the "/net2" net.

## Predefined and Waveform (Calculator) Functions

## period\_jitter

```
period_jitter(
    o_waveform
    t_crossType
    [ ?mode t_mode ]
    [ ?threshold n_threshold ]
    [ ?binSize n_binSize ]
    [ ?xName t_xName ]
    [ ?outputType t_outputType ]
    )
    => o_waveform / f_val / nil
```

## **Description**

Computes the period jitter. It returns a waveform or a value representing deviation from the average period.

#### Predefined and Waveform (Calculator) Functions

## **Arguments**

o waveform Name of the signal, expression, or a family of

waveforms.

t crossType The points at which the curves of the waveform

intersect with the threshold. While intersecting, the

curve may be either rising or falling.

Valid values: rising and falling.

Default value: rising.

?mode t\_mode The mode for calculating the threshold. Valid values:

auto and user.

If set to user, an n threshold value needs to be

specified by you.

If set to auto, n threshold is calculated as:

Auto Threshold Value = integral of the waveform

divided by the X range

Default value: auto

?threshold n threshold The threshold value against which the period is to be

calculated. It needs to be specified only when the

mode selected is user.

?binSize *n binSize* The width of the moving average window.The

deviation of value at the particular point from the

average of this window is the jitter.

If binsize=0, all periods are used to calculate the

average.

If binsize=N, the last N periods are used to calculate

the average.

Default value: 0

?xName t xName The X-axis of the output waveform. It specifies

whether you want to retrieve the period jitter against time (or another X-axis parameter for non-transient data) or cycle. Cycle numbers refer to the n'th

occurrence of the delay event in the input waveform.

Valid values: time and cycle.

Default value: time

## Predefined and Waveform (Calculator) Functions

?outputType t\_outputType Type of output.Valid values: sd and plot.

If set to plot, the output is a jitter waveform.

If set to sd, the output is a standard deviation of the

jitter waveform.

Default value: plot

#### Value Returned

o waveform Returns the period jitter values as a function of time or

cycle when the output Type is set to plot.

f val Returns the standard deviation value when the

output Type is set to sd.

nil Returns nil otherwise.

### **Example**

period\_jitter( wave1 "rising" ?mode "user" ?threshold 1 ?binSize 2 ?xName "cycle"
?outputType "sd" )
=> 1.695467

Returns the standard deviation for the period jitter of wave1 with the threshold of 1 against the cycle on the x-axis.

## Predefined and Waveform (Calculator) Functions

## phase

## **Description**

Gets the phase of the waveform or number. The phase command is similar to the phaseDegUnwrapped command and returns the unwrapped phase in degrees.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_number	Number.

#### Value Returned

o_waveform	Returns a waveform object if the input argument is a waveform object or returns a family of waveforms if the input argument is a family of waveforms.
n_number	Returns a number if the input argument is a number.
nil	Returns nil and an error message otherwise.

## **Example**

```
phase( v( "5" ) )
```

Gets the phase of the waveform representing the voltage at net 5. You can also use the vp alias to perform the same command, as in vp ( "5" ).

```
phase( i( "VFB" ) )
```

Gets the phase of the waveform representing the current through the VFB component. You can also use the ip alias to perform the same command, as in ip ( "VFB" ).

```
phase(-2.0) => 180.0
```

Gets the phase of -2.

## Predefined and Waveform (Calculator) Functions

## phaseDeg

## **Description**

Calculates the wrapped phase in degrees of a waveform and returns a waveform.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_number	Number.

## **Value Returned**

o_waveform	Returns a waveform object representing the wrapped phase in degrees of the input waveform. Returns a family of waveforms if the input argument is a family of waveforms.
n_number	Returns a number if the input argument is a number.
nil	Returns nil and an error message otherwise.

## **Example**

```
phaseDeg( v( "vout" ) )
```

Takes the input waveform, representing the voltage of the "vout" net, and returns the waveform object representing the wrapped phase in degrees.

## Predefined and Waveform (Calculator) Functions

## phaseDegUnwrapped

## **Description**

Calculates the unwrapped phase in degrees of a waveform and returns a waveform.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_number	Number.

### Value Returned

o_waveform	Returns a waveform object representing the unwrapped phase in degrees of the input waveform. Returns a family of waveforms if the input argument is a family of waveforms.
n_number	Returns a number if the input argument is a number.
nil	Returns nil and an error message otherwise.

## **Example**

```
phaseDegUnwrapped( v( "vout" ) )
```

Takes the input waveform, representing the voltage of the "vout" net, and returns the waveform object representing the unwrapped phase in degrees.

# Predefined and Waveform (Calculator) Functions

## phaseMargin

```
phaseMargin(
     o waveform
    => o waveform / n value / nil
```

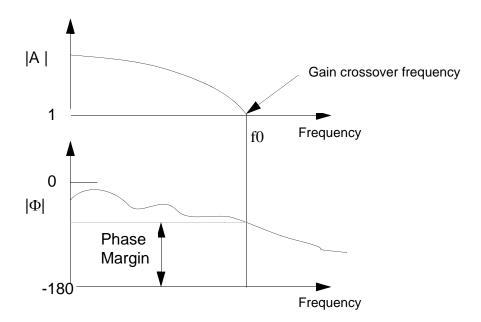
## **Description**

Computes the phase margin of the loop gain of an amplifier.

You supply a waveform representing the loop gain of interest over a sufficiently large frequency range.

```
phaseMargin(gain) = 180 + phase(value(gain f0))
```

The phase margin is calculated as the difference between the phase of the gain in degrees at f0 and at -180 degrees. The frequency f0 is the lowest frequency where the gain is 1. For stability, the phase margin must be positive.



## Predefined and Waveform (Calculator) Functions

## **Arguments**

o\_waveform Waveform object representing simulation results that can be

displayed as a series of points on a grid. (A waveform object

identifier looks like this: srrWave: XXXXX.)

#### Value Returned

o\_waveform Returns a waveform representing the phase margin of the loop

gain of an amplifier for a family of waveforms if the input

argument is a family of waveforms.

n value Returns the value (in degrees) equivalent to the phase margin

of the input waveform.

nil Returns nil and an error message otherwise.

#### **Example**

```
phaseMargin( v( "/OUT" ) )
```

Returns the phase margin for the waveform representing the voltage of the "/OUT" net.

## Predefined and Waveform (Calculator) Functions

## phaseRad

## **Description**

Calculates the wrapped (discontinuous) phase in radians of a waveform.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_number	Number.

## Value Returned

o_waveform	Returns a waveform representing a discontinuous value (in radians) for the phase of the input waveform. Returns a family of waveforms if the input argument is a family of waveforms.
n_number	Returns a number when the input argument is a number.
nil	Returns nil and an error message otherwise.

## **Example**

```
plot( phaseRad( v( "/OUT" ) ) )
```

Returns the wrapped phase of the waveform representing the voltage of the "/OUT" net.

## Predefined and Waveform (Calculator) Functions

## phaseRadUnwrapped

```
phaseRadUnwrapped(
    o_waveform
)
=> o waveform / nil
```

## **Description**

Calculates the unwrapped (continuous) phase in radians of a waveform and returns a waveform.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be
	displayed as a series of points on a grid. (A waveform object

identifier looks like this: srrWave: XXXXX.)

#### Value Returned

o waveform	Ret	urns a wavet	torm represent	ıng the	unwrapped	(continuous)

value for the phase of the input waveform in radians. Returns a

family of waveforms if the input argument is a family of

waveforms.

nil Returns nil and an error message otherwise.

### Example

```
plot( phaseRadUnwrapped( v( "/OUT" ) )
```

Returns the unwrapped phase of the waveform representing the voltage of the "/OUT" net.

## Predefined and Waveform (Calculator) Functions

## PN

```
PN(
    o_waveform
    t_crossType
    n_threshold 1.0
    [?windowName t_windowName]
    [?smooth x_smooth]
    [?windowsize x_windowsize]
    [?detrending t_detrending]
    [?cohGain f_cohGain]
)
    => o waveform / nil
```

## **Description**

Calculates the transient phase noise of the input waveforms in decibels (dBc/Hz). Phase noise is defined as the power spectral density of the absolute jitter of an input waveform.

#### Predefined and Waveform (Calculator) Functions

#### Arguments

o waveform

Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)

?crossType t crossType

The points at which the curves of the waveform intersect with the threshold. While intersecting, the curve may be either rising or falling.

Valid values: rising and falling, respectively.

Default crossType is rising.

?windowName t windowName

The window type. Valid values: 'Blackman, 'Cosine2,

'Cosine4, 'ExtCosBell, 'HalfCycleSine,

'HalfCycleSine3 or 'HalfCycleSine6, 'Hamming,

'Hanning, 'Kaiser, 'Parzen, 'Rectangular,

'Triangular, or 'Nuttall.

Default value: 'Rectangular

?smooth x smooth

The Kaiser window smoothing parameter. The 0 value requests no smoothing. Valid values:  $0 \le x \mod th \le 15$ .

Default value: 1

?windowsize x windowsize

The number of frequency domain points to use in the Fourier analysis. A larger window size results in an expectation operation over fewer samples, which leads to larger variations in the power spectral density. A small window size can smear out sharp steps in the power spectral density that might really be present.

Default value: 256

?detrending t detrending

The detrending mode to use. Valid values: 'None, 'mean,

'Linear

Default value: 'Mean

?cohGain f cohGain

## Predefined and Waveform (Calculator) Functions

A scaling parameter. A non-zero value scales the power spectral density by  $1/(f\_cohGain)$ . Valid values: none, default, magnitude, dB20, or dB10

Default value: db20

#### Value Returned

o\_waveform The power spectral density waveform returned when the

command is successful.

nil Returns nil when the command fails.

## **Example**

```
PN(v("net9") "rising" 1.0 ?windowName "Rectangular" ?smooth 1 ?windowSize 256 ?detrending "Mean" ?cohGain (10**(/20)))
```

Returns the Phase Noise waveform, net9, for the window type rectangular at threshold value 1.0.

## Predefined and Waveform (Calculator) Functions

## pow

## **Description**

Takes the exponent of a given waveform or number.

## **Arguments**

o_waveformBase	Waveform object to be used as the base for the expression.
o_waveformExpn	Waveform object to be used as the exponent for the expression.
n_numberBase	Number to be used as the base for the expression.
$n\_numberExpn$	Number to used as the exponent for the expression.

#### Value Returned

o_waveform	Returns a family of waveforms if one of the input arguments is a family of waveforms or returns a waveform if one of the input arguments is a waveform (and none is a family).
n_result	Returns a number if both the input arguments are numbers.
nil	Returns nil and an error message otherwise.

## **Example**

```
pow( average( v( "/net9" ) ) 0.5 )
```

Gets the square root of the average value of the voltage at "/net9".

```
pow(23)
```

Gets the value of 2 to the third power, or 8.

```
pow( -2 2 )
=> 4
```

Gets the value of -2 to the second power.

## Predefined and Waveform (Calculator) Functions

pow(2.5 -1.2) => 0.3330213

Gets the value of 2.5 to the power of -1.2.

## Predefined and Waveform (Calculator) Functions

#### prms

## **Description**

Computes the periodic root mean square of a family of signals for each time point, which is the square root of the periodic average of the square of the input waveform.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like srrWave:XXXXX.).
n_from	Starting numeric value for the range on the X-axis.
n_to	Ending numeric value for the range on the X-axis.
n_period	Numeric value for the period of the input waveform.
n_sfactor	Sampling factor. This can be increased in order to increase the accuracy of the output.
	Default value: 1

#### **Values Returned**

o_waveform	Returns a waveform representing the periodic root mean square of a family of signals.
nil	Returns nil and an error message otherwise.

## **Example**

```
prms v("/net8") ?from 1n ?to 20n ?period 2n ?sfactor 1)
```

Returns the value of the periodic root mean square for the family of waveforms representing the voltage of "/net8".

## Predefined and Waveform (Calculator) Functions

## psd

```
psd(
    o_waveform
    f_timeStart
    f_timeEnd
    x_num
    [?windowName t_windowName]
    [?smooth x_smooth]
    [?cohGain f_cohGain]
    [?windowsize x_windowsize]
    [?detrending t_detrending]
)
    => o_waveformReal / nil
```

## **Description**

Returns an estimate for the power spectral density of  $o\_waveform$ . If  $x\_windowsize$  is not a power of 2, it is forced to the next higher power of 2. If  $x\_num$  is less than x windowsize, x num is forced to x windowsize.

#### Predefined and Waveform (Calculator) Functions

#### **Arguments**

o waveform Time domain waveform object with units of volts or amps.

f timeStart Starting time for the spectral analysis interval. Use this

parameter and  $f\_timeEnd$  to exclude part of the interval. For example, you might set these values to discard initial transient

data.

f timeEnd Ending time for the spectral analysis interval.

x num The number of time domain points to use. The maximum

frequency in the Fourier analysis is proportional to x num and

inversely proportional to the difference between

f timeStart and f timeEnd.

Default value: 512

?windowName t windowName

The window to be used for applying the moving window FFT.

Valid values: Blackman, Cosine2, Cosine4, ExtCosBell, HalfCycleSine, Half3CycleSine or HalfCycleSine3, Half6CycleSine or HalfCycleSine6, Hamming, Hanning, Kaiser, Parzen, Rectangular, Triangle, Triangular,

or Nuttall.

Default value: Hanning

?smooth  $x\_smooth$  The Kaiser window smoothing parameter. The 0 value requests

no smoothing. Valid values:  $0 \le x \mod h \le 15$ .

Default value: 1

?cohGain f cohGain

A scaling parameter. A non-zero value scales the power

spectral density by  $1/(f\_cohGain)$ .

Valid values:  $0 < f\_cohGain < 1$  (You can use 1 if you do not

want the scaling parameter to be used)

Default value: 1

?windowSize x windowsize

### Predefined and Waveform (Calculator) Functions

The number of frequency domain points to use in the Fourier analysis. A larger window size results in an expectation operation over fewer samples, which leads to larger variations in the power spectral density. A small window size can smear out sharp steps in the power spectral density that might really be present.

Default value: 256

?detrending t\_detrending

The detrending mode to use.

Valid values: mean, linear, none

Default value: none

The psd function works by applying a moving windowed FFT to time-series data. If there is a deterministic trend to the underlying data, you might want to remove the trend before performing the spectral analysis. For example, consider analyzing phase noise in a VCO model. Without the noise, the phase increases more or less linearly with time, so it is appropriate to set the detrending mode to 'linear. To subtract an average value, set the detrending mode to 'mean. Where the spectrum of raw data is desired, set the detrending mode to none.

## Value Returned

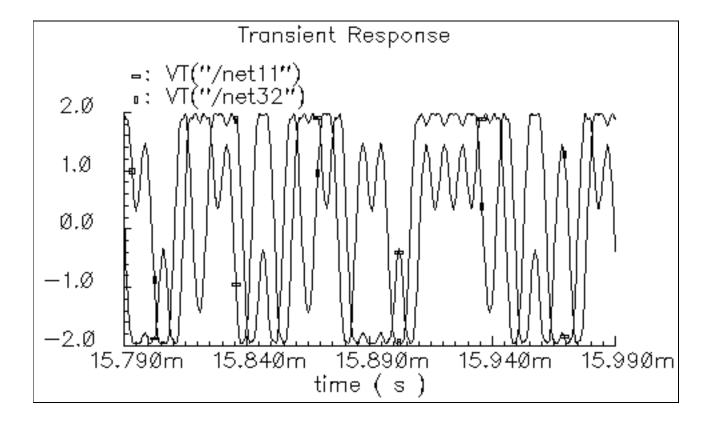
o\_waveformReal The power spectral density waveform returned when the command is successful.

nil Returns nil when the command fails.

#### **Example**

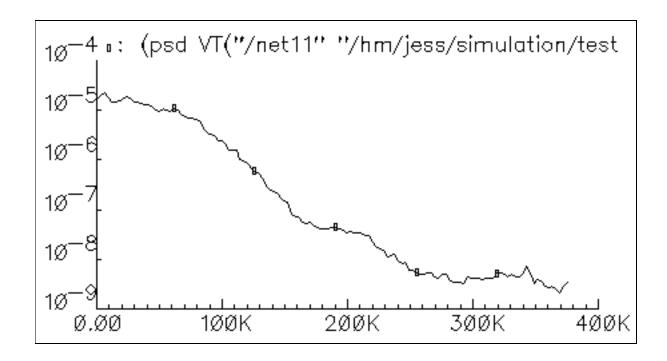
Predefined and Waveform (Calculator) Functions

Consider applying this command to one of the waveforms in the following illustration.



# OCEAN Reference Predefined and Waveform (Calculator) Functions

The result is the following spectrum, which is displayed with a logarithmic vertical scale.



## Predefined and Waveform (Calculator) Functions

## psdbb

```
psdbb(
    o_waveform1
    o_waveform2
    f_timeStart
    f_timeEnd
    x_num
    [?windowName t_windowName]
    [?smooth x_smooth]
    [?cohGain f_cohGain]
    [?windowsize x_windowsize]
    [?detrending t_detrending]
)
    => o_waveformReal / nil
```

## **Description**

Returns an estimate for the power spectral density of  $o\_waveform1+j*o\_waveform2$ . If  $x\_windowsize$  is not a power of 2, it is forced to the next higher power of 2. If  $x\_num$  is less than x windowsize, x num is forced to x windowsize.

### Predefined and Waveform (Calculator) Functions

#### **Arguments**

o\_waveform1 Time domain waveform object with units of volts or amps.

o waveform2 Time domain waveform object with units of volts or amps.

f timeStart Starting time for the spectral analysis interval. Use this

parameter and  $f\_timeEnd$  to exclude part of the interval. For example, you might set these values to discard initial

transient data.

f timeEnd Ending time for the spectral analysis interval.

x num The number of time domain points to use. The maximum

frequency in the Fourier analysis is proportional to x num and

inversely proportional to the difference between

f timeStart and f timeEnd.

?windowName t windowName

The window to be used for applying the moving window FFT.

Valid values: Blackman, Cosine2, Cosine4, ExtCosBell, HalfCycleSine, Half3CycleSine Or HalfCycleSine3, Half6CycleSine Or HalfCycleSine6, Hamming, Hanning, Kaiser, Parzen, Rectangular, Triangle,

Triangular, or Nuttall.

Default value: Hanning

?smooth x smooth The Kaiser window smoothing parameter. 0 requests no

smoothing.

Valid values:  $0 \le x \mod h \le 15$ .

Default value: 1

?cohGain f cohGain A scaling parameter. A non-zero value scales the power

spectral density by 1/(f\_cohGain). Valid values: 0 <

f cohGain < 1 (You can use 1 if you do not want the scaling

parameter to be used)

Default value: 1

?windowSize x windowsize

### Predefined and Waveform (Calculator) Functions

The number of frequency domain points to use in the Fourier analysis. A larger window size results in an expectation operation over fewer samples, which leads to larger variations in the power spectral density. A small window size can smear out sharp steps in the power spectral density that might really be present.

Default value: 256

?detrending t\_detrending

The detrending mode to use.

Valid values: mean, linear, none

Default value: none

The psd function works by applying a moving windowed FFT to time-series data. If there is a deterministic trend to the underlying data, you might want to remove the trend before performing the spectral analysis. For example, consider analyzing phase noise in a VCO model. Without the noise, the phase increases more or less linearly with time, so it is appropriate to set the detrending mode to 'linear. To subtract an average value, set the detrending mode to 'mean. Where the spectrum of raw data is desired, set the detrending mode to 'none.

#### Value Returned

o\_waveformReal The power spectral density waveform returned when the

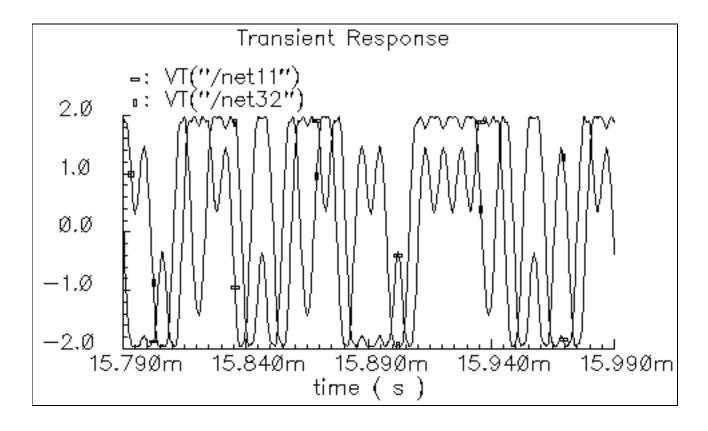
command is successful.

nil Returns nil when the command fails.

#### Example

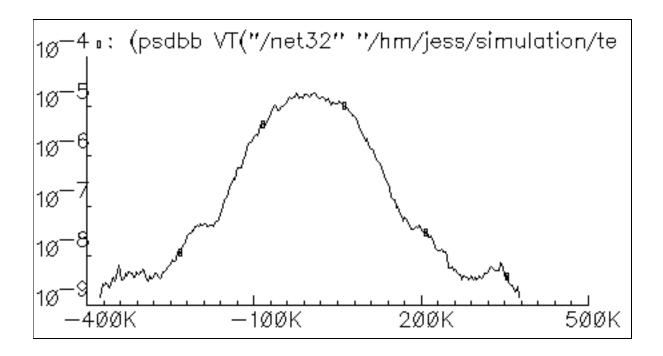
## Predefined and Waveform (Calculator) Functions

Consider applying this command to both of the waveforms in the following illustration.



## Predefined and Waveform (Calculator) Functions

The result is the following spectrum, which is displayed with a logarithmic vertical scale.



## Predefined and Waveform (Calculator) Functions

## pstddev

```
pstddev(
    o_waveform
    n_from
    n_to
    [ n_period [ n_sfactor ] ]
    )
    => o_waveform / nil
```

### **Definition**

Computes the periodic standard deviation of a family of signals for each time point.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like srrWave:XXXXX.).
$n\_from$	Starting numeric value for the range on the X-axis.
n_to	Ending numeric value for the range on the X-axis.
n_period	Numeric value for the period of the input waveform.
n_sfactor	Sampling factor. This can be increased in order to increase the accuracy of the output. Default value: 1

#### **Values Returned**

o_waveform	Returns a waveform representing the periodic standard deviation of a family of signals.
nil	Returns nil and an error message otherwise.

## **Example**

```
pstddev( v("/net8") ?from 1n ?to 20n ?period 2n ?sfactor 1)
```

Returns the value of the periodic standard deviation for the family of waveforms representing the voltage of "/net8"

# Predefined and Waveform (Calculator) Functions

# pzbode

```
pzbode(
    f_transferGain
    f_minfrequency
    f_maxfrequency
    x_nponits
    [ ?poles o_waveform1 ]
    [ ?zeros o_waveform2 ]
    )
    => o waveform / nil
```

# **Description**

Calculates and plots the transfer function of a circuit from pole zero simulation data.

**Note:** This command also works for the parametric or sweep data.

### Predefined and Waveform (Calculator) Functions

### **Arguments**

f transferGain The transfer gain constant.

f\_minfrequency The minimum frequency for the bode plot.

f\_maxfrequency The maximum frequency for the bode plot.

 $x_npoints$  The frequency interval for the bode plot, in points per decade.

?poles waveform1 Poles from the dumped simulation data.

Default value: all

?zeros o waveform2

Zeros from the dumped simulation data.

Default value: all

#### Value Returned

o waveform Waveform containing the X and Y points of the transfer function.

The scale of the Y-axis will be db20.

nil Returns nil and error message otherwise.

#### **Example**

pzbode( 1.0 1M 1G 20 ?poles complexPoleWave ?zeros complexZeroWave )

## Predefined and Waveform (Calculator) Functions

# pzfilter

```
pzfilter(
    [ o_PoleWaveform ]
    [ o_ZeroWaveform ]
    [ ?maxfreq t_maxfreq ]
    [ ?reldist n_reldist ]
    [ ?absdist n_absdist ]
    [ ?minq n_minq ]
    [ ?output_type o_output ]
)
=> o waveform / nil
```

# **Description**

Returns the filtered Pole and Zero waveforms.

**Note:** If you do not specify values for  $o\_PoleWaveform\ and\ o\_ZeroWaveform$  arguments, you should have run pz analysis prior to using this function. This command also works for the parametric or sweep data.

### Predefined and Waveform (Calculator) Functions

### **Arguments**

o\_PoleWaveform Input Pole waveform (complex points).

Default value: Poles of the simulator pz-analysis dump

o ZeroWaveform Input Zero waveform (complex points).

Default value: Zeros of the simulator pz-analysis dump

t maxfreq Maximum frequency.

Default value: 1e10

?reldist n\_reldist

Relative distance to be considered while filtering.

Default value: 0.05

?absdist  $n_absdist$ 

Absolute distance to be considered while filtering.

Default value: 1e-6

 $?minq n_minq$  Minimum q factor to be allowed while filtering.

?output type o output

Specifies the type of the output. If this argument is not passed, the output is a family of waves with two child waveforms, representing poles and zeros respectively, with the real component of each waveform as the X values and the imaginary components as the Y values.

Valid value: complexwave. The output is a family of waves with two child waves, both of which are complex and represent poles and zeros, respectively.

#### Value Returned

o\_waveform Returns a family (waveform) of Pole and Zero waveforms.

nil Returns nil otherwise.

#### Example

pzfilter( complexPoleWave complexZeroWave )
=> srrWave:175051584

# Predefined and Waveform (Calculator) Functions

Returns a family of filtered Pole and Zero waveforms, which correspond to the sweep values of "Pole" and "Zero", respectively.

## Predefined and Waveform (Calculator) Functions

# rapidIPNCurves

```
rapidIPNCurves(
    o_result
    [ ?resultsDir t_resultsDir ]
    [ ?resistance n_resistance ]
    @Rest args
)
    => o waveformReal / nil
```

### **Description**

Plots IPN curves.

# **Arguments**

 $o\_result$  Object representing simulation results that can be displayed as

a series of points on a grid.

?resultsDir t\_resultsDir

Name of the directory where results are saved.

?resistance n\_resistance

Value of resistance

Default value: 50

@Rest 1 args List of arguments to be used by the value function on the results

data. Refer to the value function for more details.

#### Value Returned

o waveformReal Returns a waveform.

nil Returns nil or an error message otherwise.

### **Example**

```
w2 = rapidIPNCurves("ac-ip3" ?resultsDir "./simulation/amplifier/spectre/
schematic/psf" ?r 50)
```

### Predefined and Waveform (Calculator) Functions

# rapidIIPN

```
rapidIIPN(
    o_result
    [ ?resultsDir t_resultsDir ]
    [ ?resistance n_resistance ]
    @Rest args
)
    => o waveform / nil
```

### **Description**

Plots the input IPN curves.

# **Arguments**

o\_result Object representing simulation results that can be displayed as

a series of points on a grid.

?resultsDir t\_resultsDir

Name of the directory where results are saved.

?resistance n resistance

Value of resistance

Default value: 50

1 args List of arguments to be used by the value function on the results

data. Refer to the value function for more details.

#### Value Returned

o waveform Returns a waveform.

nil Returns nil or an error message otherwise.

### **Example**

rapidIIPN("hbac ip3")

### Predefined and Waveform (Calculator) Functions

### real

```
real(
     { o_waveform | n_input }
)
=> o waveformReal / n numberReal / nil
```

### **Description**

Returns the real part of a waveform representing a complex number, or returns the real part of a complex number.

### **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_input	Complex number.

#### Value Returned

o_waveformReal	Returns a waveform when the input argument is a waveform.
n_numberReal	Returns a number when the input argument is a number.
nil	Returns nil and an error message otherwise.

### **Example**

```
real( v( "/net8" ) )
```

Returns a waveform representing the real part of the voltage of "/net8". You also can use the vr alias to perform the same command, as in vr ( "net8").

```
x=complex(-1,-2) => complex(-1,-2)
real(x) => -1.0
```

Creates a variable  $\mathbf{x}$  representing a complex number, and returns the real portion of that complex number.

## Predefined and Waveform (Calculator) Functions

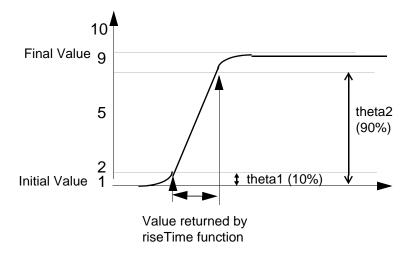
## riseTime

```
riseTime(
    o_waveform
    n_initVal
    g_initType
    n_finalVal
    g_finalType
    n_theta1
    n_theta2
    [ g_multiple [ s_Xname ] [ g_histoDisplay ] [ x_noOfHistoBins ] ]
    )
    => o waveform / n value / nil
```

# **Description**

Returns the rise time measured between theta1 (percent low) to theta2 (percent high) of the difference between the initial value and the final value.

The riseTime function can also be used to compute the fall time if initVal is higher than finalVal.



# Predefined and Waveform (Calculator) Functions

# **Arguments**

J	
o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_initVal	Initial value at which to start the computation.
g_initType	Specifies how n_initVal functions.
	Valid values: a non-nil value specifies that the initial value is taken to be the value of the waveform, interpolated at $n\_initVal$ , and the waveform is clipped from below as follows:
	<pre>o_waveform = clip( o_waveform g_initVal nil )</pre>
	where nil specifies that $n\_initVal$ is defined by the X value entered. (The command gets the Y value for the specified X value and uses that value for $n\_initVal$ .)
n_finalVal	Final value at which to end the computation.
g_finalType	Specifies how the $n_{finalVal}$ argument functions.
	Valid values: a non-nil value specifies that the final value is taken to be the value of the waveform, interpolated at $n\_finalVal$ , and the waveform is clipped from above, as follows:
	<pre>o_waveform = clip(o_waveform nil n_finalVal)</pre>
	where nil specifies that the $n\_finalVal$ argument is defined by the X value entered. (The command gets the Y value for the specified X value and uses that value for $n\_finalVal$ .)
n_theta1	Percent low.
n_theta2	Percent high.
g_multiple	An optional boolean argument that takes the value $\tt nil$ by default. If set to $\tt t$ , the function returns multiple occurrences of the riseTime event.
s_xName	An optional argument that is used only when $g_{multiple}$ is set to t. It takes the value time by default. It controls the contents of the x vector of the waveform object returned by the function.
	Valid values: 'time, 'cycle

### Predefined and Waveform (Calculator) Functions

 $g_histoDisplay$  When set to t, returns a waveform that represents the statistical

distribution of the riseTime data in the form of a histogram. The

height of the bars (bins) in the histogram represents the frequency of the occurrence of values within the range of

riseTime data.

Valid values: t nil

Default value: nil

x noOfHistoBins Denotes the number of bins represented in the histogram

representation.

Valid values: Any positive integer

Default value: nil

**Note:**  $g_histoDisplay$  and  $x_noOfHistoBins$  are added for backward compatibility only. It will be deprecated in future releases. Use the histo function for plotting the histogram of the resulting function.

#### **Value Returned**

o_waveform	Returns a waveform representing the rise time for a family of waveforms if the input argument is a family of waveforms or if $g_{multiple}$ is set to t.
n_value	Returns a value for the rise time if the input is a single waveform.
nil	Returns nil and an error message otherwise.

### **Example**

```
riseTime( v( "/net8" ) 0 t 2 t 10 90 )
```

Computes the rise time for the waveform representing the voltage of "/net8" from 0 to 2.

For the next example, assume that v is the following sinusoidal waveform:

```
sin( 2 * pi * time)
riseTime( v 0.25 t 0.5 t 10 90)
```

Computes the fall time of the first falling edge from 1 to 0.

```
riseTime(VT("/out") 0.5 nil 4.5 nil 10 90 t "time") (s)
```

# Predefined and Waveform (Calculator) Functions

Returns multiple occurrences of riseTime specified against time-points at which each riseTime event occurs.

```
riseTime(VT("/out") 0.5 nil 4.5 nil 10 90 t "cycle") (s)
```

Returns multiple occurrences of riseTime specified against cycle numbers, where a cycle number refers to the n'th occurrence of the riseTime event in the input waveform.

# Predefined and Waveform (Calculator) Functions

#### rms

```
rms(
    o_waveform
)
=> o_waveform / n_value / nil
```

# **Description**

Returns the root-mean-square value of a waveform.

# **Arguments**

o_waveform	Waveform object representing simulation results that can be	
	displayed as a series of points on a grid. (A waveform object	
	identifier looks like this: srrWave: XXXXX.)	

### Value Returned

o_waveform	Returns a waveform representing the root-mean-square value for a family of waveforms if the input argument is a family of waveforms.
n_value	Returns a value for the root-mean-square value for the specified waveform if the input is a single waveform.
nil	Returns nil and an error message otherwise.

# **Example**

```
rms( v( "/out" ) )
```

Returns the root-mean-square value of the waveform representing the voltage of the "/out" net.

# Predefined and Waveform (Calculator) Functions

# rmsNoise

```
rmsNoise(
    n_from
    n_to
)
    => o waveform / n value / nil
```

# **Description**

Computes the integrated root-mean-square noise over the specified bandwidth.

# **Arguments**

n_from	Frequency in hertz that specifies the minimum value for the bandwidth.
n_to	Frequency in hertz that specifies the maximum value for the bandwidth.

## **Value Returned**

o_waveform	Returns a waveform (or a family of waveforms) representing the integrated root-mean-square noise if the data being analyzed is parametric.
n_value	Returns a value for the integrated root-mean-square noise if the data being analyzed is from a single simulation run.
nil	Returns nil and an error message otherwise.

# Example

```
rmsNoise( 100 100M) => 250e-6
```

Computes the integrated root-mean-square noise from 100 to 100M.

# Predefined and Waveform (Calculator) Functions

# rmsVoltage

```
rmsVoltage(
    t_net
    [ t_net1 ]
)
=> f voltage / nil
```

# **Description**

Calculates the root-mean-square voltage between two nets for fast and regular envelop analysis.

# **Arguments**

t_net	Name of the net selected in the schematic.
t_net1	Name of the second net selected in the schematic. This argument is optional. If not specified, the default value is assumed as gnd.

#### Value Returned

f_voltage	Returns a value in terms of voltage.
nil	Returns nil and an error message otherwise.

### **Example**

```
rmsVoltage( "net1" "!gnd")
=> 120
```

Calculates the root-mean-square voltage between net1 and qnd.

### Predefined and Waveform (Calculator) Functions

# rmsTerminalVoltage

```
rmsTerminalVoltage(
    t_terminal
    [ t_terminal1 ]
    )
    => f voltage / nil
```

## **Description**

Calculates the root-mean-square voltage between two terminals for fast and regular envlp analysis.

## **Arguments**

t_terminal	Name of the terminal selected in the schematic.
t_terminal1	Name of the second terminal selected in the schematic. This argument is optional. If not specified, the default value is assumed as gnd.

#### Value Returned

f_voltage	Returns a value in terms of voltage.
nil	Returns nil or an error message.

### **Example**

If the following expression is created and plotted:

The rms terminal voltage function creates the following expression:

```
rmsTerminalVoltage("IO/M1/D")
```

If the result is re-evaluated, the scalar value is added to the ADE output.

ı			
	testcase:mixer	RMS Voltage	1.506

# Predefined and Waveform (Calculator) Functions

### root

```
root(
    o_waveform
    n_rootVal
    x_n )
=> o waveform / n value / l value / nil
```

# Description

Returns the *n*th X value at which the Y value equals the specified Y value (root Val).

## **Arguments**

o_waveform	Waveform object representing simulation results that can be
	displayed as a series of points on a grid. (A waveform object
	identifier looks like this: srrWave: XXXXX.)

n root Val Y value of interest.

 $x_n$  Number that specifies which X value to return. If n equals 1, the

first X value that crosses over the Y rootVal is returned. If n equals 2, the second X value that crosses over the Y rootVal is returned, and so on. If you specify a negative integer for n, the X values that cross the rootVal are counted from right to left (from maximum to minimum). If you specify n as 0, the list of

root values is returned.

#### **Value Returned**

o_waveform	Returns a waveform if the input argument is a family of
	waveforms.

n value Returns an X value when the input argument is a single

waveform.

1 value Returns a list of all the root values when n is 0.

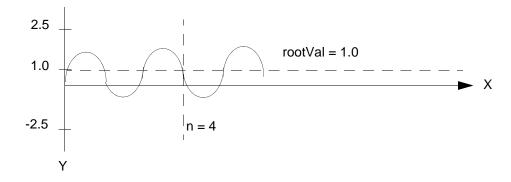
nil Returns nil and an error message otherwise.

# **Example**

```
root( v( "vout" ), 1.0, 4 )
```

# Predefined and Waveform (Calculator) Functions

Returns the X value for the point at which the waveform curve crosses the 1.0 Y value for the fourth time.



# Predefined and Waveform (Calculator) Functions

# rshift

```
rshift(
    o_waveform
    n_delta
)
=> o waveform / nil
```

# **Description**

Shifts the waveform to the right by the n delta value.

This command is the inverse of the lshift command.

### **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_delta	Value by which the waveform is to be shifted.

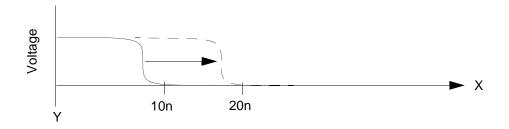
# **Value Returned**

o_waveform	Returns a waveform object. Returns a family of waveforms if the input argument is a family of waveforms.
nil	Returns nil and an error message otherwise.

### **Example**

```
rshift( v( "vout" ) ) 10n )
```

Shifts the waveform representing the voltage through the "vout" net to the right by 10n.



### Predefined and Waveform (Calculator) Functions

## sample

```
sample(
    o_waveform
    n_from
    n_to
    t_type
    n_by
)
=> o_waveform / n_number / nil
```

### **Description**

Samples a waveform at the specified interval.

You can use this function to reduce the time it takes to plot waveforms that have many data points. If you sample a waveform beyond its range, you get the final value of the waveform. You can use this function to demodulate a signal. Consider an AM modulated sine wave. Assume the carrier frequency is 1 GHz, and the modulation frequency is 1 MHz. If the waveform is sampled every 1 ns, the resulting signal is cleanly demodulated (the 1 GHz carrier is completely eliminated by the sampling).

**Note:** The function can be used to sample both a waveform object as well as a family of waveforms. If the family is of dimension m, the arguments  $n_from$ ,  $n_to$ , and  $n_by$  would be of dimension m-1.

# Predefined and Waveform (Calculator) Functions

# **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
$n\_from$	Starting value for the sampling.
n_to	Ending value for the sampling.
t_type	Type of the sampling.
	Valid values: "linear" or "log"

Interval at which to sample.

#### Value Returned

n by

o_waveform	Returns a waveform representing the sampling you specified.
n_number	Returns a number if the output contains only one point.
nil	Returns nil and an error message otherwise.

# **Example**

```
sample( v( "vout" ) 0 50n "linear" 0.1n )
```

Takes a linear sample of the waveform representing the voltage of the "vout" net.

```
sample( v( "vout" ) 0 100m "log" 10 )
```

Takes a logarithmic sample of the waveform representing the voltage of the "vout" net.

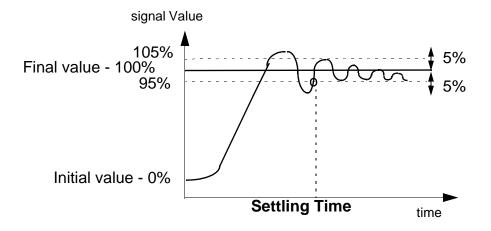
## Predefined and Waveform (Calculator) Functions

# settlingTime

```
settlingTime(
    o_waveform
    n_initVal
    g_initType
    n_finalVal
    g_finalType
    n_theta
    [ g_multiple [ s_Xname ] ]
    )
    => o_waveform / n_value / nil
```

# **Description**

The settling time is the time by which the signal settles within the specified Percent of step (theta) of the difference between the Final Value and Initial Value from the Final Value.



**Note:** The above graph represents the Initial value of the signal as 0% and Final value as 100%. The Percent of Step is taken as 5%.

## Predefined and Waveform (Calculator) Functions

### **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_initVal	Initial value at which to start the computation.
g_initType	Specifies whether the values entered are X values or Y values.
	Valid values: t specifies that $initVal$ is defined by the X value entered; nil specifies that $initVal$ is defined by the Y value entered
n_finalVal	Final value at which to start the computation.
g_finalType	Specifies whether the values entered are X values or Y values.
	Valid values: t specifies that $finalVal$ is defined by the X value entered; nil specifies that $finalVal$ is defined by the Y value entered
n_theta	Percent of the total step.g_multiple
	An optional boolean argument that takes the value $\mathtt{nil}$ by default. If set to $\mathtt{t}$ , the function returns multiple occurrences of the settlingTime event.
s_xName	An optional argument that is used only when $g_{multiple}$ is set to t. It takes the value time by default. It controls the contents of the x vector of the waveform object returned by the function.
	Valid values: 'time, 'cycle

#### **Additional Information**

The equation used to calculate maximum delta value is:

```
maxDeltaY = ((theta/100.0)*abs(FinalVal-InitVal))
```

Firstly, check if the absolute difference between the last element of the waveform and finalVal is less than maxDeltaY. If yes, then compute settlingTime, else returns nil.

To compute settlingTime, subtract finalVal from the waveform, get the subtracted-wave and calculate settling time as first cross on subtracted-wave at maxDeltaY (from opposite direction for falling edge). If no such crossing exists, then return 0.0.

```
maxDeltaY = ((theta/100.0) * abs(FinalVal -InitVal))
```

## Predefined and Waveform (Calculator) Functions

```
if( abs(last_Y_element_of_waveform - finalVal) < maxDeltaY then
    or( cross( abs(waveform - finalVal) maxDeltaY -1 -1) 0.0)
else
    nil
)</pre>
```

### Value Returned

o_waveform	Returns a waveform representing the settling time for a family of waveforms if the input argument is a family of waveforms or if $g_{multiple}$ is set to t.
n_value	Returns a value for the settling time for the specified waveform if the input is a single waveform.
nil	Returns nil and an error message otherwise.

### **Example**

```
settlingTime( v("/out" ) 0 t 2 t 90 )
```

Computes the time required for the waveform representing the voltage of the "/out" net to settle within 90 percent of the step from 0 to 2.

```
settlingTime(VT("/out") 0.5 nil 4.95 nil 5 t "time") (s)
```

Returns multiple occurrences of settlingTime specified against time-points at which each settlingTime event occurs.

```
settlingTime(VT("/out") 0.5 nil 4.95 nil 5 t "cycle") (s)
```

Returns multiple occurrences of settlingTime specified against cycle numbers, where a cycle number refers to the n'th occurrence of the settlingTime event in the input waveform.

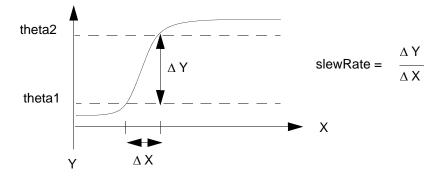
### Predefined and Waveform (Calculator) Functions

# slewRate

```
slewRate(
    o_waveform
    n_initVal
    g_initType
    n_finalVal
    g_finalType
    n_theta1
    n_theta2
    [ g_multiple [ s_Xname ] ]
    [ g_histoDisplay ] [ x_noOfHistoBins ]
)
    => o_waveform / n_value / nil
```

# **Description**

Computes the average rate at which an expression changes from theta1 (percent low) to theta2 (percent high) of the difference between the initial value and final value.



# Predefined and Waveform (Calculator) Functions

# **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_initVal	Initial X-axis value at which to start the computation.
g_initType	Specifies whether the values entered are X values or Y values.
	Valid values: t specifies that $initVal$ is defined by the X value entered; nil specifies that $initVal$ is defined by the Y value entered
n_finalVal	Final value at which to end the computation.
g_finalType	Specifies whether the values entered are X values or Y values.
	Valid values: t specifies that $finalVal$ is defined by the X value entered; nil specifies that $finalVal$ is defined by the Y value entered
n_theta1	Percent low (percentage of the total step).
n_theta2	Percent high (percentage of the total step).
g_multiple	An optional boolean argument that takes the value $\tt nil$ by default. If set to $\tt t$ , the function returns multiple occurrences of the slewRate event.
s_xName	An optional argument that is used only when $g_{multiple}$ is set to t. It takes the value time by default. It controls the contents of the x vector of the waveform object returned by the function.
	Valid values: 'time, 'cycle
g_histoDisplay	When set to t, returns a waveform that represents the statistical distribution of the riseTime data in the form of a histogram. The height of the bars (bins) in the histogram represents the frequency of the occurrence of values within the range of riseTime data.
	Valid values: t nil
	Default value: nil

### Predefined and Waveform (Calculator) Functions

 $x_noOfHistoBins$  Denotes the number of bins represented in the histogram

representation.

Valid values: Any positive integer

Default value: nil

**Note:**  $g_histoDisplay$  and  $x_noOfHistoBins$  are added for backward compatibility only. It will be deprecated in future releases. Use the histo function for plotting the histogram of the resulting function.

#### Value Returned

o_waveform	Returns a waveform representing the slew rate for a family of waveforms if the input argument is a family of waveforms or if $g_{multiple}$ is set to t.
n_value	Returns a value for the slew rate for the specified waveform if the input is a single waveform.
nil	Returns nil or an error message otherwise.

### **Example**

```
slewRate( v( "vout" ) 10n t 30n t 10 90 )
```

Computes the slew rate for the waveform representing the voltage of the "vout" net from 10n to 30n.

```
slewRate( v( "vout" ) 0 nil 10 nil 5 95 )
```

Computes the slew rate for the waveform representing the voltage of the "vout" net from 0 to 10. In this example, the initial value and final value are entered as Y values.

```
slewRate(VT("/out") 0.5 nil 4.5 nil 10 90 t 'time)
```

Return multiple occurrences of slewRate values, computed at different time-points.

```
slewRate(VT("/out") 0.5 nil 4.5 nil 10 90 t 'cycle)
```

Returns multiple occurrences of slewRate values specified against cycle numbers (where cycle number refers to the n'th occurrence of slewRate computation).

# Predefined and Waveform (Calculator) Functions

# smithType

```
smithType(
    x_mode
)
    => t / nil
```

## **Description**

Sets the Smith display mode type for the active graph.

### Arguments.

x mode

Type of Smith display.

Valid Values:

- impedance: Circular graph, also known as Z Smith, where the region above the x-axis repre-sents inductive impedances and the region below the x-axis represents capacitive impedances.
- admittance: Circular graph, also known as Y Smith, where the region above the x-axis repre-sents capacitive admittances and the region below the x-axis represents inductive admittances.
- polar: plot graph, representing data using the polar coordinates system.

For more information on circular graphs, see the <u>Creating a Circular Graph</u> section in *Virtuoso Visualization and Analysis XL User Guide*.s

### Value Returned

t Returns t when the specified Smith display id set.

nil Returns nil if there is an error.

# **Example**

```
smithType("impedance")
=>t
```

# Predefined and Waveform (Calculator) Functions

Sets the Smith display to impedance.

### Predefined and Waveform (Calculator) Functions

# spectralPower

```
spectralPower(
    o_current
    o_voltage
)
    => o power / nil
```

# **Description**

Returns the spectral power given the spectral current and voltage.

To obtain a list of the harmonic frequencies, use harmonicList.

# **Arguments**

o current Waveform representing the current. The current can be

obtained by calling the i data access function for the desired

terminal.

#### Value Returned

o power Waveform representing the power of the net.

nil Returns nil if there is an error.

#### **Example**

```
plot(db10(spectralPower(i("/PORT0/PLUS") v("/net28"))))
```

Plots power of the output "/net28". "/PORT0/PLUS" is a member of "/net28".

## Predefined and Waveform (Calculator) Functions

# spectrumMeas

```
spectrumMeas(
    o_waveform
    n_from
    n_to
    x_numSamples
    x_noiseBins
    n_startFreq
    n_endFreq
    t_windowName
    n_adcSpan
    t_measType
)
    => o_spectrumWaveform / g_value / nil
```

# **Description**

Calculates Signal-to-Noise-and-Distortion Ratio (SINAD), Spurious Free Dynamic Range (SFDR), Effective Number of Bits (ENOB), and Signal-to-Noise Ratio (without distortion) by using discrete fourier transform of any given input signal.

The spectrum measure is used for characterizing A-to-D converters and is typically supported for transient simulation data.

# Predefined and Waveform (Calculator) Functions

# Arguments

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.).
n_from	The X-axis start value of the portion of input o_waveform to be used for FFT and subsequent calculations.
n_to	The X-axis end value of the portion of input $o\_waveform$ to be used for FFT and subsequent calculations.
$x\_numSamples$	Optional number of sampled points used for the FFT.
	Valid values: Any integer power of two greater than zero.
	Default value: Number of data points in the signal o_waveform.
x_noiseBins	Optional number of noise bins, where the size of one bin is the reciprocal of the data window width. For example, 1 ms of transient data creates a bin size of 1 kHz.
	Valid values: Any integer power of two greater than or equal to zero.
	Default value: 0, implying that no signal is spilling into the bins.
	A frequency band of bin-size times the number of bins is calculated and adjusted as a function of the selected window. Frequency components in each band to the left and right of the fundamental or the harmonics are set to zero and do not contribute to any output result.
n_startFreq	Optional lower limit of frequency range for the spectrum measures.
	Default value: First frequency point of the FFT.
$n\_endFreq$	Optional upper limit of frequency range for the spectrum measures.
	Default value: Last frequency point of the FFT.
t_windowName	Optional windowing function applied to o_waveform.
	Valid values: Blackman, Cosine2, Cosine4, ExtCosBell, HalfCycleSine, HalfCycleSine3, HalfCycleSine6, Hamming, Kaiser, Parzen, Rectangular, and Triangular.
	Default value: Rectangular.

## Predefined and Waveform (Calculator) Functions

n adcSpan Optional full-scale span, ignoring any DC offsets. This is used in

ENOB calculation. Valid values: Any floating point number.

Default value: If  $n\_adcSpan$  is not specified or is nil, it is assumed to be 0 and is taken to be the peak-to-peak value of

the fundamental.

t measType Result specifier.

Valid values: sinad, sfdr (db), enob, and snhr.

### Value Returned

 $o\_spectrum \verb|Wavefor| Returns a waveform of spectrum measures.$ 

m

g value Returns the spectrum measure specified by the t measType

argument.

nil Returns nil and an error message otherwise.

# Example

```
spectrumMeas( VT("/vcoOut") 1K nil 1K 10G "Rectangular" nil "snhr") => -4.948
```

Returns the value of the spectrum measure snhr, as specified by the spectrumMeas function.

## Predefined and Waveform (Calculator) Functions

# **spectrumMeasurement**

```
spectrumMeasurement(
    o_waveform
    g_isTimeWave
    n_from
    n_to
    x_numSamples
    n_startFreq
    n_endFreq
    x_signalBins
    t_windowName
    n_satLvl
    g_isNoiseAnalysis
    x_noOfHarmonics
    t_measType
)
    => g value / nil
```

# **Description**

Calculates Signal-to-Noise-and-Distortion Ratio (SINAD), Spurious Free Dynamic Range (SFDR), Effective Number of Bits (ENOB), and Signal-to-Noise Ratio (without distortion) by using Fast Fourier Transform (FFT) of any given input signal.

The spectrum measure is used for characterizing A-to-D converters and is typically supported for transient simulation data.

# Predefined and Waveform (Calculator) Functions

# **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
g_isTimeWave	Boolean that specifies whether the input wave type is time domain waveform or frequency domain waveform.
n_from	The X-axis start value of the portion of input o_waveform to be used for FFT and subsequent calculations.
n_to	The X-axis end value of the portion of input $o\_waveform$ to be used for FFT and subsequent calculations.
$x\_numSamples$	Number of sampled points used for the FFT.
	Valid values: Any integer power of two greater than zero. For a value that is not a power of two, the function rounds it up to the next closest power of two.
	Default value: Number of data points in the Signal.
n_startFreq	Lower limit of frequency range for the spectrum measures.
	Default value: First frequency point of the FFT.
n_endFreq	Upper limit of frequency range for the spectrum measures.
	Default value: Last frequency point of the FFT.
$x\_signalBins$	Number of signal bins. When you select a window type, this field displays the default number of bins for the selected window type.
	For example, if you select the <i>Window Type</i> as Kaiser that has two signal bins, this field displays 2. You can increase the number of signal bins to up to half the value of the sample count.
	For example, if the sample count is 16 for the window type Kaiser, you can increase the signal bin count in the Signal Bins field up to 8. You cannot decrease the displayed signal bin

Default value: 0.

value. Valid values: 0 to 99.

### Predefined and Waveform (Calculator) Functions

t\_windowName Windowing function applied to o\_wave while applying the FFT

for measurement calculations.

Valid values: Blackman, Cosine2, Cosine4, ExtCosBell, HalfCycleSine, HalfCycleSine3, HalfCycleSine6, Hanning, Hamming, Kaiser, Parzen, Rectangular, and

Triangular.

Default value: Rectangular.

n satLv1 Peak saturation level of the FFT waveform. Magnitude of the

FFT wave is divided by the Peak Sat Level before using it in calculations. Peak sat level is the full-scale span ignoring any DC offsets and used in ENOB calculation. Valid values: Any

floating point number.

Default value: 0

g isNoiseAnalysis Boolean that specifies whether the analysis type is Signal

Analysis Or Noise Analysis.

 $x_noOfHarmonics$  Number of harmonics for the waveform that you want to plot.

For example, If this variable is n, where n should be greater than 1 and the fundamental frequency is harmonic 1, the n harmonics are considered for the harmonic power calculation. The signal bins are used for calculating the harmonic power.

For example, to calculate the total harmonic distortion (THD), if you set the Harmonics value to n, where n is greater than 1, and the fundamental frequency is harmonic 1, the number of harmonics used to calculate THD is 2,...,n. If n=3, the 2nd and

3rd harmonics are used to calculate THD.

t measType Result specifier.

Valid values: sinad, sfdr (db), enob, and snhr.

Default value: sinad

# Predefined and Waveform (Calculator) Functions

#### Value Returned

g_value	Returns the spectrum measure specified by the t_measType argument.
nil	Returns nil and an error message otherwise.

## **Example**

```
spectrumMeasurement(v("/OUT" ?result "tran") t 0 3e-08 1024 0 1.25e+08 0 "Rectangular" 0 0 1 "sinad") => -0.07218201
```

Returns the value of the spectrum measure sinad, as specified by the spectrumMeasurement function.

#### **Additional Information**

When you send the computed measurement values from the Spectrum toolbox to ADE Outputs and create an expression for them using ADE, the <code>spectrumMeasurement</code> function is used in the expression. For more information about <u>Spectrum</u> toolbox, see Spectrum in *Virtuoso Visualization and Analysis XL User Guide*.

The spectrumMeas function uses the same algorithm to calculate measurement values as that of the spectrumMeasurement SKILL function. The following table displays the mapping in the arguments for spectrumMeas and spectrumMeasurement functions:

spectrumMeas	spectrumMea- surement	Description
waveform	waveform	Specifies the waveform object.
NA	isTimeWave	This argument is available only in spectrum-Measurement function. The value of this argument is nil if the waveform sweep vector is of frequency domain, and the value is t if it is of time domain.  In spectrumMeas function, internally the unit of X-Vector is checked for Hz to know whether it is frequency domain or not.
from	from	The X-axis start value of the portion of input <code>o_waveform</code> to be used for FFT and subsequent calculations.

# Predefined and Waveform (Calculator) Functions

spectrumMeas	spectrumMea- surement	Description
to	to	The X-axis end value of the portion of input $o\_waveform$ to be used for FFT and subsequent calculations.
numSamples	numSamples	Number of sampled points used for the FFT. Valid values: Any integer power of two greater than zero. Default value: Number of data points in the Signal.
noiseBins	signalBins	In spectrumMeas, Number of Noise bins is the number of noise bins where the size of one bin is the reciprocal of the data window width. For example, 1 ms of transient data creates a bin size of 1 kHz.  Valid values: Any integer power of two greater than or equal to zero.  Default value: 0, implying that no signal is spilling into the bins
		In spectrumMeasurement, signalBins specifies the number of signal bins. When you select a window type, this field displays the default number of bins for the selected window type.  Default value: 0 to indicate the rectangular window type.
startFreq	startFreq	Lower limit of frequency range for the spectrum measures. Default value: First frequency point of the FFT.
endFreq	endFreq	Upper limit of frequency range for the spectrum measures. Default value: Last frequency point of the FFT.
windowName	windowName	Windowing function applied to o_wave while applying the FFT for measurement calculations. Valid values: Blackman, Cosine2, Cosine4, ExtCosBell, HalfCycleSine, HalfCycleSine3, HalfCycleSine6, Hanning, Hamming, Kaiser, Parzen, Rectangular, and Triangular.  Default value: Rectangular

# Predefined and Waveform (Calculator) Functions

spectrumMeas	spectrumMea- surement	Description
adcSpan	satLvl	In spectrumMeas, ADC Span is the full-scale span ignoring any DC offsets. This is used in ENOB calculation. Valid values: Any floating point number.
		In spectrumMeasurement, <code>satLv1</code> specifies the peak saturation level of the FFT waveform. Magnitude of the FFT wave is divided by the Peak Sat Level before using it in calculations. Peak sat level is the full-scale span ignoring any DC offsets and used in ENOB calculation. Valid values: Any floating point number.
NA	isNoiseAnaly- sis	This argument is present only in the spectrum-Measurement function. It specifies whether the analysis type is Noise Analysis.
NA	noOfHarmonics	This argument is available only in spectrum- Measurement function. This specifies the num- ber of harmonics for the waveform that you want to plot. Default value: 1
measType	measType	Result specifier. This argument is common for both the functions, but includes the following differences:
		■ sfdr(db) of spectrumMeas is same as sfdr of spectrumMeasurement or Spectrum assistant
		■ snhr of spectrumMeas is same as snr of spectrumMeasurement or Spectrum assistant.
		■ spectrumMeas supports the following measurements—sinad, sfdr(db), v, enob, thd. However, spectrumMeasurement supports more measurements in addition to the measurements supported by spectrumMeas.

# Predefined and Waveform (Calculator) Functions

# ssb

```
ssb(
    o_s11
    o_s12
    o_s21
    o_s22
    g_frequency
)
    => o_waveform / nil
```

# **Description**

Computes the source stability circles.

## Predefined and Waveform (Calculator) Functions

## **Arguments**

o_s11	Waveform object representing s11.
o_s12	Waveform object representing s12.
o_s21	Waveform object representing s21.
o_s22	Waveform object representing s22.

g\_frequency Frequency. It can be specified as a scalar or a linear range. The

frequency is swept if it is specified as a linear range. The linear range is specified as a list with three values: the start of the range, the end of the range, and the increment. For example, list (100M 1G 100M) specifies a linear range with the

following values:

```
{ 100M, 200M, 300M, 400M, 500M, 600M, 700M, 800M, 900M, 1G }
```

In that case, a source stability circle is calculated at each one of the 10 frequencies.

#### Value Returned

o waveform Waveform object representing the source stability circles.

nil Returns nil and an error message otherwise.

#### **Example**

plot(ssb(s11 s12 s21 s22 list(800M 1G 100M)))

## Predefined and Waveform (Calculator) Functions

## stddev

```
stddev(
    o_waveform
    [ ?overall overall ]
    )
    => n stddev / o waveformStddev / nil
```

# **Description**

Computes the standard deviation of a waveform (or a family of waveforms) over its entire range. Standard deviation (stddev) is defined as the square-root of the variance where variance is the integral of the square of the difference of the expression f(x) from average (f(x)), divided by the range of x.

For example, if y=f(x)

$$stddev(y) = \sqrt{\frac{from}{\int (y - average(y))^2}}$$

## Predefined and Waveform (Calculator) Functions

## **Arguments**

o\_waveform Waveform object or family of waveforms representing simulation

results that can be displayed as a series of points. (A waveform

object identifier looks like this: srrWave:XXXXX)

?overall

#### Value Returned

n stddev Returns a number representing the standard deviation value of

the input waveform.

o waveformStddev Returns a waveform representing the average value if the input

is a family of waveforms.

nil Returns nil or an error message.

# **Example**

```
stddev( v( "/net9" ) )
```

Gets the standard deviation of the voltage (Y-axis value) of /net9 over the entire time range specified in the simulation analysis.

# Predefined and Waveform (Calculator) Functions

# tangent

```
tangent(
    o_waveform
    [ ?x n_x ]
    [ ?y n_y ]
    [ ?slope n_slope ]
    [ ?ckm ckm ]
    )
    => o_waveform / nil
```

# **Description**

Returns the tangent to a waveform through the point  $(n \times, n \times)$  with the given slope.

# **Arguments**

o_waveform	Waveform object representing the wave.
?x n_x	X coordinate of the point. The default value is the X coordinate of the first point on the wave.
?y n_y	Y coordinate of the point. The default value is the Y coordinate at the given or default X coordinate.
?slope n_slope	Slope of the line. Default value: 1.0
?ckm ckm	

#### Value Returned

o_waveform	Wave object representing the line.
nil	Returns nil if there is an error.

# **Example**

```
refLine
=> tangent(refWave ?x -25 ?slope 1.0)
```

## Predefined and Waveform (Calculator) Functions

## thd

```
thd(
    o_waveform
    n_from
    n_to
    x_num
    n_fund
)
=> o_waveform / n_thdValue / nil
```

## **Description**

The thd function computes the percentage of total harmonic content of a signal with respect to the fundamental frequency expressed as a voltage percentage.

The computation uses the <u>dft</u> function. Assume that the *dft* function returns complex coefficients  $A_0$ ,  $A_1$ ...,  $A_f$ , .... Please note that fundamental frequency f is the frequency contributing to the largest power in the signal.  $A_0$  is the complex coefficient for the DC component and  $A_i$  is the complex coefficient for the ith harmonic where  $i \neq \theta, f$ . Then, total harmonic distortion is computed as:

$$\frac{\sqrt{\sum_{i=1, i \neq 0, f} |A_i|^2}}{|A_f|} \times 100 \%$$

## Predefined and Waveform (Calculator) Functions

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
n_from	Starting time for the DFT sample window.
n_to	Ending time for the DFT sample window.
x_num	Number of timepoints.
n_fund	Fundamental Frequency of the signal. If it is nil or zero then the non-zero frequency contributing to the largest power in the signal is used as the fundamental frequency. Otherwise, the harmonic frequency nearest to its value is used as the fundamental frequency.

#### Value Returned

o_waveform	Returns a waveform representing the absolute value of the total harmonic distortion if the input argument is a family of waveforms.
n_thdValue	Returns the absolute value of the total harmonic distortion of the input waveform.
nil	Returns nil and an error message otherwise.

#### Example

```
plot( thd( v( "/net8" ) 10u 20m 64 0 ) )
```

Computes the absolute value of the total harmonic distortion for the waveform representing the voltage of "/net8". The computation is done from 10u to 20m with 64 time points using the non-zero frequency contributing to the largest power in the signal as the fundamental frequency. The resulting waveform is plotted.

```
plot( thd( v( "/net8" ) 10u 20m 64 90 ) )
```

Computes the absolute value of the total harmonic distortion for the waveform representing the voltage of "/net8". The computation is done from 10u to 20m with 64 timepoints using a harmonic frequency, whose absolute difference w.r.t 90 is minimum, as the fundamental frequency. The resulting waveform is plotted.

# Predefined and Waveform (Calculator) Functions

# thd\_fd

```
thd_fd(
    t_name
    t_result
)
=> n thdValue / nil
```

# **Description**

The thd\_fd function returns the total harmonic distortion of the input waveform.

# **Arguments**

t_name	Name of the node for which total harmonic distorted is to be computed.
t_result	Name of the result of the specified node.

#### **Value Returned**

n_thdValue	Return a value of total harmonic distortion of the input waveform.
nil	Returns nil and an error message otherwise.

# **Example**

```
thd_fd( "Plo" ?result "FOURO-tran.test_fourier" )
```

Computes the total harmonic distortion for the "FOUR0-tran.test\_fourier" dataset in the results, where Fourier is connected to node "Plo".

# Predefined and Waveform (Calculator) Functions

# unityGainFreq

```
unityGainFreq(
    o_gainFreqWaveform
)
=> n frequency / nil
```

# **Description**

Computes and reports the frequency at which the gain is unity.

# **Arguments**

```
o\_gainFreqWavefor Gain frequency waveform.
```

## **Value Returned**

n_frequency	Returns a scalar value representing the frequency at which the gain of the input waveform is unity.
nil	Returns nil otherwise.

# **Example**

```
unityGainFrequency( VF("/out") )
```

# Predefined and Waveform (Calculator) Functions

# value

```
value(
    o_waveform
    [ ?scale scale ]
    [ ?period n_period ]
    [ ?xName s_xName]
    [ ?histoDisplay g_histoDisplay ]
    [ ?noOfHistoBins x_noOfHistoBins ]
    @Rest args
)
    => o_waveform / g_value / nil
```

# **Description**

Returns the Y value of a waveform for a given X value.

# Predefined and Waveform (Calculator) Functions

# **Arguments**

Arguments	
o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
?scale <i>scale</i>	Specifies the interpolation scale for WREAL signals. Valid values: linear, log, nil. Note that all scale elements after the second element are ignored.
	X is interpolated in linear mode if scale=nil or the first element of scale is linear.
	X is interpolated in log mode if the first element of scale is log.
	Y is interpolated in linear mode if scale has a single element linear or second element is linear.
	Y is interpolated in log mode if scale has a single element log, or second element is log
s_name	The name of the innermost or outermost sweep variable. If the sweep variable name is not supplied, the innermost sweep variable is used.
g_value	Value (X value) at which to provide the Y value. If a string has been defined for a value or set of values, the string may be used instead of the value.
?period <i>n_period</i>	The interval or period after which the value needs to be computed.
g_multiple	An optional boolean argument that takes the value nil by default. If set to t, the function returns multiple occurrences of the interpolated value.
?xName s_xName	An optional argument that is used only when $g_{multiple}$ is set to t. It takes the value time by default. It controls the contents of the x vector of the waveform object returned by the function.
	Valid values: time, cycle

?histoDisplay  $g_histoDisplay$ 

## Predefined and Waveform (Calculator) Functions

When set to t, returns a waveform that represents the statistical distribution of the riseTime data in the form of a histogram. The height of the bars (bins) in the histogram represents the frequency of the occurrence of values within the range of riseTime data.

Valid values: t nil
Default value: nil

?noOfHistoBins x\_noOfHistoBins

Denotes the number of bins represented in the histogram representation.

Valid values: Any positive integer

Default value: 1

@Rest args

**Note:**  $g_histoDisplay$  and  $x_noOfHistoBins$  are added for backward compatibility only. It will be deprecated in future releases. Use the histo function for plotting the histogram of the resulting function.

For the simplest calls to the function, which specify only the given waveform ( $o\_waveform$ ) and the X value ( $g\_value$ ), the given waveform can be a family of waveforms. If the family is of dimension m,  $g\_value$  can be either of dimension m-1 or a scalar. If  $g\_value$  is scalar, the function returns the Y value of all the components of the family at the specified  $g\_value$ .

## Predefined and Waveform (Calculator) Functions

#### Value Returned

o_waveform	Returns a waveform or a family of waveforms if the input argument is a family of waveforms or if values are expected at multiple points.
g_value	Returns the Y value if the input argument is a single waveform.
	For parametric sweeps, the value might be a waveform that can be printed with the ocnPrint command.
nil	Returns nil and an error message if the value cannot be printed.

## **Example**

```
value( v( "/net18" ) 4.428e-05 )
```

Prints the value of "/net18" at time=4.428e-05. This is a parametric sweep of temperature over time.

```
value( v( "/OUT" )'TEMPDC 20.0 )
```

Returns srrWave: XXXXX, indicating that the result is a waveform.

```
print( value( v( "/OUT" )'TEMPDC 20.0 ) )
```

Prints the value of v ( "/OUT" ) at every time point for TEMPDC=20.

```
print( value( v( "/OUT" ) 200n ?period 100n) )
```

Prints the value of v ( "/OUT" ) at 200n, 300n and so on at intervals of 100n until the end of the waveform.

```
value(VT("/out") 2e-07 ?period 2e-07 ?xName "time") (V)
```

Returns multiple occurrences of the value specified against time-points at which each interpolated value occurs.

```
value(VT("/out") 2e-07 ?period 2e-07 ?xName "cycle") (V)
```

Returns multiple occurrences of value specified against cycle numbers, where a cycle number refers to the n'th occurrence of the value event in the input waveform.

## Predefined and Waveform (Calculator) Functions

#### **xmax**

```
xmax(
    o_waveform
    x_numberOfPeaks
)
=> o waveform / g value / l value / nil
```

# **Description**

Computes the value of the independent variable (X) at which the Y value attains its maximum value.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be
	displayed as a series of points on a grid. (A waveform object
	identifier looks like this: srrWave: XXXXX.)
x numberOfPeaks	Specifies the <i>n</i> th X value corresponding to the maximum Y

value. For example, if  $x_numberOfPeaks$  is 3, the X value corresponding to the third maximum Y value is returned. If you specify a negative integer for  $x_numberOfPeaks$ , the X values are counted from right to left (from maximum to minimum). If  $x_numberOfPeaks$  is 0, xmax returns a list of X

locations.

#### Value Returned

o_waveform	Returns a waveform (or a family of waveforms) if the input argument is a family of waveforms.
g_value	Returns the X value corresponding to the peak specified with $x\_numberOfPeaks$ if the input argument is a single waveform.
l_value	Returns a list of X locations when $x\_numberOfPeaks$ is 0 and the input argument is a single waveform.
nil	Returns nil and an error message otherwise.

# Example

```
xmax(v("/net9")1)
```

# Predefined and Waveform (Calculator) Functions

Gets the time value (X-axis value) at which the voltage of "/net9" attains its first peak value.

Gets the list of time values (X-axis values) at which the voltage of "/net9" attains each of its peak values.

# Predefined and Waveform (Calculator) Functions

# xmin

```
xmin(
    o_waveform
    x_numberOfValleys
)
=> o_waveform / g_value / l_value / nil
```

# **Description**

Computes the value of the independent variable (X) at which the Y value attains its minimum value.

# **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
x_numberOfValleys	Specifies the $n$ th X value corresponding to the minimum Y value. For example, if $x\_numberOfValleys$ is 3, the X value corresponding to the third minimum Y value is returned. If you specify a negative integer for $x\_numberOfValleys$ , the X-values are counted from right to left (from maximum to minimum). If $x\_numberOfValleys$ is 0, xmin returns a list of X locations.

#### Value Returned

o_waveform	Returns a waveform (or a family of waveforms) if the input argument is a family of waveforms.
g_value	Returns the X value corresponding to the valley specified with $x\_numberOfValleys$ if the input argument is a single waveform.
l_value	Returns a list of X locations when $x\_numberOfValleys$ is 0 and the input argument is a single waveform.
nil	Returns nil and an error message otherwise.

# Example

```
xmin( v( "/net9" ) 1 )
```

# Predefined and Waveform (Calculator) Functions

Gets the time value (X axis) at which the voltage of "/net9" has its first low point or valley. xmin(v("/net9")0)

Gets the list of time values (X axis) at which the voltage of "/net9" has low points or valleys.

## Predefined and Waveform (Calculator) Functions

## xval

```
xval(
    o_waveform
)
=> o_waveform / nil
```

# **Description**

Returns a waveform whose X vector and Y vector are equal to the input waveform's X vector.

## **Arguments**

o_waveform	Waveform object representing sir	mulation results that can be
	displayed as a series of points or	n a grid. (A waveform object

identifier looks like this: srrWave: XXXXX.)

#### **Value Returned**

o_waveform	Returns a waveform if the input argument is a single waveform.
	Returns a family of waveforms if the input argument is a family
	of waveforms.

nil Returns nil and an error message otherwise.

## **Example**

```
xval( v( "/net8" ))
```

Returns a waveform in which the X vector for the voltage of "/net8" is also used for the Y vector.

## Predefined and Waveform (Calculator) Functions

# ymax

```
ymax(
    o_waveform
    [ ?overall overall ]
    )
    => n max / o waveformMax / nil
```

# **Description**

Computes the maximum value of the waveform's Y vector.

A waveform consists of an independent-variable X vector and a corresponding Y vector.

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
?overall overall	

# Value Returned

n_max	Returns a number representing the maximum value of Y if the input argument is a single waveform.
o_waveformMax	Returns a waveform (or family of waveforms) representing the maximum value of Y if the input argument is a family of waveforms.
nil	Returns nil and an error message otherwise.

# **Example**

```
ymax( v( "/net9" ) )
```

Gets the maximum voltage (Y value) of "/net9".

# Predefined and Waveform (Calculator) Functions

# ymin

```
ymin(
    o_waveform
    [?overall overall]
)
=> n min / o waveformMin / nil
```

# **Description**

Computes the minimum value of a waveform's Y vector.

(A waveform consists of an independent-variable X vector and a corresponding Y vector.)

## **Arguments**

o_waveform	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave: XXXXX.)
?overall overall	

#### Value Returned

n_min	Returns a number representing the minimum value of Y if the input argument is a single waveform.
o_waveformMin	Returns a waveform (or family of waveforms) representing the minimum value of Y if the input argument is a family of waveforms.
nil	Returns nil and an error message otherwise.

# **Example**

```
ymin( v( "/net9" ) )
```

Gets the minimum voltage (Y value) of "/net9".

# Predefined and Waveform (Calculator) Functions

# **Spectre RF Calculator Functions**

This section describes the following calculator functions used for Spectre RF data analysis:

- ifreq
- **■** <u>ih</u>
- <u>itime</u>
- pir
- pmNoise
- pn
- <u>pvi</u>
- <u>pvr</u>
- spm
- totalNoise
- vfreq
- <u>vfreqterm</u>
- <u>vh</u>
- vhterm
- vtime
- **■** <u>vtimeterm</u>
- <u>ypm</u>
- zpm

## Predefined and Waveform (Calculator) Functions

# ifreq

```
ifreq(
    s_ana
    t_terminal
    [ freq n_freq ]
    )
    => o waveform / nil
```

# Description

Returns the current of the terminal at a specified frequency or at all frequencies in the frequency domain.

## **Arguments**

s_ana	Analysis type or analysis name. The available analyses are hb,
	pss, qpss, pac, hbac, qpac, <b>and</b> ac.

Default value: hb

t\_terminal Terminal name on the schematic or signal name from the

Results Browser.

n freq Frequency for which you want to plot the results. It is an optional

field.

Valid values: Any integer or floating point number.

Default value: nil

When you specify nil, current on all the frequency points are

returned.

#### Value Returned

o\_waveform Returns a waveform representing current at a specified

frequency or at all frequency points.

nil Returns nil and an error message otherwise.

## **Example**

```
ifreq("hb" "/load/PLUS" 50 )
```

# Predefined and Waveform (Calculator) Functions

Returns the current for /load/PLUS signal, which is obtained from hb analysis, at frequency=50.

## Predefined and Waveform (Calculator) Functions

## ih

```
ih(
    s_ana
    t_terminal
    [ harmonic x_hlist ]
)
=> o waveform / nil
```

## **Description**

Returns the current of the terminal at a specified harmonic or at all harmonics in the frequency domain.

## **Arguments**

	s ana	Analysis type or analysis name. The available analyses are	e hb,
--	-------	--	-------

pss, qpss, pac, hbac, and qpac.

Default value: hb

t\_terminal Terminal name on the schematic or signal name from the

Results Browser.

x hlist Harmonics for which you want to plot the results. It is an

optional field. For analyses, such as hb, pss, pac, and hbac, you can add either single harmonic value or an available list of

harmonic values in this field.

Valid values: Any integer or a list from the available list of harmonics. You can find the available harmonics by using the

harmonicList function.

Default value: nil.

#### Value Returned

o waveform Returns a waveform representing current at a specified

harmonic or at all harmonic points.

nil Returns nil and an error message otherwise.

# **Example**

```
ih("hb" "/rf/PLUS" 2 )
```

# Predefined and Waveform (Calculator) Functions

Returns the current for /rf/PLUS signal, which is obtained from hb analysis, at harmonic= 2.

## Predefined and Waveform (Calculator) Functions

# itime

```
itime(
    s_ana
    t_terminal
    [ time n_time ]
)
    => o_waveform / nil
```

# **Description**

Returns the current of the terminal at a specified time point or at all time points in the time domain.

## **Arguments**

s_ana	Analysis type or analysis name. The available analyses are ${\tt hb}$ , ${\tt pss}$ , and ${\tt tran}$ .
	Default value: hb
t_terminal	Terminal name on the schematic or signal name from the Results Browser.
n_time	Time points for which you want to plot the results. If you specify a time point in this field, the result of the specified time is returned. It is an optional field.
	Valid values: Any integer or floating point number.
	Default value: nil.

#### Value Returned

o_waveform	Returns a waveform representing current at a specified time point or at all time points.
nil	Returns nil and an error message otherwise.

# Example

```
itime("hb" "/load/PLUS" 4 )
```

Returns the current for /load/PLUS signal, which is obtained from hb analysis, at time=4s.

# Predefined and Waveform (Calculator) Functions

# pir

```
pir(
    s_ana
    t_branch1
    t_branch2
    n_resistance
    [ harmonic x_hlist ]
    )
    => o_waveform / nil
```

# **Description**

Returns the spectral power from current and resistance for a specified harmonic list or for all harmonic points.

## Predefined and Waveform (Calculator) Functions

## **Arguments**

s ana Analysis type or analysis name. The available analyses are hb,

pss, qpss, pac, hbac and qpac.

Default value: hb

t branch1 First branch name on the schematic or signal name from the

Results Browser.

t branch2 Second branch name on the schematic or signal name from the

Results Browser.

n resistance The resistance value.

Valid values: Any integer or floating point number.

harmonic x\_hlist

Harmonics for which you want to plot the results. It is an optional field. For analyses, such as hb, pss, pac, and hbac, you can add either single harmonic value or an available list of

harmonic values in this field.

Valid values: Any integer or a list from the available list of harmonics. You can find the available harmonics by using the

harmonicList function.

Default value: nil.

#### **Value Returned**

o waveform Returns a waveform representing spectral power from current

and resistance for a specified harmonic list.

nil Returns nil and an error message otherwise.

#### **Example**

```
pir("hb" "/V1/PLUS" "/rf/PLUS" 2 5 )
```

This example returns the spectral power for /V1/PLUS and /rf/PLUS, which are obtained from the hb analysis, at resistance=2 ohms and harmonic=5.

# Predefined and Waveform (Calculator) Functions

# pmNoise

```
pmNoise(
    s_ana
    [ freq n_freq ]
    s_modifier
    g_dsb
)
    => o_waveform / n_pnoise / nil
```

# **Description**

Returns the modulated phase noise at a specified frequency or for the entire spectrum.

## Predefined and Waveform (Calculator) Functions

## **Arguments**

s ana Analysis type or analysis name.

Valid values: pnoise, and hbnoise.

Default value: pnoise

n freq Frequency for which you want to calculate the modulated phase

noise.

Valid values: Any integer or floating point number

Default value: nil, which means the frequency at all points are

calculated.

s modifier Modifier to be used.

Valid values: dBc, normalized, Power, Magnitude, and dBV

Default value: dBc.

g\_dsb Specifies whether you want to include the double side band.

Valid values: t and nil

Default value: t

#### **Value Returned**

n pnoise Returns the modulated phase noise at the specified frequency

point.

o waveform Returns a waveform representing the modulated phase noise at

all frequency points.

nil Returns nil and an error message otherwise.

#### **Example**

```
pmNoise("hbnoise" 50 "dBc" t )
```

This example returns the modulated phase noise for hbnoise analysis at frequency=50 and modifier=dBc and double side bands included.

## Predefined and Waveform (Calculator) Functions

#### pn

# **Description**

Returns the phase noise at a specified frequency or at all frequency points.

## **Arguments**

s	ana	Analysis type or analysis name.

Valid values: pnoise, hbnoise, and qpnoise.

Default value: pnoise

 $n\_freq$  Frequency for which you want to calculate the phase noise.

Valid values: Any integer or floating point number

Default value: nil, which means the frequency at all points are

calculated.

#### Value Returned

n_pn	Returns the phase noise at a specified frequency point.
o_waveform	Returns a waveform representing the phase noise at all frequency points.
nil	Returns nil and an error message otherwise.

#### Example

```
pn("hbnoise" 50 )
```

This example returns the phase noise for hbnoise analysis at frequency=50.

# Predefined and Waveform (Calculator) Functions

# pvi

```
pvi(
    s_ana
    t_pos
    t_neg
    t_branch1
    t_branch2
    [ harmonic x_hlist ]
)
    => o waveform / nil
```

# **Description**

Returns the spectral power from voltage and current for a specified harmonic list or for all harmonics.

# Predefined and Waveform (Calculator) Functions

# **Arguments**

s_ana	Analysis type or analysis name. The available analyses are hb, pss, qpss, pac, hbac and qpac.
	Default value: hb
t_pos	Positive node or net from the schematic or from the Results Browser. This field can also contain an explicit voltage value.
t_neg	Negative node or net from the schematic or from the Results Browser. This field can also contain an explicit voltage value.
t_branch1	First branch name on the schematic or signal name from the Results Browser.
t_branch2	Second branch name on the schematic or signal name from the Results Browser.
$x_hlist$	Harmonics for which you want to plot the results. It is an optional field. For analyses, such as hb, pss, pac, and hbac, you can add either single harmonic value or available list of harmonic values in this field.
	Valid values: Any integer or a list from the available list of harmonics. You can find the available harmonics by using the <u>harmonicList</u> function.

#### **Value Returned**

o_waveform	Returns a waveform representing the spectral power from voltage and current for a specified harmonic list or for all harmonics.
nil	Returns nil and an error message otherwise.

Default value: nil

# **Example**

```
pvi("hb" "/RFin" "/RFout" "/V1/PLUS" "/V2/PLUS" 2)
```

This example returns the spectral power for the following values:

- Analysis Type is hb
- Positive node is /RFin
- Negative node is /RFout

# Predefined and Waveform (Calculator) Functions

- Branch name 1 /V1/PLUS
- Branch name 2 /V2/PLUS
- Harmonic List is 2

### Predefined and Waveform (Calculator) Functions

### pvr

```
pvr(
    s_ana
    t_pos
    t_neg
    n_resistance
    [ harmonic x_hlist ]
    )
    => o_waveform / nil
```

# **Description**

Returns the spectral power at a specified harmonic list or at all harmonics with resistor and voltage on the positive and negative nodes.

#### Predefined and Waveform (Calculator) Functions

### **Arguments**

s_ana	Analysis type or analysis name.	The available analyses are hb,
-------	---------------------------------	--------------------------------

pss, qpss, pac, hbac and qpac.

Default value: hb

t pos Positive node or net from the schematic or from the Results

Browser. This field can also contain an explicit voltage value.

*t\_neg* Negative node or net from the schematic or from the Results

Browser. This field can also contain an explicit voltage value.

n resistance The resistance value.

Valid values: Any integer or floating point number

 $x_hlist$  Specify the harmonics for which you want to plot the results. It is

an optional field. For analyses, such as hb, pss, pac, and hbac, you can add either single harmonic value or available list

of harmonic values in this field.

Valid values: Any integer or a list from the available list of harmonics. You can find the available harmonics by using the

harmonicList function.

Default value: nil.

#### Value Returned

o waveform Returns a waveform representing the spectral power on

specified harmonic list or on all harmonics with resistor and

voltage on the positive and negative nodes

nil Returns nil and an error message otherwise.

#### Example

```
pvr("hb" "/RFin" "/RFout" 2 2 )
```

This example returns the spectral power for the following values:

- Analysis Type is hb
- Positive node is /RFin
- Negative node is /RFout
- Resistance is 2

# Predefined and Waveform (Calculator) Functions

■ Harmonic List is 2

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# Predefined and Waveform (Calculator) Functions

# spm

```
spm(
    s_ana
    x_index1
    x_index2
    [ ?port1 x_port1 ]
    [ ?port2 x_port2 ]
    )
    => o_waveform / nil
```

# **Description**

Returns the waveform for s-parameters.

### Predefined and Waveform (Calculator) Functions

#### **Arguments**

 $s_ana$  Analysis type or analysis name.

Valid values: sp, psp, qpsp, and hbsp

Default value: sp

x index1 Port index for sp simulation. By default, this field is set to blank.

Valid values: Available port index, such as 1, 2.

x index1 Port index for sp simulation. By default, this field is set to blank.

Valid values: Available port index, such as 1, 2.

?port1 x\_port1

Port instance. The port instance can be specified only for the differential s-parameter analysis and not applicable for psp,

qpsp and hbsp analyses.

Valid values: Predefined values "c" and "d" for Spectre

simulator.

?port2 x port2

Port instance. The port instance can be specified only for the differential s-parameter analysis and not applicable for psp,

gpsp and hbsp analyses.

Valid values: Predefined values "c" and "d" for Spectre

simulator.

#### Value Returned

o waveform Returns a waveform representing the s-parameters.

nil Returns nil and an error message otherwise.

#### Example

spm("sp" 1 1 ?port1 nil ?port2 nil)

This example plots the s-parameter waveform for sp analysis with index1=1 and index 2=1.

### Predefined and Waveform (Calculator) Functions

### totalNoise

```
totalNoise(
    s_ana
    n_sfreq
    n_efreq
    [ instances l_instances ]
)
    => n_totalNoise / nil
```

### **Description**

Returns the total noise in a specified frequency limit.

### **Arguments**

s_ana	Analysis type or analysis name. The available analyses are noise, pnoise, qpnoise, and hbnoise.
	Default value: noise.
n_sfreq	The start frequency.
	Valid values: Any integer or floating point number
n_efreq	The end frequency.
	Valid values: Any integer or floating point number
l_instances	List of instances or instance names. The noise contributed by the instances specified in this field is ignored while calculating the total noise. This is an optional field.

#### Value Returned

$n\_$ totalNoise	Returns the total noise in a specified frequency limit.
nil	Returns nil and an error message otherwise.

### Example

```
totalNoise("hbnoise" 1k 100k out )
```

This example returns the total noise for hbnoise analysis in the frequency range 1k to 100k with instance out being excluded.

### Predefined and Waveform (Calculator) Functions

# vfreq

# **Description**

Returns the voltage of a net at a specified frequency or at all frequencies in the frequency domain.

### **Arguments**

s_ana	Analysis type or analysis name. The available analyses are hb, pss, qpss, pac, hbac, qpac, and ac.
	Default value: hb
t_net	Net name from the schematic or signal name from the Results Browser.
x_freq	Frequency for which you want to plot the results. It is an optional field.
	Valid values: Any integer value
	Default Value: nil

#### Value Returned

o_waveform	Returns a waveform representing the voltage of net at a specified frequency
nil	Returns nil and an error message otherwise

### Example

```
vfreq("hb" "/outp" 50 )
```

This example returns the voltage of /outp net from hb analysis at frequency=50.

# Predefined and Waveform (Calculator) Functions

# vfreqterm

```
vfreqterm(
    s_ana
    t_terminal
    [ freq x_freq ]
    )
    => o_waveform | g_value | nil
```

# **Description**

Returns the voltage of a terminal at a specified frequency or at all frequencies in the frequency domain.

### **Arguments**

s_ana	Analysis type or analysis name. The available analyses are ac, pss, hb, pac, hbac, qpss, and qpac.
	Default value: hb
t_terminal	Terminal name in schematic from the Results Browser.
freq x_freq	Frequency for which you want to plot the results. It is an optional field.
	Valid value: Any integer value
	Default value: nil

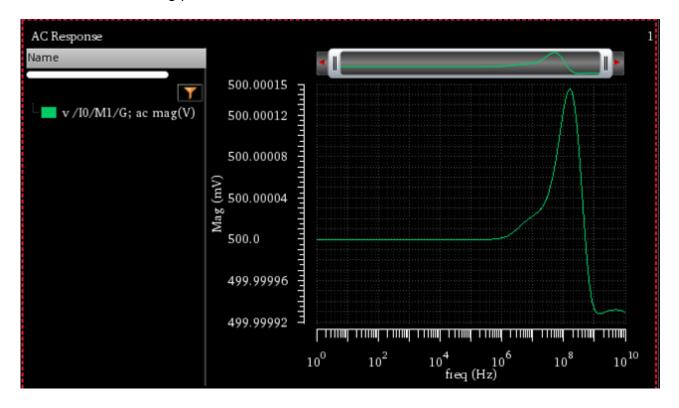
#### Value Returned

o_waveform	Returns a waveform representing the voltage at a terminal on the specified frequency.
g_value	Returns the value of voltage at a terminal on the specified frequency.
nil	Returns nil or an error message.

```
vfreqterm( 'ac "/IO/M1/G" )
```

# Predefined and Waveform (Calculator) Functions

It returns the following plot.



### Predefined and Waveform (Calculator) Functions

#### vh

```
vh(
     s_ana
     t_net
     [ harmonic x_hlist ]
)
     => o waveform / nil
```

### **Description**

Returns the voltage on a net at a specified harmonic or at all harmonics in the frequency domain.

### **Arguments**

s_ana I	٩na	lysis type oı	analysis name.	The available analyses are hb,
---------	-----	---------------	----------------	--------------------------------

pss, qpss, pac, hbac, and qpac.

Default value: hb

t net Net name on the schematic or signal name from the Results

Browser.

x hlist Harmonics for which you want to plot the results. It is an

optional field. For analyses, such as hb, pss, pac, and hbac, you can add either single harmonic value or available list of

harmonic values in this field.

Valid values: Any integer or a list from the available list of harmonics. You can find the available harmonics by using the

harmonicList function.

Default value: nil

#### Value Returned

o\_waveform Returns a waveform representing the voltage on a net on the

specified harmonic

nil Returns nil and an error message otherwise

```
vh("hb" "/outp" 5 )
```

# Predefined and Waveform (Calculator) Functions

This example returns the voltage of	/outp	net from ha	o <b>analysis</b>	at harmonic=5.
-------------------------------------	-------	-------------	-------------------	----------------

#### Predefined and Waveform (Calculator) Functions

#### vhterm

```
vhterm(
    s ana
    t_terminal
     [ harmonic x hlist ]
    => o waveform | g value | nil
```

### **Description**

Returns the voltage on a terminal at the specified harmonic or at all the harmonics in the frequency domain.

### **Arguments**

s ana Analysis type or analysis name. The availab
---

analyses are pss, hb, pac, hbac, qpss, and qpac.

Default value: hb

Terminal name in schematic from the Results t terminal

Browser.

Harmonics for which you want to plot the results. It is harmonic x hlist

> an optional field. For analyses, such as hb, pss, pac, and hbac, you can add either single harmonic value

or available list of harmonic values in this field.

Valid values: Any integer or a list from the available

list of harmonics. You can find the available harmonics by using the harmonic list function.

Default value: nil

#### Value Returned

0	waveform	Rε	eturns a	a wave	torm r	epresent	ting t	the vol	tage at	: a

terminal on the specified harmonic.

Returns the value of voltage at a terminal on the g value

specified harmonic.

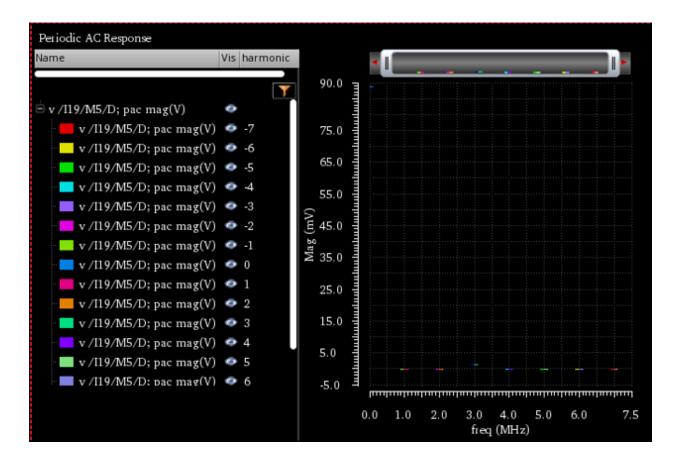
nil Returns nil or an error message.

# Predefined and Waveform (Calculator) Functions

### **Example**

vhterm( 'pac "/I19/M5/D" )

It returns the following plot.



### Predefined and Waveform (Calculator) Functions

## vtime

```
vtime(
    s_ana
    t_net
    [ time n_time ]
)
    => o_waveform / nil
```

# **Description**

Returns the voltage of a net at a specified time point or at all time points in the time domain.

# **Arguments**

s_ana	Analysis type or analysis name. The available analyses are hb, pss, and tran.
	Default value: hb
t_net	Net name from the schematic or signal name from the Results Browser.
n_time	Time points for which you want to plot the results. If you specify a time point in this field, the result of the specified time is returned. Otherwise, It is an optional field.
	Valid values: Any integer or floating point number.

#### Value Returned

o_waveform	Returns a waveform representing the voltage of net at a specified time point
nil	Returns nil and an error message otherwise

Default value: nil

### **Example**

```
vtime("hb" "/outm" 20)
```

This example returns the voltage of /outp net from hb analysis at time=20s.

### Predefined and Waveform (Calculator) Functions

### vtimeterm

```
vtimeterm(
    s_ana
    t_terminal
    [ time n_time ]
    )
    => o_waveform | g_value | nil
```

# **Description**

Returns the voltage of a terminal at a specified time point or at all time points in the time domain.

### **Arguments**

s_ana	Analysis type or analysis name. The available analyses are tran, pss, and hb.
	Default value: hb
t_terminal	Terminal name in schematic from the Results Browser.
time n_time	Time points for which you want to plot the results. If you specify a time point in this field, the result of the specified time point is returned. Otherwise, it is an optional field.
	Valid value: Any integer or floating point number
	Default value: nil

#### Value Returned

o_waveform	Returns a waveform representing the voltage at a terminal on the specified time point.
g_value	Returns the value of voltage at a terminal on the specified time point.
nil	Returns nil or an error message.

### **OCEAN Reference** Predefined and Waveform (Calculator) Functions

# **Example**

vtimeterm( 'tran "/IO/M1/D" )

It returns the following plot.



### Predefined and Waveform (Calculator) Functions

### ypm

```
ypm(
    s_ana
    x_index1
    x_index2
)
    => o_waveform / nil
```

# **Description**

Returns the waveform for y-parameters.

# **Arguments**

s_ana	Analysis type or analysis name. Valid values: sp, psp, qpsp, and hbsp
	Default value: sp
x_index1	Port index for sp simulation. By default, this field is set to blank.
	Valid values: Available port index, such as 1, 2
x_index1	Port index for sp simulation. By default, this field is set to blank.
	Valid values: Available port index, such as 1, 2

### Value Returned

o_waveform	Returns a waveform representing the y-parameters
nil	Returns nil and an error message otherwise

# Example

```
ypm("sp" 1 1)
```

This example returns the waveform for y-parameters when index1=1 and index2=1.

### Predefined and Waveform (Calculator) Functions

# zpm

# **Description**

Returns the waveform for z-parameters.

# **Arguments**

s_ana	Analysis type or analysis name.
	Valid values: sp, psp, qpsp, and hbsp
	Default value: sp
x_index1	Port index for sp simulation. By default, this field is set to blank.
	Valid values: Available port index, such as 1, 2
x_index1	Port index for sp simulation. By default, this field is set to blank.
	Valid values: Available port index, such as 1, 2

### **Value Returned**

o_waveform	Returns a waveform representing the z-parameters
nil	Returns nil and an error message otherwise

# Example

```
zpm("sp" 1 1)
```

This example returns the waveform for z-parameters when index1=1 and index2=1.

### Predefined and Waveform (Calculator) Functions

# **RF Functions**

This section describes the OCEAN commands for the following RF functions:

- B1f
- gac\_freq
- gac gain
- Gmax
- Gmin
- Gmsg
- GP
- gpc\_freq
- gpc\_gain
- GT
- Gmux
- Kf
- <u>loadStability</u>
- nc\_freq
- <u>nc\_gain</u>
- NF
- NFmin
- **■** <u>rn</u>
- sourceStability
- <u>s11</u>
- <u>s12</u>
- <u>s21</u>
- <u>s22</u>

### Predefined and Waveform (Calculator) Functions

### B<sub>1</sub>f

```
Blf(
     [ dataDir t_dataDir ]
   )
     => o waveform / nil
```

### **Description**

Returns the alternative stability factor in terms of the specified parameters.

### **Arguments**

dataDir t\_dataDir Results directory path.

### **Values Returned**

o\_waveform Waveform object representing the alternative stability factor.

nil Returns nil and an error message otherwise.

### Predefined and Waveform (Calculator) Functions

# gac\_freq

```
gac_freq(
    n_gain
    n_startFreq
    n_stopFreq
    n_step
    [ ?resultsDir t_resultsDir ]
    )
    => o_waveform / nil
```

### **Description**

Returns the available power gain circles where the gain is fixed and frequency is swept.

# **Arguments**

n_gain	Gain value in dB.
n_startFreq	Starting frequency.
n_stopFreq	Ending frequency.
n_step	Frequency step size to be used.
?resultsDir t_resultsDir	Results directory path.

### **Values Returned**

o\_waveformnilReturns nil and an error message otherwise.

```
gac freq(16 2G 3G 100M)
```

# Predefined and Waveform (Calculator) Functions

# gac\_gain

```
gac_gain(
    n_freq
    n_startGain
    n_stopGain
    n_step
    [ ?resultsDir t_resultsDir ]
    )
    => o_waveform / nil
```

### **Description**

Returns the available power gain circles where the frequency is fixed and gain is swept.

# **Arguments**

n_freq	Frequency in Hz.
n_startGain	Starting gain value.
n_stopGain	End gain value.
n_step	Gain step size to be used.
?resultsDir t_resultsDir	Results directory path.

#### **Values Returned**

o_waveform	Waveform object.
nil	Returns nil and an error message otherwise.

```
gac_gain(2.4G 14 18 0.5)
```

# Predefined and Waveform (Calculator) Functions

### **Gmax**

```
Gmax(
      [ dataDir t_dataDir ]
   )
      => o waveform / nil
```

### **Description**

Returns the maximum available gain for a two port.

### **Arguments**

dataDir t\_dataDir Results directory path.

### **Values Returned**

o\_waveform Waveform object.

### Predefined and Waveform (Calculator) Functions

# **Gmin**

```
Gmin(
     [ dataDir t_dataDir ]
   )
     => o waveform / nil
```

### **Description**

Returns the optimum noise reflection coefficient for NFmin.

# **Arguments**

dataDir t\_dataDir Results directory path.

### **Values Returned**

o\_waveform Waveform object.

# Predefined and Waveform (Calculator) Functions

# **Gmsg**

```
Gmsg(
     [ dataDir t_dataDir ]
)
=> o waveform / nil
```

### **Description**

Returns the maximum stable power gain for a two port.

### **Arguments**

dataDir t dataDir Results directory path.

### **Values Returned**

o\_waveform Waveform object.

### Predefined and Waveform (Calculator) Functions

### **GP**

```
GP(
     [ dataDir t_dataDir ]
)
=> o waveform / nil
```

### **Description**

Returns the power gain. Operating power gain, GP, is defined as the ratio between the power delivered to the load and the power input to the network.

### **Arguments**

dataDir t\_dataDir Results directory path.

### **Values Returned**

o\_waveform Waveform object representing the power gain.

nil Returns nil and an error message otherwise.

### Predefined and Waveform (Calculator) Functions

# gpc\_freq

### **Description**

Returns the operating power gain circles where the gain is fixed and frequency is swept.

### **Arguments**

n_gain	Gain value in dB.
n_startFreq	Starting frequency.
$n\_stopFreq$	Ending frequency.
n_step	Frequency step size to be used.
?resultsDir t resultsDir	Results directory path.

### **Values Returned**

o\_waveformnilReturns nil and an error message otherwise.

```
gpc freq(16 2.4G 2.5G 10M)
```

# Predefined and Waveform (Calculator) Functions

# gpc\_gain

### **Description**

Returns the operating power gain circles where the frequency is fixed and gain is swept.

# **Arguments**

n_freq	Frequency value in Hz.
n_start	Starting gain value.
n_stop	Ending gain value.
n_step	Gain step size to be used.
?resultsDir t_resultsDir	Results directory path.

#### **Values Returned**

o_waveform	Waveform object.
nil	Returns nil and an error message otherwise.

```
gpc_gain(2.4G 14 18 0.5)
```

### Predefined and Waveform (Calculator) Functions

### **GT**

```
GT(
     [ dataDir t_dataDir ]
)
=> o waveform / nil
```

### **Description**

Returns the transducer gain. Transducer power gain, GT, is defined as the ratio between the power delivered to the load and the power available from the source.

### **Arguments**

dataDir t dataDir Results directory path.

### **Values Returned**

o\_waveform Waveform object.

### Predefined and Waveform (Calculator) Functions

### **Gmux**

```
Gmux(
      [ dataDir t_dataDir ]
   )
      => o waveform / nil
```

### **Description**

Returns the maximum unilateral power gain for a two port.

Maximum unilateral transducer power gain, Gumx, is the transducer power gain when S12 is zero, and the source and load impedances conjugate are matching.

### **Arguments**

dataDir t dataDir Results directory path.

#### **Values Returned**

o\_waveform Waveform object.

### Predefined and Waveform (Calculator) Functions

### Kf

```
Kf(
     [ dataDir t_dataDir ]
)
=> o waveform / nil
```

### **Description**

Returns the Stern stability factor.

# **Arguments**

dataDir t\_dataDir Results directory path.

### **Values Returned**

o\_waveform Waveform object.

### Predefined and Waveform (Calculator) Functions

# loadStability

```
loadStability(
    n_startFreq
    n_stopFreq
    n_step
    [ ?resultsDir x_resultsDir ]
    )
    => o_waveform / nil
```

# **Description**

Computes the load stability circles.

### **Arguments**

n_startFreq	Start of the frequency range.
n_StopFreq	End of the frequency range.
n_step	Frequency step size to be used.
?resultsDir x resultsDir	Results directory path.

#### Value Returned

o_waveform	Waveform object representing the load stability circles.
nil	Returns nil and an error message otherwise.

```
loadStability(2G 3G 0.2G)
loadStability(2G 3G 0.2G?resultsDir "./psf" )
```

# Predefined and Waveform (Calculator) Functions

# nc\_freq

### **Description**

Returns the noise circles with fixed gain and swept frequency.

# **Arguments**

$n\_noiseLevel$	Noise level.
n_startFreq	Starting frequency.
n_stopFreq	Ending frequency.
n_step	Frequency step size to be used.
?resultsDir t resultsDir	Results directory path.
c	

### **Values Returned**

o_waveform	waveform object.
nil	Returns nil and an error message otherwise.

```
nc_freq(2 2G 3G 0.2G)
```

### Predefined and Waveform (Calculator) Functions

# nc\_gain

# **Description**

Returns the noise circles with fixed frequency and swept noise level.

# **Arguments**

n_freq	Fixed frequency.
n_startNoiseLvl	Starting noise level.
$n\_stopNoiseLvl$	Ending noise level.
n_step	Sweeping noise level step size
?resultsDir t_resultsDir	Results directory path.

#### **Values Returned**

o_waveform	Waveform object.
nil	Returns nil and an error message otherwise.

```
nc_gain(2.4G 1.5 2.4 0.2)
```

### Predefined and Waveform (Calculator) Functions

### NF

```
NF(
      [ dataDir t_dataDir ]
   )
=> o_waveform / nil
```

## **Description**

Returns the noise figure.

# **Arguments**

dataDir t\_dataDir Results directory path.

### **Values Returned**

o\_waveform Waveform object.

# Predefined and Waveform (Calculator) Functions

# **NFmin**

```
NFmin(
      [ dataDir t_dataDir ]
)
=> o waveform / nil
```

# **Description**

Returns the minimum noise figure.

# **Arguments**

dataDir t\_dataDir Results directory path.

#### **Values Returned**

o\_waveform Waveform object.

nil Returns nil and an error message otherwise.

# Predefined and Waveform (Calculator) Functions

#### rn

```
rn(
    [ dataDir t_dataDir ]
)
=> o waveform / nil
```

# **Description**

Returns the normalized equivalent noise resistance as a function of frequency.

# **Arguments**

dataDir t\_dataDir Results directory path.

#### **Values Returned**

o\_waveformnilReturns nil and an error message otherwise.

# Predefined and Waveform (Calculator) Functions

# sourceStability

```
sourceStability(
    n_startFreq
    n_stopFreq
    n_step
    [ ?resultsDir x_resultsDir ]
    )
    => o_waveform / nil
```

# **Description**

Computes the source stability circles.

# **Arguments**

n_startFreq	Start of the frequency range.
n_StopFreq	End of the frequency range.
n_step	Frequency step size to be used.
?resultsDir x resultsDir	Results directory path.

#### Value Returned

o_waveform	Waveform object representing the source stability circles.
nil	Returns nil and an error message otherwise.

# **Example**

```
sourceStability(2G 3G 0.2G)
sourceStability(2G 3G 0.2G?resultsDir "./psf" )
```

# Predefined and Waveform (Calculator) Functions

# s11

```
s11(
     [ ?resultsDir t_resultsDir ]
   )
     => o waveform / nil
```

# **Description**

Returns the response at port 1 due to a signal at port 1.

# **Arguments**

?resultsDir Results directory path.
t resultsDir

#### **Values Returned**

o\_waveform Waveform object

nil Returns nil and an error message otherwise.

# Predefined and Waveform (Calculator) Functions

# s12

```
s12(
     [ ?resultsDir t_resultsDir ]
   )
     => o waveform / nil
```

# **Description**

Returns the response at port 1 due to a signal at port 2.

# **Arguments**

?resultsDir Results directory path. t resultsDir

#### **Values Returned**

o\_waveformnilReturns nil and an error message otherwise.

# Predefined and Waveform (Calculator) Functions

# **s21**

```
s11(
     [ ?resultsDir t_resultsDir ]
   )
     => o waveform / nil
```

# **Description**

Returns the response at port 2 due to a signal at port 1.

# **Arguments**

?resultsDir Results directory path.
t resultsDir

#### **Values Returned**

o\_waveform Waveform object

nil Returns nil and an error message otherwise.

# Predefined and Waveform (Calculator) Functions

# **s22**

```
s22(
     [ ?resultsDir t_resultsDir ]
   )
     => o waveform / nil
```

# **Description**

Returns the response at port 2 due to a signal at port 2.

# **Arguments**

?resultsDir Results directory path. t resultsDir

#### **Values Returned**

o\_waveformnilReturns nil and an error message otherwise.

Predefined and Waveform (Calculator) Functions

11

# **Parametric Analysis Commands**

These commands set up a parametric analysis. When you run a parametric analysis, you can plot the resulting data as a family of curves.

This chapter contains information on the following commands:

- paramAnalysis
- paramRun

# Parametric Analysis Commands

# paramAnalysis

```
paramAnalysis(
     t desVar
     [ ?start n_start ]
     [ ?stop n stop ]
     [ ?center n center ]
     [ ?span n_span ]
     [ ?step f step ]
     [ ?lin n_lin ]
     [ ?log n log ]
     [ ?dec n dec ]
     [ ?oct n oct ]
     [ ?times n times ]
     [ ?spanPercent n spanPercent ]
     [ ?sweepType t_sweepType ]
     [ ?values 1 values ]
     [ o paramAnalysis ]
     => undefined / nil
```

# **Description**

Sets up a parametric analysis.

Groups the PSF data so that it can be plotted as a family of curves when the analysis is finished. The commands can be nested as shown in the syntax of the command.

If you specify more than one range, the OCEAN environment uses the following precedence to select a single range to use.

Similarly, if you specify more than one step control, the OCEAN environment uses the following precedence.

f step highest precedence

# Parametric Analysis Commands



To run the analysis, use the paramRun command described in "paramRun" on page 663.

# Parametric Analysis Commands

# **Arguments**

t_desVar	Name of the design variable to be swept.
?start n_start	Beginning value for the design variable.
?stop n_stop	Final value for the design variable.
?center n_center	
	Center point for a range of values that you want to sweep.
?span n_span	Range of values that you want to sweep around the center point. For example, if $n\_center$ is 100 and $n\_span$ is 20 then the sweep range extends from 90 to 110.
?step f_step	Increment by which the value of the design variable changes. For example, if $n\_start$ is 1.0, $n\_stop$ is 2.1, and $f\_step$ is 0.2, the parametric analyzer simulates at values 1.0, 1.2, 1.4, 1.6, 1.8, and 2.0.
?lin <i>n_lin</i>	The number of steps in the analysis. The parametric analyzer automatically assigns equal intervals between the steps. With this option, there is always a simulation at both $n\_start$ and $n\_stop$ . The value for the $n\_lin$ argument must be an integer greater than 0.
	For example, if $n\_start$ is 0.5, $n\_stop$ is 2.0, and $n\_lin$ is 4, the parametric analyzer simulates at values 0.5, 1.0, 1.5, and 2.0.
?log <i>n_log</i>	The number of steps between the starting and stopping points at equal-ratio intervals using the following formula:
	log multiplier = $(n_stop/n_start)^{(n_log-1)}$ The number of steps can be any positive number, such as 0.5, 2, or 6.25.

Default value: 5

For example, if  $n\_start$  is 3,  $n\_stop$  is 15, and  $n\_log$  is 5, the parametric analyzer simulates at values 3, 4.48605, 6.7082, 10.0311, and 15.

The ratios of consecutive values are equal, as shown below. 3/4.48605 = 4.48605/6.7082 = 6.7082/10.0311 = 10.0311/15 = .67

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# Parametric Analysis Commands

?dec  $n_dec$ 

The number of steps between the starting and stopping points calculated using the following formula:

decade multiplier =  $10^{1/n}$ \_dec

The number of steps can be any positive number, such as 0.5, 2, or 6.25.

Default value: 5

For example, if  $n\_start$  is 1,  $n\_stop$  is 10, and  $n\_dec$  is 5, the parametric analyzer simulates at values 1, 1.58489, 2.51189, 3.98107, 6.30957, and 10.

The values are  $10^0$ ,  $10^{.2}$ ,  $10^{.4}$ ,  $10^{.6}$ ,  $10^{.8}$ , and  $10^1$ .

?oct n oct

The number of steps between the starting and stopping points using the following formula:

The number of steps can be any positive number, such as 0.5, 2, or 6.25.

Default value: 5

For example, if  $n\_start$  is 2,  $n\_stop$  is 4, and  $n\_oct$  is 5, the parametric analyzer simulates at values 2, 2.2974, 2.63902, 3.03143, 3.4822, and 4.

These values are  $2^1$ ,  $2^{1.2}$ ,  $2^{1.4}$ ,  $2^{1.6}$ ,  $2^{1.8}$ , and  $2^2$ .

octave?multiplier =  $2^{1/(n\_oct)}$ 

?times  $n_{times}$ 

A multiplier. The parametric analyzer simulates at the points between  $n\_start$  and  $n\_stop$  that are consecutive multiples of  $n\_times$ .

For example, if  $n\_start$  is 1,  $n\_stop$  is 1000, and  $n\_times$  is 2, the parametric analyzer simulates at values 1, 2, 4, 8, 16, 32, 64, 128, 256, and 512.

?spanPercent  $n_spanPercent$ 

Range specified as a percentage of the center value. For example, if  $n\_center$  is 100 and  $n\_spanPercent$  is 40, the sweep range extends from 80 to 120.

# Parametric Analysis Commands

?sweepType t sweepType

Type of parametric analysis. Valid values are:

- paramset Runs Parametric Set analysis, specific to Spectre.
- nil Runs Sweeps & Ranges type parametric analysis.

Default value: nil

?values 1 values

List of values to be swept. You can use  $1\_values$  by itself or in conjunction with  $n\_start$ ,  $n\_stop$ , and  $f\_step$  to specify the set of values to sweep.

?paramAnalysis o paramAnalysis

Value returned from another paramAnalysis call used to achieve multidimensional parametric analysis.

#### Value Returned

undefined

The return value for this command is undefined.

nil

Returns nil and prints an error message if there are problems setting the option.

#### **Example**

Sets up a parametric analysis for the rs design variable. The swept values are 200, 400, 600, 800, 1000, 1030, 1050, and 1090.

Sets up a nested parametric analysis for the rl design variable.

```
paramAnalysis("temp" ?start -50 ?stop 100 ?step 50)
```

Sets up a parametric analysis for temperature.

#### Parametric Analysis Commands

# paramRun

```
paramRun(
     [ s_paramAnalysis ]
     )
     => t / nil
paramRun (
     [ ?jobName t jobName ]
     [ ?drmsCmd t_drmsCmd ]
     => s jobName / nil
paramRun(
     [ ?jobName t jobName ]
     [ ?host t hostName ]
     [ ?queue t queueName ]
     [ ?startTime t startTime ]
     [ ?termTime t termTime ]
     [ ?dependentOn t dependentOn ]
     [ ?mail t mailingList ]
     [ ?block s block ]
     [ ?notify s notifyFlag ]
     [ ?lsfResourceStr s lsfResourceStr ]
     => s jobName / nil
```

# Description

Runs the specified parametric analysis.

If you do not specify a parametric analysis, all specified analyses are run. Distributed processing must be enabled using the hostmode command before parametric analyses can be run in distributed mode.

When the paramRun command finishes, the PSF directory contains a file named runObjFile that points to a family of data. To plot the family, use a normal plot command. For example, you might use plot (v("/out")).

For information about specifying a parametric analysis, see the paramAnalysis command described in <u>"paramAnalysis"</u> on page 658.

# Parametric Analysis Commands

# Arguments

s_paramAnalysis	Parametric analysis.
?jobName t_jobName	Used as the basis of the job name. The value entered for $t\_jobName$ is used as the job name and return value if the run command is successful. If the name given is not unique, a number is appended to create a unique job name.
?host t_hostName	Name of the host on which to run the analysis. If no host is specified, the system assigns the analysis to an available host.
?drmsCmd t_drmsCmd	A DRMS (Distributed Resource Management System) command, such as a bsub command for LSF or a qsub command for SGE (Sun Grid Engine) used to submit a job. When this argument is used, all other arguments, except ?jobName will be ignored. Moreover, it will not be possible to call the OCEAN function wait on the jobs submitted using this argument.
	To know more about the command option, refer to the section Submitting a Job in the chapter <u>Using the Distributed</u> <u>Processing Option in the Analog Design Environment</u> of the Virtuoso Analog Distributed Processing OptionUser Guide.
?queue t_queueName	Name of the queue. If no queue is defined, the analysis is placed in the default queue (your home machine).
?startTime t_startTime	Desired start time for the job. If dependencies are specified, the job does not start until all dependencies are satisfied.
<pre>?termTime t_termTime</pre>	Termination time for job. If the job is not completed by $t\_termTime$ , the job is terminated.
?dependentOn t_dependentOn	List of jobs on which the specified analysis is dependent. The analysis is not started until after dependent jobs are complete.
?mail t_mailingList	List of users to be notified by e-mail when the analysis is complete.
?block s_block	When $s\_block$ is not nil, the OCEAN script halts until the job is complete.
	Default value: nil
?notify s_notifyFlag	When $notifyFlag$ is not nil, a job completion message is echoed to the OCEAN interactive window.
	Default value: t

#### Parametric Analysis Commands

?lsfResouceStr Specifies an LSF Resource Requirement string to submit a job.  $s\_1sfResourceStr$  It is effective only in the LSF mode.

#### Value Returned

t Returned if successful.

nil Returns nil and prints an error message if unsuccessful.

# **Example**

```
paramRun() => t
```

#### Runs all specified parametric analyses.

```
rsAnalysis = paramAnalysis("CAP" ?values '(10 20))
paramRun('rsAnalysis)
```

#### OR

```
rsAnalysis = paramAnalysis("CAP" ?values '(10 20) paramAnalysis("RES" ?values '(10 20 )))
paramRun('rsAnalysis)
```

#### Runs the rs parametric analysis.

```
paramRun(?queue "background" ?lsfResourceStr "mem>500")
```

Runs the analysis in the queue named background on a machine, if it has at least 500 MB of RAM memory.

# Parametric Analysis Commands

# OCEAN Distributed Processing Commands

The Open Command Environment for Analysis (OCEAN) distributed processing commands let you run OCEAN jobs across a collection of computer systems.

This chapter contains information on the following commands:

- deleteJob
- digitalHostMode
- <u>digitalHostName</u>
- hostMode
- hostName
- killJob
- monitor
- remoteDir
- resumeJob
- suspendJob
- wait

This chapter also provides sample OCEAN scripts that optimally use these commands. See the section <u>Sample Scripts</u>.

For detailed information on distributed processing, refer to <u>Virtuoso Analog Distributed Processing Option User Guide</u>.

# **OCEAN Distributed Processing Commands**

# deleteJob

```
deleteJob(
    t_jobName
    [ t_jobName2 t_jobName3 ... t_jobNameN ]
    )
    => t / nil
```

# **Description**

Removes a job or series of jobs from the text-based job monitor.

Deleted jobs are no longer listed in the job monitor. The deleteJob command applies only to ended jobs.

# **Arguments**

t_jobName	Name used to identify the job.
t_jobname2t_jobn	Additional jobs that you want to delete.
ameN	

#### Value Returned

t	Returns t if successful.
nil	Returns nil and prints an error message if unsuccessful.

# Example

```
deleteJob( 'myckt)
=> t
```

Deletes the myckt job.

# **OCEAN Distributed Processing Commands**

# digitalHostMode

```
digitalHostMode(
     { 'local | 'remote }
)
     => t / nil
```

#### **Description**

For mixed-signal simulation, specifies whether the digital simulator will run locally or on a remote host.

#### **Arguments**

'local	Sets the simulation to run locally on the user's machine.
--------	---

'remote Sets the simulation to run on a remote host. If you use this

argument, you must specify the host name by using the

digitalHostName command.

#### Value Returned

t Returns t if successful.

nil Returns nil and prints an error message if unsuccessful.

#### **Example**

```
digitalHostMode( 'local )
```

Sets the digital simulator to run locally on the user's host.

# **OCEAN Distributed Processing Commands**

# digitalHostName

```
digitalHostName(
    t_name
)
    => t / nil
```

# **Description**

For mixed-signal simulation, specifies the name of the remote host for the digital simulator.

When you use the digitalHostMode ('remote) command, use this command to specify the name of the remote host.

# **Arguments**

t name Name used to identify the host for the digital simulator.

#### Value Returned

t Returns t if successful.

nil Returns nil and prints an error message if unsuccessful.

# **Example**

```
digitalHostName( "digitalhost" )
```

Indicates that the digital simulator runs on the host called digitalhost.

# OCEAN Distributed Processing Commands

# hostMode

# **Description**

Sets the simulation host mode.

The default value for hostMode is specified in the asimenv.startup file with the hostMode environment variable.

# **Arguments**

'local	Sets the simulation to run locally on the user's machine.
'remote	Sets the simulation to run on a remote host queue. For this release, the remote host is specified in the .cdsenv file.
'distributed	Sets the simulation to run using the distributed processing software.

# **Value Returned**

t	Returns t if successful.
nil	Returns nil and prints an error message if unsuccessful.

# **Example**

```
hostMode( 'distributed )
=> t
```

Enables distributed processing on the current host.

# **OCEAN Distributed Processing Commands**

# hostName

# **Description**

Specifies the name of the remote host.

When you use the hostMode('remote) command, use this command to specify the name of the remote host.

# **Arguments**

*t\_name* Name used to identify the remote host.

#### Value Returned

t Returns t if successful.

nil Returns nil and prints an error message if unsuccessful.

# **Example**

```
hostName( "remotehost" )
```

Specifies that the host called remotehost is to be used for remote simulation.

# **OCEAN Distributed Processing Commands**

#### killJob

```
killJob(
    t_jobName [ t_jobName2 t_jobName3 ... t_jobNameN ]
   )
   => t / nil
```

# **Description**

Stops processing of a job or a series of jobs.

The job might still show up in the job monitor, but it cannot be restarted. Use the deleteJob command to remove the job name from the job server and job monitor.

# **Arguments**

```
t\_jobName Name used to identify the job. t\_jobname2...t\_jobnameN Additional jobs that you want to stop.
```

#### Value Returned

t	Returns t if successful.
nil	Returns nil and prints an error message if unsuccessful.

# **Example**

```
killJob( 'myckt )
=> t
```

Aborts the job called myckt. If the job is in the queue and has not started running yet, it is deleted from the queue.

# **OCEAN Distributed Processing Commands**

#### monitor

```
monitor(
     [ ?taskMode s_taskMode ]
)
     => t / nil
```

# **Description**

Monitors the jobs submitted to the distributed system.

#### **Arguments**

s\_taskMode When not nil, multitask jobs are expanded to show individual

jobs. A multitask job is one that contains several related jobs.

#### Value Returned

t Returns t if successful.

nil Returns nil and prints an error message if unsuccessful.

#### **Example**

```
monitor( ?taskMode t )
```

Displays the name, host, and queue for all pending tasks sorted on a queue name.

# **OCEAN Distributed Processing Commands**

#### remoteDir

```
remoteDir(
    t_path
)
=> t / nil
```

# **Description**

Specifies the project directory on the remote host to be used for remote simulation.

When you use the hostMode('remote) command, use this command to specify the project directory on the remote host.

# **Arguments**

t\_path Specifies the path to the project directory on the remote host to

be used for remote simulation.

#### Value Returned

t Returns t if successful.

nil Returns nil and prints an error message if unsuccessful.

#### **Example**

```
remoteDir( "~/simulation" )
```

Specifies that the project directory is  $\sim$ /simulation.

# OCEAN Distributed Processing Commands

#### resumeJob

# **Description**

Resumes the processing of a previously suspended job or series of jobs. The resumeJob command applies only to jobs that are suspended.

# **Arguments**

t_jobName	Name used to identify the job.
	Additional jobs that you want to resume.
ameN	

#### **Value Returned**

t	Returns t if successful.
nil	Returns nil and prints an error message if unsuccessful.

# **Example**

```
resumeJob( 'myckt )
=> t
```

Resumes the myckt job that was halted with the suspendJob command.

# **OCEAN Distributed Processing Commands**

# suspendJob

```
suspendJob(
    t_jobName [ t_jobName2 t_jobName3 ... t_jobNameN ]
   )
   => t / nil
```

# **Description**

Suspends the processing of a job or series of jobs. The suspendJob command applies only to jobs that are pending or running.

# **Arguments**

```
t\_jobName Name used to identify the job. 
 t\_jobName2...t\_jobnameN Additional jobs that you want to suspend.
```

#### Value Returned

t Returns t if successful.

nil Returns nil and prints an error message if unsuccessful.

# **Example**

```
suspendJob( 'myckt )
=> t
```

Suspends the job called myckt.

# **OCEAN Distributed Processing Commands**

#### wait

```
wait(
    [ ?queue t_queueName ]
    jobName [ jobName2 jobName3 ... jobNameN ]
)
    => t / nil
```

# Description

Postpones processing of a script until the specified jobs complete. This command is ignored if distributed processing is not available.

The wait command is highly useful when you use the non-blocking mode of distributed processing and you want to do some post-processing, such as selecting and viewing results after a job is completed. The wait command is not required when you use the blocking mode of distributed processing. To know more about blocking and non-blocking modes of DP, refer to <u>Virtuoso Analog Distributed Processing Option User Guide</u>.

# **Arguments**

?queue t queueName

The name of queue on which job launched by wait is

submitted.

t\_jobName Name used to identify the job. The job name is user defined or

system generated, depending on how the user submitted the

job.

 $t_{jobName2...}t_{jobnameN}$ 

Additional jobs that you want to postpone.

#### Value Returned

t Returns t if successful.

nil Returns nil and prints an error message if unsuccessful.

#### **Examples**

```
wait( 'myckt1 )
=> t
```

# **OCEAN Distributed Processing Commands**

Postpones execution of all subsequent OCEAN commands until the job myckt1 completes.

```
wait( ?queue "lnx64" 'job0 )
=> t
```

Job launched by wait is submitted on lnx64 queue that postpones the execution of all subsequent OCEAN commands until the job job0 completes.

# **OCEAN Distributed Processing Commands**

# **Sample Scripts**

This section provides sample scripts for the following:

- To submit multiple jobs and show the use of the dependentOn argument in one job
- To set up and run a simple analysis in blocking mode and select results
- To set up and run a parametric analysis in blocking mode and select results
- To submit multiple jobs without using wait or selecting results
- To submit multiple jobs using wait and selection of results

#### To submit multiple jobs and show the use of the dependentOn argument in one job

This script can be used to submit multiple jobs while using the dependent on argument in one of these jobs.

```
; set up the environment for the jobs
simulator( 'spectre )
hostMode( 'distributed )
design( "/home/simulation/test2/spectre/schematic/netlist/netlist")
resultsDir( "/home/simulation/test2/spectre/schematic" )
analysis('tran ?stop "5u" )
temp(27)
jobList = nil
; starting first job
jobList = append1( jobList run( ?queue "test" ?host "menaka" ) )
analysis ('tran ?stop "50u")
; starting second job
jobList = append1( jobList run(?jobName "job_2" ?queue "test" ?host "menaka"))
analysis('tran ?stop "10u")
; starting third job, which is dependent on job 2
jobList= append1(jobList run(?jobName "disable" ?queue "test" ?dependentOn
                symbolToString(car(last(jobList)))))
; wait for all the jobs to complete
```

# **OCEAN Distributed Processing Commands**

```
wait((append1 last(jobList) nil))
; open and plot the results of the jobs
openResults( car(last(jobList)))
selectResult( 'tran )
newWindow()
plot(getData("/net61") )

openResults( nth(1 jobList))
selectResult('tran)
newWindow()
plot(getData("/net61") )
```

#### To set up and run a simple analysis in blocking mode and select results

```
; set up the environment for Simple Analysis
simulator( 'spectre )
hostMode( 'distributed )
design (
"/home/amit/Artist446/simulation/ampTest/spectre/schematic/netlist/netlist" )
resultsDir( "/home/Artist446/simulation/ampTest/spectre/schematic" )
modelFile(
    '("/home/Artist446/Models/myModels.scs" "")
analysis('tran ?stop "3u")
desVar( "CAP" 0.8p )
temp(27)
; submit the job in blocking mode, to the queue test and machine menaka
run(?queue "test" ?host "menaka" ?block t)
; select and plot the results
selectResult( 'tran )
plot(getData("/out"))
```

# To set up and run a parametric analysis in blocking mode and select results

```
; set up the environment for parametric analysis.
simulator('spectre')
hostMode('distributed')
design(
```

#### **OCEAN Distributed Processing Commands**

```
"/home/amit/Artist446/simulation/ampTest/spectre/schematic/netlist")
resultsDir( "/home/amit/Artist446/simulation/ampTest/spectre/schematic"
)
modelFile(
        '("/home/amit/Artist446/Models/myModels.scs" "")
)
analysis('tran ?stop "3u" )
desVar( "CAP" 0.8p )
temp( 27 )
paramAnalysis("CAP" ?values '(1e-13 2.5e-13 4e-13 ))
; submit the job in blocking mode, to the queue test and machine menaka
paramRun(?queue "fast" ?host "menaka" ?block t)

; select and plot the results
selectResult( 'tran )
plot(getData("/out") )
```

# To submit multiple jobs without using wait or selecting results

```
; set up the environment for the jobs
simulator( 'spectre )
hostMode( 'distributed )
design (
"/home/Artist446/simulation/ampTest/spectre/schematic/netlist/netlist")
resultsDir( "/home/Artist446/simulation/ampTest/spectre/schematic" )
modelFile(
    '("/home/Artist446/Models/myModels.scs" "")
; setup and submit first job
analysis('tran ?stop "3u" )
desVar(
        "CAP" 0.8p
temp(27)
run(?queue "SUN5 5032" ?host "menaka")
; setup and submit second job
analysis('ac ?start "1M" ?stop "2M" )
analysis('tran ?stop "3u" )
desVar( "CAP" 0.8p )
temp(27)
```

#### **OCEAN Distributed Processing Commands**

```
run(?queue "SUN5_5032" ?host "menaka")
```

#### To submit multiple jobs using wait and selection of results

```
; set up the environment for the jobs
simulator( 'spectre )
hostMode( 'distributed )
design (
"/home/Artist446/simulation/ampTest/spectre/schematic/netlist/netlist")
resultsDir( "/home/Artist446/simulation/ampTest/spectre/schematic" )
modelFile(
    '("/home/Artist446/Models/myModels.scs" "")
; initialize jobList to nil
jobList = nil
; setup and submit first job
analysis('tran ?stop "3u" )
desVar(
        "CAP" 0.8p
temp(27)
jobList = append1( jobList run(?queue "SUN5 5032" ?host "menaka") )
; setup and submit second job
analysis('ac ?start "1M" ?stop "2M" )
analysis('tran ?stop "3u" )
desVar(
        "CAP" 0.8p
temp( 27 )
jobList = append1( jobList run(?queue "SUN5 5032" ?host "menaka"))
; wait for both the jobs to finish
wait( (append1 jobList nil) )
; open and plot the result of first job
openResults( (car jobList))
selectResult( 'tran )
plot(getData("/out") )
; open and plot the result of second job
openResults( (cadr jobList))
selectResult( 'tran )
```

# **OCEAN Distributed Processing Commands**

```
plot(getData("/out") )
selectResult( 'ac )
plot(getData("/out") )
; delete the jobs
foreach( x jobList deleteJob( x ) )
```

# **Language Constructs**

There are three types of SKILL language constructs:

Conditional statements

Conditional statements test for a condition and perform operations when that condition is found. These statements are if, unless, and when.

Selection statements

A selection statement allows a list of elements, each with a corresponding operation. A variable can then be compared to the list of elements. If the variable matches one of the elements, the corresponding operation is performed. These statements include for, foreach, and while.

Iterative statements

Iterative statements repeat an operation as long as a certain condition is met. These statements include case and cond.

This chapter contains information on the following statements

case if

cond unless

for when

foreach while

### Language Constructs

#### if

```
if(
    g_condition
    g_thenExpression
    [ g_elseExpression ]
)
    => g_result
    if(
    g_condition
    then g_thenExpr1 ...
    [ else g_elseExpr1 ... ]
)
    => g result
```

Evaluates  $g\_condition$ , typically a relational expression, and runs  $g\_thenExpression$  if the condition is true (that is, its value is non-nil); otherwise, runs  $g\_elseExpression$ .

The value returned by if is the value of the corresponding expression evaluated.

Additionally, if also can also be used with the keywords then and else to group sequences of expressions for conditional execution. If the  $g\_condition$  is true, the sequence of expressions between then and else (or the end of the if form) is evaluated, with the value of the last expression evaluated returned as the value of the form.

# **Arguments**

g_condition	Any Virtuoso® SKILL language expression
$g\_thenExpression$	Any SKILL expression
g_elseExpression	Any SKILL expression

#### Value Returned

 $g\_result$  Returns the value of  $g\_thenExpression$  if  $g\_condition$  has a non-nil value. The value of  $g\_elseExpression$  is returned otherwise.

### **Example**

```
x = 2
if (( x > 5) 1 0)
=> 0
```

### Language Constructs

Returns 0 because x is less than 5.

```
a ="npn"
if(( a == "npn" ) print( a ) ) "npn"
=> nil
```

Prints the string npn and returns the result of print.

```
x = 5
if( x "non-nil" "nil" )
=> "non-nil"
```

Returns "non-nil" because x was not nil. If x was nil, "nil" would be returned.

```
x = 7
if((x > 5) then 1 else 0)
=> 1
```

Returns 1 because x is greater than 5.

### Language Constructs

#### unless

```
unless(
    g_condition
    g_expr1 ...
)
=> g result / nil
```

# **Description**

Evaluates a condition. If the result is true (non-nil), it returns nil; otherwise it evaluates the body expressions in sequence and returns the value of the last expression.

The semantics of this function can be read literally as "unless the condition is true, evaluate the body expressions in sequence."

### **Arguments**

g_condition	Any SKILL expression
g_expr1	Any SKILL expression

#### Value Returned

g_result	Returns the value of the last expression of the sequence $g expr1$ if $g condition$ evaluates to nil.
nil	Returns nil if g_condition evaluates to non-nil.

# **Example**

```
x = -123 unless( x \ge 0 println( "x is negative" ) -x ) => 123 
Prints "x is negative" as a side effect.
```

```
unless( x < 0 println( "x is positive ") x) => nil
```

Returns nil.

# Language Constructs

### when

```
when(
    g_condition
    g_expr1 ...
)
=> g result / nil
```

# **Description**

Evaluates a condition.

If the result is non-nil, evaluates the sequence of expressions and returns the value of the last expression. Otherwise, returns nil.

# **Arguments**

g_condition	Any SKILL expression
g_expr1	Any SKILL expression

#### Value Returned

g_result	Returns the value of the last expression of the sequence $g_{expr1}$ if $g_{condition}$ evaluates to non-nil.
nil	Returns nil if the $g\_condition$ expression evaluates to nil.

### **Example**

```
x = -123
when(x < 0 println("x is negative") -x)
=> 123

Prints "x is negative" as a side effect.
when(x >= 0 println("x is positive") x)
=> nil
```

#### Returns nil.

### Language Constructs

# for

```
for(
    s_loopVar
    x_initialValue
    x_finalValue
    g_expr1
    [ g_expr2 ... ]
    )
    => t
```

### **Description**

Evaluates the sequence  $g_{expr1} g_{expr2} \dots$  for each loop variable value, beginning with  $x_{initialValue}$  and ending with  $x_{finalValue}$ .

First evaluates the initial and final values, which set the initial value and final limit for the local loop variable named  $s\_loopVar$ . Both  $x\_initialValue$  and  $x\_finalValue$  must be integer expressions. During each iteration, the sequence of expressions  $g\_exprl$   $g\_exprl$  ... is evaluated and the loop variable is then incremented by one. If the loop variable is still less than or equal to the final limit, another iteration is performed. The loop ends when the loop variable reaches a value greater than the limit. The loop variable must not be changed inside the loop. It is local to the for loop and would not retain any meaningful value upon exit from the for loop.

**Note:** Everything that can be done with a for loop can also be done with a while loop.

# Language Constructs

# **Arguments**

s_loopVar	Name of the local loop variable that must not be changed inside the loop.
x_initialValue	Integer expression setting the initial value for the local loop variable.
$x_finalValue$	Integer expression giving final limit value for the loop.
g_expr1	Expression to evaluate inside loop.
g_expr2	Additional expressions to evaluate inside loop.

#### Value Returned

t

This construct always returns t.

# **Example**

```
sum = 0
for( i 1 10
    sum = sum + i
    printf( "%d" sum ))
=> t
```

#### Prints 10 numbers and returns t.

```
sum = 0
for( i 1 5
    sum = sum + i
    println( sum )
    )
=> t
```

Prints the value of sum with a carriage return for each pass through the loop:

```
1
3
6
10
15
```

### foreach

# **Description**

Evaluates one or more expressions for each element of a list of values.

#### The first syntax form,

```
foreach( s_formalVar g_exprList g_expr1 [g_expr2 ...] )
=> l_valueList
```

evaluates  $g\_exprList$ , which returns a list  $l\_valueList$ . It then assigns the first element from  $l\_valueList$  to the formal variable  $s\_formalVar$  and processes the expressions  $g\_expr1$   $g\_expr2$  ... in sequence. The function then assigns the second element from  $l\_valueList$  and repeats the process until  $l\_valueList$  is exhausted.

### The second syntax form,

```
\label{local_condition} \begin{tabular}{ll} for each ( & s_formalVar1...s_formalVarN) & g_exprList1... & g_exprListN & g_expr1 & [g_expr2 & ...] & ) => 1 & valueList & \\ \end{tabular}
```

can iterate over multiple lists to perform vector operations. Instead of a single formal variable, the first argument is a list of formal variables followed by a corresponding number of expressions for value lists and the expressions to be evaluated.

### The third syntax form,

```
foreach( s_formalVar g_exprTable g_expr1 [g_expr2 ...])
=> o valueTable
```

### Language Constructs

can be used to process the elements of an association table. In this case,  $s\_formalVar$  is assigned each key of the association table one by one, and the body expressions are evaluated each iteration. The syntax for association table processing is provided in this syntax statement.

### **Arguments**

$s\_formalVar$	Name of the variable.
g_exprList	Expression whose value is a list of elements to assign to the formal variable $s\_formalVar$ .
g_expr1 g_expr2	Expressions to execute.
g_exprTable	Association table whose elements are to be processed.

#### Value Returned

$1\_valueList$	Returns the value of the second argument, $g_{exprList}$ .
o_valueTable	Returns the value of $g_{exprTable}$ .

#### **Example**

```
foreach( x '( 1 2 3 4 ) println( x ) )
1
2
3
4
=> ( 1 2 3 4 )
```

Prints the numbers 1 through 4 and returns the second argument to foreach.

```
foreach( key myTable printf( "%L : %L" key myTable[key] ) )
```

Accesses an association table and prints each key and its associated data.

```
( foreach ( x y ) '( 1 2 3 ) '( 4 5 6 ) ( println x+y ) )
5
7
9
=> ( 1 2 3 )
```

Uses foreach with more than one loop variable.

# Language Constructs

# **Errors and Warnings**

The error messages from foreach might at times appear cryptic because some foreach forms get expanded to call the mapping functions mapc, mapcar, mapcan, and so forth.

# Language Constructs

# while

```
while(
    g_condition
    g_expr1 ...
)
    => t
```

# **Description**

Repeatedly evaluates  $g\_condition$  and the sequence of expressions g expr1 ... if the condition is true.

This process is repeated until  $g\_condition$  evaluates to false (nil). As this form always returns t, it is principally used for its side effects.

**Note:** Everything that can be done with a for loop can also be done with a while loop.

# **Arguments**

g_condition	Any SKILL expression
g exprl	Any SKILL expression

#### Value Returned

t Always returns t.

# **Example**

```
i = 0
while( (i <= 10) printf("%d" i++) )
=> t.
```

Prints the digits 0 through 10.

### Language Constructs

#### case

```
case(
    g_selectionExpr
    l_clause1 [ l_clause2 ... ]
)
    => g result / nil
```

# **Description**

Evaluates the selection expression, matches the resulting selector values sequentially against comparators defined in clauses, and runs the expressions in the matching clause.

Each  $1\_clause$  is a list of the form ( $g\_comparator g\_expr1$  [ $g\_expr2...$ ]), where a comparator is either an atom (that is, a scalar) of any data type or a list of atoms. Comparators are always treated as constants and are never evaluated. The  $g\_selectionExpr$  expression is evaluated and the resulting selector value is matched sequentially against comparators defined in  $1\_clause1$   $1\_clause2...$ . A match occurs when either the selector is equal to the comparator or the selector is equal to one of the elements in the list given as the comparator. If a match is found, the expressions in that clause and that clause only (that is, the first match) are run. The value of case is then the value of the last expression evaluated (that is, the last expression in the clause selected). If there is no match, case returns nil.

The symbol t has special meaning as a comparator: it matches anything. It is typically used in the last clause to serve as a default case when no match is found with other clauses.

# Language Constructs

### **Arguments**

 $g\_selectionExpr$  An expression whose value is evaluated and tested for equality

against the comparators in each clause. When a match is

found, the rest of the clause is evaluated.

1 clause1 An expression whose first element is an atom or list of atoms to

be compared against the value of  $g\_selectionExpr$ . The remainder of the l clause is evaluated if a match is found.

1\_clause2... Zero or more clauses of the same form as 1\_clause1.

#### Value Returned

g result Returns the value of the last expression evaluated in the

matched clause.

nil Returns nil if there is no match.

# **Example**

Sets path to ./min.

#### cond

### **Description**

Examines conditional clauses from left to right until either a clause is satisfied or there are no more clauses remaining.

This command is useful when there is more than one test condition, but only the statements of one test are to be carried out. Each clause is of the form  $(g\_condition g\_expr1...)$ . The cond function examines a clause by evaluating the condition associated with the clause. The clause is satisfied if  $g\_condition$  evaluates to non-nil, in which case expressions in the rest of the clause are evaluated from left to right, and the value returned by the last expression in the clause is returned as the value of the cond form. If  $g\_condition$  evaluates to nil, however, cond skips the rest of the clause and moves on to the next clause.

# **Arguments**

1	c1	а	и	S	e	1
_	$\sim$	~	~	$\sim$	$\overline{}$	_

Each clause must be of the form  $(g\_condition g\_expr1...)$ . When  $g\_condition$  evaluates to non-nil, all the succeeding expressions are evaluated.

#### Value Returned

g\_result
nil

Returns the value of the last expression of the satisfied clause.

Returns nil if no clause is satisfied.

# **Example**

# Language Constructs

```
=> nil; Prints "Arg is a number".
test('sym)
=> nil; Prints "Arg is an unknown type".
```

Tests each of the arguments according to the conditions specified with cond.

Language Constructs

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# **File Commands and Functions**

This chapter contains information on the following commands:

close

**fscanf** 

<u>gets</u>

infile

<u>load</u>

<u>newline</u>

<u>outfile</u>

<u>pfile</u>

printf

<u>println</u>

#### File Commands and Functions

# close

```
close(
    p_port
)
=> t
```

# **Description**

Drains, closes, and frees a port.

When a file is closed, it frees the FILE\* associated with  $p\_port$ . Do not use this function on piport, stdin, poport, stdout, or stderr.

# **Arguments**

p port

Name of port to close.

#### **Value Returned**

t

The port closed successfully.

# **Example**

```
p = outfile( "~/test/myFile" ) => port:"~/test/myFile"
close( p )
=> t
```

Drains, closes, and frees the /test/myFile port.

#### File Commands and Functions

### fscanf

```
fscanf(
    p_inputPort
    t_formatString
    [ s_var1 ... ]
)
    => x items / nil
```

# Description

Reads input from a port according to format specifications and returns the number of items read in.

The results are stored into corresponding variables in the call. The fscanf function can be considered the inverse function of the fprintf output function. The fscanf function returns the number of input items it successfully matched with its format string. It returns nil if it encounters an end of file.

The maximum size of any input string being read as a string variable for fscanf is currently limited to 8 K. Also, the function lineread is a faster alternative to fscanf for reading Virtuoso® SKILL objects.

The common input formats accepted by fscanf are summarized below. Common Input Format Specifications

Format Specification	Types of Argument	Scans for
%d	fixnum	An integer
%f	flonum	A floating-point number
%s	string	A string (delimited by spaces) in the input

#### File Commands and Functions

### **Arguments**

 $p\_inputPort$  Input port to read from.  $t\_formatString$  Format string to match against in the reading.  $s\_var1...$  Name of the variable in which to store results.

#### Value Returned

x_items	Returns the number of input items it successfully read in. As a side effect, the items read in are assigned to the corresponding variables specified in the call.
nil	Returns nil if an end of file is encountered.

# **Example**

```
fscanf( p "%d %f" i d )
```

Scans for an integer and a floating-point number from the input port p and stores the values read in the variables i and d, respectively.

Assume a file testcase with one line:

```
hello 2 3 world
x = infile("testcase")
=> port:"testcase"
fscanf( x "%s %d %d %s" a b c d )
=> 4
(list a b c d) => ("hello" 2 3 "world")
```

#### File Commands and Functions

# gets

```
gets(
    s_variableName
    [ p_inputPort ]
)
    => t string / nil
```

# **Description**

Reads a line from the input port and stores the line as a string in the variable. This is a macro.

The string is also returned as the value of gets. The terminating newline character of the line becomes the last character in the string.

# **Arguments**

$s\_variableName$	Variable in which to store the input string.
p_inputPort	Name of input port.
	Default value: piport

#### Value Returned

t_string	Returns the input string when successful.
nil	Returns nil when the end of file is reached.
	(s variableName maintains its last value.)

#### **Example**

Assume the test1.data file has the following first two lines:

```
#This is the data for test1
0001 1100 1011 0111
p = infile("test1.data") => port:"test1.data"
gets(s p) => "#This is the data for test1"
gets(s p) => "0001 1100 1011 0111"
s => "0001 1100 1011 0111"
```

Gets a line from the test1.data file and stores it in the variable s. The s variable contains the last string stored in it by the gets function.

#### File Commands and Functions

### infile

```
infile(
    S_fileName
)
    => p inport / nil
```

# **Description**

Opens an input port ready to read a file.

Always remember to close the port when you are done. The file name can be specified with either an absolute path or a relative path. In the latter case, the current SKILL path is used if it is not nil.

### **Arguments**

S fileName

Name of the file to be read; it can be either a string or a symbol.

#### Value Returned

p\_inportReturns the port opened for reading the named file.nilReturns nil if the file does not exist or cannot be opened for

reading.

### Example

```
in = infile( "~/test/input.il" ) => port:"~/test/input.il"
close( in )
=> t
```

Closes the /test/input.il port.

Opens the input port /test/input.il.

```
infile("myFile") => nil
```

Returns nil if myFile does not exist according to the current setting of the SKILL path or exists but is not readable.

#### File Commands and Functions

### load

```
load(
     t_fileName
     [ t_password ]
    )
     => t
```

# Description

Opens a file and repeatedly calls lineread to read in the file, immediately evaluating each form after it is read in.

This function uses the file extension to determine the language mode (.il for SKILL, .ils for SKILL++, and .ocn for a file containing OCEAN commands) for processing the language expressions contained in the file. For a SKILL++ file, the loaded code is always evaluated in the top-level environment.

load closes the file when the end of file is reached. Unless errors are discovered, the file is read in quietly. If load is interrupted by pressing Control-c, the function skips the rest of the file being loaded.

SKILL has an autoload feature that allows applications to load functions into SKILL on demand. If a function being run is undefined, SKILL checks to see if the name of the function (a symbol) has a property called autoload attached to it. If the property exists, its value, which must be either a string or an expression that evaluates to a string, is used as the name of a file to be loaded. The file should contain a definition for the function that triggered the autoload. Processing proceeds normally after the function is defined.

#### File Commands and Functions

### **Arguments**

t fileName

File to be loaded. Uses the file name extension to determine the language mode to use. Valid values:

- .ils Means the file contains SKILL++ code.
- .ocn Means the file contains OCEAN commands (with SKILL or SKILL++ commands)

t password

Password, if t fileName is an encrypted file.

#### Value Returned

t

Returns t if the file is successfully loaded.

# **Example**

```
load( "test.ocn" )
```

Loads the test.ocn file.

```
procedure( trLoadSystem()
    load( "test.il" ) ;;; SKILL code
    load( "test.ils" );;; SKILL++ code
) ; procedure
```

You might have an application partitioned into two files. Assume that test.il contains SKILL code and test.ils contains SKILL/SKILL++ code. This example loads both files.

# File Commands and Functions

# newline

```
newline(
        [ p_outputPort ]
    )
    => nil
```

# **Description**

Prints a newline (backslash n) character and then flushes the output port.

# **Arguments**

p\_outputPort Output port.

Defaults value: poport

# Value Returned

nil Prints a newline and then returns nil.

# **Example**

```
print( "Hello" ) newline() print( "World!" )
"Hello"
"World!"
=> nil
```

Prints a newline character after Hello.

#### File Commands and Functions

### outfile

```
outfile(
    S_fileName
    [ t_mode ]
    )
    => p outport / nil
```

### **Description**

Opens an output port ready to write to a file.

Various print commands can write to this file. Commands write first to a character buffer, which writes to the file when the character buffer is full. If the character buffer is not full, the contents are not written to the file until the output port is closed or the drain command is entered. Use the close or drain command to write the contents of the character buffer to the file. The file can be specified with either an absolute path or a relative path. If a relative path is given and the current SKILL path setting is not nil, all directory paths from SKILL path are checked in order, for that file. If found, the system overwrites the first updatable file in the list. If no updatable file is found, it places a new file of that name in the first writable directory.

# **Arguments**

S	fileName	Name of the file to open or create.
$\nu_{-}$	_ I I I CIVainC	riamo er mo mo to open er ereate.

t\_mode Mode in which to open the file. If a, the file is opened in append

mode; If w, a new file is created for writing (any existing file is

overwritten).

Default value: w

#### Value Returned

p\_outport An output port ready to write to the specified file.

nil returns nil if the named file cannot be opened for writing. An

error is signaled if an illegal mode string is supplied.

### **Example**

```
p = outfile( "/tmp/out.il" "w" )
=> port:"/tmp/out.il"
```

Opens the /tmp/out.il port.

# File Commands and Functions

```
outfile( "/bin/ls" )
=> nil
```

Returns nil, indicating that the specified port could not be opened.

#### File Commands and Functions

# pfile

# **Description**

Opens an output port ready to write to a file or returns the name of an existing port indicating that it is available.

This command is similar to the outfile command when a valid S\_fileName is specified. When p\_port is specified, it returns the file port that is currently being used by p\_port. When no argument is specified, it opens the stdout port.

# **Arguments**

$S_{\_}$ fileName	Name of the file to open or create.
p_port	Retrieves the name of the file port that is being used.

#### Value Returned

p_port	The ID of the port that was opened, or stdout.
nil	Returns nil if the named file cannot be opened for writing.

#### **Example**

```
p = pfile( "/tmp/out.il" "w" )
=> port:"/tmp/out.il"

Opens the /tmp/out.il port.

pfile( "/bin/ls" )
=> nil
```

Returns nil, indicating that the specified port could not be opened.

```
p = pfile()
=> port:"*stdout*"
```

Returns stdout as the file port indicating that stdout has been opened.

```
pfile( p )
=> port:"/tmp/out.il"
```

# File Commands and Functions

Returns the file port.

#### File Commands and Functions

# printf

# **Description**

Writes formatted output to *poport*, which is the standard output port.

The optional arguments following the format string are printed according to their corresponding format specifications. Refer to the "Common Output Format Specifications" table for fprintf in the Cadence SKILL Language User Guide.

printf is identical to fprintf except that it does not take the  $p\_port$  argument and the output is written to poport.

# **Arguments**

t_formatString	Characters to be printed verbatim, intermixed with format specifications prefixed by the "%" sign.
g_arg1	Arguments following the format string are printed according to their corresponding format specifications.

#### Value Returned

t Prints the formatted output and returns t.

### Example

```
x = 197.9687 \Rightarrow 197.9687
printf( "The test measures %10.2f." x )
```

Prints the following line to poport and returns t.

```
The test measures 197.97. \Rightarrow t
```

#### File Commands and Functions

# println

```
println(
    g_value
    [ p_outputPort ]
)
    => nil
```

# **Description**

Prints a SKILL object using the default format for the data type of the value, and then prints a newline character.

A newline character is automatically printed after printing  $g_value$ . The println function flushes the output port after printing each newline character.

# **Arguments**

g_value	Any SKILL value.
---------	------------------

p\_outputPort
Port to be used for output.

Default value: poport

#### **Value Returned**

nil Prints the given object and returns nil.

# Example

```
for( i 1 3 println( "hello" ))
"hello"
"hello"
"hello"
=> t
```

Prints hello three times. for always returns t.

File Commands and Functions

# **OCEAN 4.4.6 Issues**

For the 4.4.6 release of OCEAN, there are some restrictions and requirements.

The netlist file that you specify for the Spectre<sup>®</sup> circuit simulator interface with the design command must be netlist. The full path can be specified. For example, /usr/netlist is acceptable. The netlistHeader and netlistFooter files are searched in the same directory where the netlist is located. Cadence recommends that you use the netlist generated from the Virtuoso® Analog Design Environment. Netlists from other sources can also be used, as long as they contain only connectivity. You might be required to make slight modifications.

- Cadence recommends full paths for the Spectre simulator model files, definition files, and stimulus files.
- The Cadence SPICE circuit simulator is still used to parse netlists for socket interfaces (spectreS and cdsSpice, for example). Therefore, the netlist that you specify with the design command must be in Cadence SPICE syntax. Cadence recommends that you use the raw netlist generated from the Virtuoso® Analog Design Environment. Netlists from other sources can also be used, as long as they can pass through Cadence SPICE. You might be required to make slight modifications.
- Any presimulation commands that you specify are appended to the final netlist (as is currently the case in the design environment). Therefore, if you have control cards already in your netlist, and specify simulation setup commands, you might duplicate control cards, which causes a warning or an error from the simulator. You might want to remove control cards from your netlist file to avoid the warnings.
- Models, include files, stimulus files, and PWLF files must be found according to the path specified with the <u>path</u> command.

# OCEAN Reference OCEAN 4.4.6 Issues

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