

OCEAN Reference

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Preface

The preface discusses the following:

- [Related Documents](#) on page 15
- [Typographic and Syntax Conventions](#) on page 15

Related Documents

The Open Command Environment for Analysis (OCEAN) is based on the Virtuoso® SKILL programming language. The following manuals give you more information about the SKILL language and other related products.

- The [SKILL Language User Guide](#) describes how to use the SKILL language functions, the SKILL++ functions, and the SKILL++ object system (for object-oriented programming).
- The [SKILL Language Reference](#) provides descriptions, syntax, and examples for the SKILL and SKILL++ functions.
- The [SKILL++ Object System Functions Reference](#) provides descriptions, syntax, and examples for the object system functions.
- The [Virtuoso® Analog Design Environment User Guide](#) explains how to design and simulate analog circuits.
- The [Virtuoso® Mixed-Signal Circuit Design Environment User Guide](#) explains how to design and simulate mixed-signal circuits.
- The [Virtuoso® Analog Distributed Processing Option User Guide](#) explains how to set up and run distributed processing for OCEAN and other Virtuoso® Analog Design Environment applications.

Typographic and Syntax Conventions

This list describes the syntax conventions used for the Virtuoso® Analog Design Environment SKILL functions.

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Preface

<code>literal</code>	Nonitalic words indicate keywords that you must type literally. These keywords represent command (function, routine) or option names.
<code>argument (z_argument)</code>	Words in italics indicate user-defined arguments for which you must substitute a name or a value. (The characters before the underscore (<code>_</code>) in the word indicate the data types that this argument can take. Names are case sensitive. Do not type the underscore (<code>z_</code>) before your arguments.) For a listing of data types, see “Data Types Used in OCEAN” on page 24.
<code> </code>	Vertical bars (OR-bars) separate possible choices for a single argument. They take precedence over any other character.
<code>[]</code>	Brackets denote optional arguments. When used with OR-bars, they enclose a list of choices. You can choose one argument from the list.
<code>{ }</code>	Braces are used with OR-bars and enclose a list of choices. You must choose one argument from the list.
<code>...</code>	Three dots (<code>...</code>) indicate that you can repeat the previous argument. If you use them with brackets, you can specify zero or more arguments. If they are used without brackets, you must specify at least one argument, but you can specify more. <code>argument...</code> Specify at least one, but more are possible. <code>[argument]...</code> Specify zero or more.
<code>,...</code>	A comma and three dots together indicate that if you specify more than one argument, you must separate those arguments by commas.
<code>=></code>	A right arrow precedes the possible values that a SKILL function can return. This character is represented by an equal sign and a greater than sign.
<code>/</code>	A slash separates the possible values that can be returned by a SKILL function.

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<*yourSimulator*>

Angle brackets indicate places where you need to insert the name of your simulator. Do not include the angle brackets when you insert the simulator name.

Important

The characters included in the list above are the only characters that are not typed literally. All other characters in the SKILL language are required and must be typed literally.

SKILL Syntax Examples

The following examples show typical syntax characters used in the SKILL language.

Example 1

```
list( g_arg1 [g_arg2] ...
      )
      => l_result
```

Example 1 illustrates the following syntax characters.

<code>list</code>	Plain type indicates words that you must type literally.
<code><i>g_arg1</i></code>	Words in italics indicate arguments for which you must substitute a name or a value.
<code>()</code>	Parentheses separate names of functions from their arguments.
<code>_</code>	An underscore separates an argument type (left) from an argument name (right).
<code>[]</code>	Brackets indicate that the enclosed argument is optional.
<code>=></code>	A right arrow points to the return values of the function. Also used in code examples in SKILL manuals.
<code>...</code>	Three dots indicate that the preceding item can appear any number of times.

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Example 2

```
needNCells(  
  s_cellType | st_userType  
  x_cellCount  
)  
=> t/nil
```

Example 2 illustrates two additional syntax characters.

- | Vertical bars separate a choice of required options.
- / Slashes separate possible return values.

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](#) on page 20
- [OCEAN Online Help](#) on page 20
- [OCEAN Syntax Overview](#) on page 21
- [Parametric Analysis](#) on page 27
- [Distributed Processing](#) on page 28

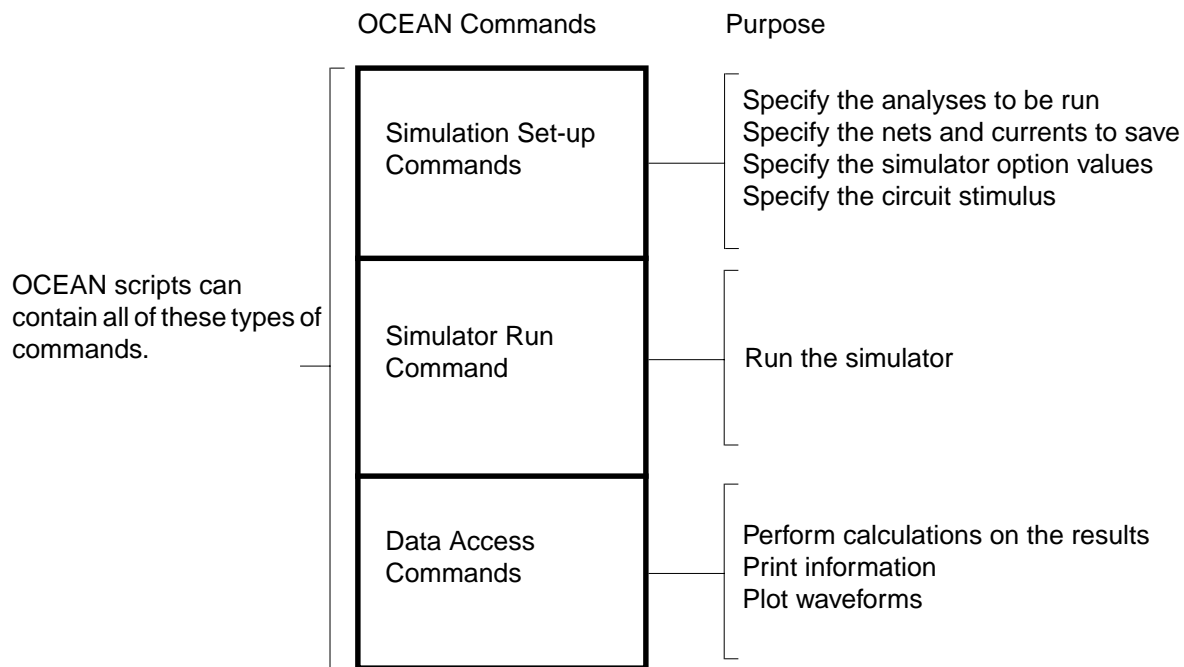
OCEAN lets you set up, simulate, and analyze circuit data. OCEAN is a text-based process that you can run from a UNIX shell or from the Command Interpreter Window (CIW). You can type OCEAN commands in an interactive session, or you can create scripts containing your commands, then load those scripts 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 easily rerun your scripts. OCEAN lets you

- Create scripts that you can run repeatedly to verify circuit performance
- Run longer analyses such as parametric analyses, Corners 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.)

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.

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

```
ocnHelp( )
```

For more information, see [“ocnHelp”](#) on page 129.

OCEAN Syntax Overview

OCEAN is based on the Virtuoso® 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 [Chapter 3, “Introduction to SKILL.”](#)

Parentheses

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

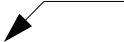
Examples

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.

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

 Syntax error.

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>
```

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 75, 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 111 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:

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

Keyword Value passed
 ↓ ↓
analysis('tran ?start 0 ?stop 1u ?step 1n)

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.

```
report([?output t_filename | p_port] [?type t_type] [?name t_name]  
      [?param t_param] [?format s_reportStyle] ) => t/nil
```

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	l
integer, floating-point number, or complex number		n

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Data Type	Internal Name	Prefix
user-defined type		o
I/O port	port	p
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
<code>a=desVar("r1" 1k)</code>	<code>a=1k</code>
<code>a=desVar("r1" 1k "r2" 2k)</code>	<code>a=2k</code>
<code>a=desVar("r1")</code>	<code>a=1k</code> , assuming <code>r1</code> was set in a <code>desVar</code> command
<code>a=desVar("r2")</code>	<code>a=2k</code> , assuming <code>r2</code> was set in a <code>desVar</code> 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 )
```

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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.

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.)

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( "r1" ?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

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.

```
Cloud = list( 2p 4p 6p 8p )
foreach( val Cloud
  desVar( "Cloud" val )
  a=resultsDir( sprintf( nil "./demo/Cloud=%g" val ) )
  printf( "%L", a )
  run( )
)
foreach( val Cloud
  openResults( sprintf( nil "<dir>/Cloud=%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 [Chapter 7, “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

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 [Virtuoso® Analog Distributed Processing](#)

Option User Guide for more information on how to configure systems for distributed processing. For information on the distributed processing commands for OCEAN, see Chapter 12, “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
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 `wait` command to block a single job. The `wait` 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.

Waveform Tool Selection

You can plot simulation results in the waveform tool of your choice. The Analog Design Environment now supports WaveScan in addition to AWD. Although WaveScan is the default, you can dynamically switch between the two waveform tools using the OCEAN function `ocnWaveformTool()` as follows:

```
ocnWaveformTool('awd)
```

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](#) on page 31
- [Using OCEAN Interactively](#) on page 32
- [Creating OCEAN Scripts](#) on page 35
- [Running Multiple Simulators](#) on page 41
- [OCEAN Tips](#) on page 41

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*
 - By hand (by you or someone else in your organization) with a text editor

For information about creating scripts, see [“Creating OCEAN Scripts”](#) on page 35.

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 [“OCEAN Online Help”](#) on page 20.

Note: The current version of OCEAN has some specific issues that are addressed in [Appendix A, “OCEAN 4.4.6 Issues.”](#) Please refer to this appendix before using OCEAN.

Using OCEAN Interactively

You can run OCEAN from a UNIX prompt or from the Virtuoso® 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 loads and reads your `.oceanrc` file.

You can place OCEAN commands in your `.oceanrc` file, which is similar to the `.cdsinit` file. OCEAN ignores your `.cdsinit` file at startup. If you want OCEAN to use any initialization options at startup, you must specify them in the `.oceanrc` file. This file can contain any valid OCEAN command, function or SKILL initialization routine (excluding graphical dfll 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 [“Interactive Session Demonstrating the OCEAN Use Model”](#) on page 34.

To quit 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:

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```
#!/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 `-nograph` option with the `ocean` 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.

The `-nograph` option must only be used to replay logfiles that have been created interactively. For example, while using OCEAN with the `-nograph` option, your `oceanScript.ocn` file must have an `exit()` statement at the end. 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 explicit `exit()` statement at the end of the OCEAN script and OCEAN exits only when it detects an `exit()` at EOF. The command is used as follows:

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

While using the `-nograph` option with `ocean`, if you find that simulation run messages are not being stored in the log file, check for the following environment variable in the `.cdsenv` 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 User Guide*.

Using OCEAN from the CIW

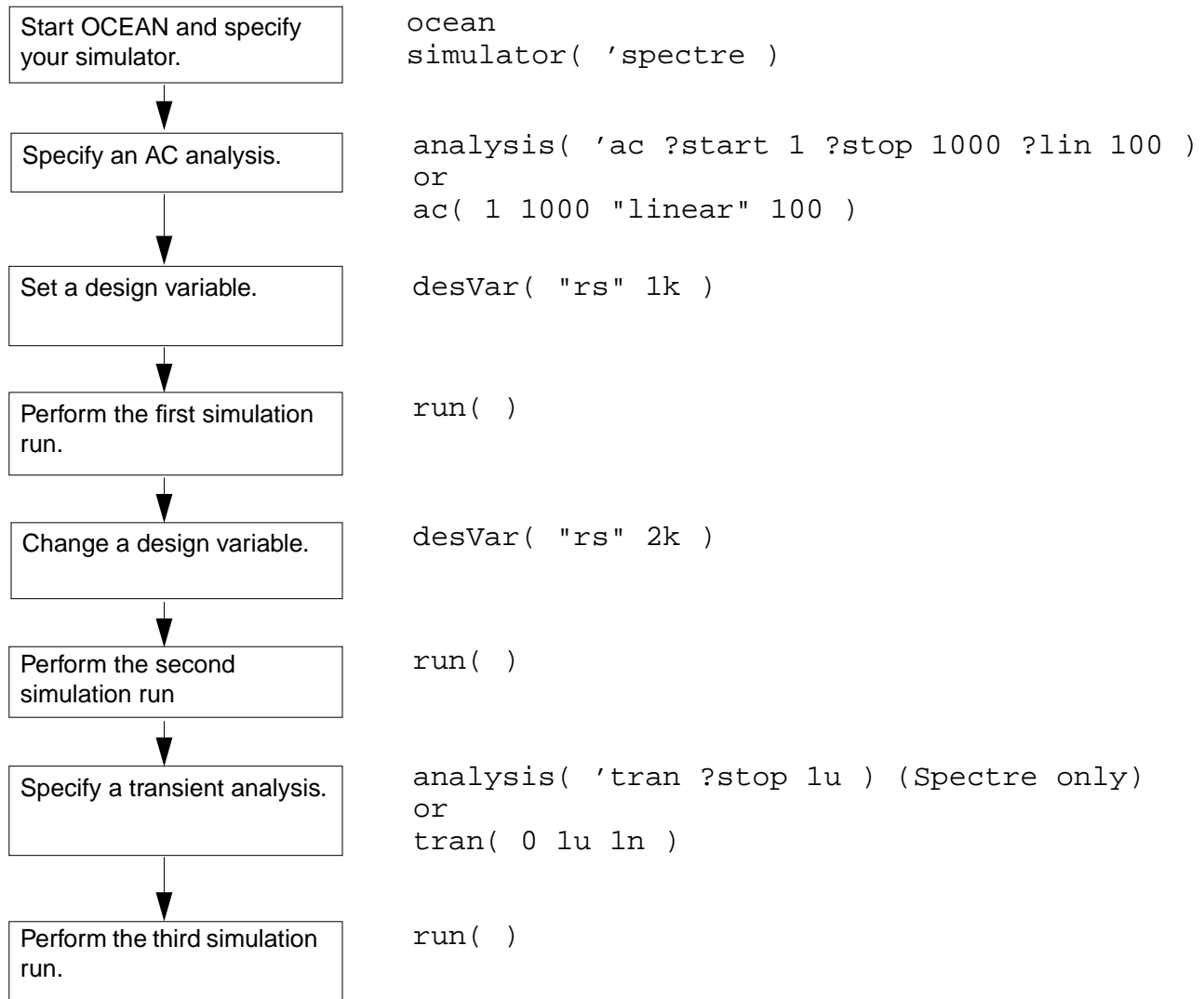
You can type OCEAN commands in the CIW after you bring up the Virtuoso® Analog Design Environment. (Starting the design environment loads the required OCEAN files.)

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

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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.

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

Creating OCEAN Scripts

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

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 `simulatorx.ocn` file in your `netlist` 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 `cdsSpice0.ocn` file is created in your `netlist` directory. You can load this `cdsSpice0.ocn` script as described in [“Loading OCEAN Scripts”](#) on page 38.

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.

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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.

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 `simulator0.ocn` script, which is automatically created in the `netlist` 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 the Virtuoso® Analog Design Environment,

1. Start the Virtuoso software with the executable you prefer; for example,

```
icms&
```

The CIW appears.
2. From the CIW, choose *Tools – Analog Environment – 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 38.

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

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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
```

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 [Virtuoso® Analog Design Environment SKILL Language Reference](#).

3. Load the functions in your `.cdsinit` file.

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

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```
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.

```
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 with the executable you prefer, for example

```
icms&
```

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.

Note: OCEAN does not read your `.cdsinit` file. If you want your `.cdsinit` file read, you must load it in your `.oceanrc` file.

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 Artist to select the results and work with features like annotation etc.

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 corners or 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 Artist.

In Artist, `psf` data is stored in:

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

where,

`runDir` is the directory where you invoke `icfb&`

`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 *Artist -> Results -> Select*

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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 on c1, c2 or c3. Alternatively, you can also single click on 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 *Artist -> Results -> Direct Plot*, *Artist -> Results -> Annotate* etc to see the results of that particular directory.

Selecting Results Run from Spectre Stand Alone

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 Artist features like *Artist -> Results -> Direct Plot* or *Artist -> Results -> Annotate*.

Consider that your data is in

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

where,

runDir is the directory where you invoke icfb&

testSchemName is your test schematic

1. Choose *Artist -> 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 on schematic directory. The psf directory should appear.
4. Click on the directory with the data in it, psf. When you click on 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".

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Note: Even though there is a confirmation message in the CIW that the select was success, Artist is not synced up to allow cross-probing and back annotation of operating points.

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

Note: You CAN use Artist -> Tools > Calculator but you CAN NOT use *Artist -> Results -> Direct Plot* or *Artist -> 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 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 `fromVal` 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:

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```
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 [“Recovering from an Omitted Quotation Mark”](#) on page 22.
- 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
```

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[®] SKILL programming language. In this chapter, you can find information about

- [The Advantages of SKILL](#) on page 43
- [Naming Conventions](#) on page 44
- [Arithmetic Operators](#) on page 44
- [Scaling Factors](#) on page 44
- [Relational and Logical Operators](#) on page 46
- [SKILL Syntax](#) on page 48
- [Arithmetic and Logical Expressions](#) on page 51

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, `dmiPurgeVersions()` or `hnlCellOutputs`. All code developed outside Cadence should name global variables by starting them with an uppercase character, such as `AcmeGlobalForm`.

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.

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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
Y	Yotta	10 ²⁴	10Y [10e+25]
Z	Zetta	10 ²¹	10Z [10e+22]
E	Exa	10 ¹⁸	10E [10e+19]
P	Peta	10 ¹⁵	10P [10e+16]
T	Tera	10 ¹²	10T [1.0e13]
G	Giga	10 ⁹	10G [10,000,000,000]
M	Mega	10 ⁶	10M [10,000,000]
K	Kilo	10 ³	10K [10,000]
%	percent	10 ⁻²	5% [0.05]
m	milli	10 ⁻³	5m [5.0e-3]
u	micro	10 ⁻⁶	1.2u [1.2e-6]
n	nano	10 ⁻⁹	1.2n [1.2e-9]
p	pico	10 ⁻¹²	1.2p [1.2e-12]
f	femto	10 ⁻¹⁵	1.2f [1.2e-15]
a	atto	10 ⁻¹⁸	1.2a [1.2e-18]
z	zepto	10 ⁻²¹	1.2z [1.2e-21]
y	yocto	10 ⁻²⁴	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 10⁶.

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	t
			3 < 2	nil
<=	numeric	leqp	3 <= 4	t
>	numeric	greaterp	5 > 3	t
>=	numeric	geqp	4 >= 3	t
==	numeric	equal	3.0 == 3	t
	string list		"abc" == "ABc"	nil
!=	numeric	nequal		
	string list		"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	<code>3 && 5</code>	5
			<code>5 && 3</code>	3
			<code>t && nil</code>	nil
			<code>nil && t</code>	nil
	general	or	<code>3 5</code>	3
			<code>5 3</code>	5
			<code>t nil</code>	t
			<code>nil t</code>	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

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
<code>\</code>	backslash	Escape for special characters
<code>()</code>	parentheses	Grouping of list elements, function calls
<code>[]</code>	brackets	Array index, super right bracket
<code>'</code>	single quotation mark	Specifies a symbol (quoting the expression prevents its evaluation)
<code>"</code>	quotation mark	String delimiter
<code>,</code>	comma	Optional delimiter between list elements

Special Characters in SKILL

Character	Name	Meaning
;	semicolon	Line-style comment character
+, -, *, /	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.  
*/
```

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

```
x = 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 be very careful to note.

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(arg1 arg2) 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)`.
- 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 simply 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" .

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.

```
lineCount = 4           => 4  
lineCount               => 4
```

```
lineCount = "abc"      => "abc"  
lineCount           => "abc"
```

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: $1+2$, $a*b+c$, $x>y$.

You can form arbitrarily complicated expressions by combining any number of the primitive expressions described above.

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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](#) on page 55
- [Data Types](#) on page 55
- [Arrays](#) on page 58
- [Concatenating Strings \(Lists\)](#) on page 58
- [Declaring a SKILL Function](#) on page 60
- [Skill Function Return Values](#) on page 62
- [Syntax Functions for Defining Functions](#) on page 62

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

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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	a
boolean		b
floating-point number	flonum	f
any data type	general	g
linked list	list	l
floating-point number or integer		n
user-defined type		o
I/O port	port	p
symbol	symbol	s
symbol or character string		S
character string (text)	string	t
window type		w
integer number	fixnum	x

Numbers

SKILL supports the following numeric data types:

- Integers
- Floating-point

Both integers and floating-point numbers may use scaling factors to scale their values. For information on scaling factors, see [“Scaling Factors” on page 44](#).

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 [“Arithmetic and Logical Expressions” 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.

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.

```
declare( week[7] )           => array[7]:9780700
week                        => array[7]:9780700
type( week )                => array
days = '(monday tuesday wednesday
         thursday friday saturday sunday)
for( day 0 length(week)-1
     week[day] = nth(day days))
```

- 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.

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```
buildString( '( "test" "il" ) "." )      => "test.il"
buildString( '( "usr" "mnt" ) "/" )      => "usr/mnt"
buildString( '( "a" "b" "c" ) )         => "a b c"
buildString( '( "a" "b" "c" ) "" )      => "abc"
```

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.

```
strncat( "abcd" "efghi" 2)      => "abcdef"
strncat( "abcd" "efghijk" 5)    => "abcdefghi"
```

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:

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

Comparing Two Strings Alphabetically (strcmp)

`strcmp` compares two strings. (To simply 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	<i>string1</i> is alphabetically greater than <i>string2</i> .
0	<i>string1</i> is alphabetically equal to <i>string2</i> .
-1	<i>string1</i> is alphabetically less than <i>string2</i> .

```
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	<i>arg1</i> is alphanumerically greater than <i>arg2</i> .
0	<i>arg1</i> is alphanumerically identical to <i>arg2</i> .
-1	<i>arg2</i> is alphanumerically greater than <i>arg1</i> .

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.

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`.

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
```

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 actual parameters bound to the formal parameters. 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.

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The actual 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.

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OCEAN Environment Commands

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

[appendPath](#) on page 66

[path](#) on page 67

[prependPath](#) on page 68

[setup](#) on page 69

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

<i>t_dirName1</i>	Directory path.
<i>t_dirNameN</i>	Additional directory path.

Value Returned

<i>t_dirNameN</i>	Returns the last path specified.
<i>nil</i>	Returns <i>nil</i> 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.

path

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

Description

Sets the search path for included files.

This command overrides the path set earlier using any of these commands: [path](#), [appendPath](#), or [prependPath](#).

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

<code>t_dirName1</code>	Directory path.
<code>t_dirNameN</code>	Additional directory path.

Value Returned

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

Examples

```
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.

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

<i>t_dirName1</i>	Directory path.
<i>t_dirNameN</i>	Additional directory path.

Value Returned

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

Examples

```
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.

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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.

Arguments

s_numberNotation

Specifies the notation for printed information.

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.

x_precision

Specifies the number of significant digits that are printed.

Valid values: 1 through 16

Default value: 6

s_reportStyle

Specifies the format of the output of the report command.

Valid values: spice, paramValPair

Default value: paramValPair

The spice format is:

	Param1	Param2	Param3
Name1	value	value	value
Name2	value	value	value
Name3	value	value	value

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OCEAN Environment Commands

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

x_charsPerLine Specifies the number of characters per line output to the display.
Default value: 80

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.

Examples

```
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`.

Simulation Commands

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

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OCEAN Reference 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.

Arguments

<i>g_fromValue</i>	Starting value for the AC analysis.
<i>g_toValue</i>	Ending value.
<i>g_ptsPerDec</i>	Points per decade.
<i>t_incType</i>	Increment type. Valid values: For the Spectre® circuit simulator, "Linear", "Logarithmic", or "Automatic". For other simulators, "Linear" or "Logarithmic".
<i>g_points</i>	Either the linear or the logarithmic value, which depends on <i>t_incType</i> .

Value Returned

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

Examples

```
ac(1 10000 2)
```

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

```
ac(1 10000 "Linear" 100)
```

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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.

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 *analysisOption1* 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 [?from 0].

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.

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*

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 [Spectre Circuit Simulator Reference](#) manual for analysis options you can use.

g_analysisOptionValue1

Value for the analysis option.

?*analysisOptionN*

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

OCEAN Reference Simulation Commands

g_analysisOptionValueN

Value for the analysis option.

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

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

For the Spectre® circuit simulator, specifies that an AC analysis be performed.

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

Specifies that a transient analysis be performed.

```
analysis('dc ?oppoint "rawfile" ?save "allpub"  
?param "temp" ?start -50 ?stop 100 )
```

Sweeps temperature for the Spectre® circuit simulator.

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

Saves the DC operating point information for the Spectre® circuit simulator.

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

Sets the Spectre transfer function analysis.

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

Sets the Spectre sensitivity analysis.

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

Sets the Spectre noise analysis.

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

Sets the Spectre dcmatch analysis.

```
analysis('pz ?freq "2" ?readns "./abc" ?oppoint "rawfile" ?fmax "4500000000"  
?zeroonly "no" ?prevoppoint "no" ?restart "no" ?annotate "no" ?stats "no" )
```

Sets the Spectre pz analysis.

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```
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" )
```

Sets the Spectre stability analysis.

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

Sets the Spectre pss RF analysis.

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

Sets the Spectre pnoise RF analysis.

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

Sets the Spectre pac RF analysis.

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

Sets the Spectre pxf RF analysis.

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

Sets the Spectre qpss RF analysis.

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

Sets the Spectre qpac analysis.

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

Sets the Spectre sp (S - parameter) analysis.

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, such as spectreS. For direct simulators, such as spectre, use `createNetlist` instead.

Arguments

None.

Value Returned

<code>t</code>	Returns <code>t</code> if the final netlist is created.
<code>nil</code>	Returns <code>nil</code> and prints an error message otherwise.

Example

```
createFinalNetlist()
```

Creates the final netlist for the current simulation run.

createNetlist

```
createNetlist( [?recreateAll b_recreateAll] [?display b_display] )  
=> t_filename/nil
```

Description

Creates the simulator input file.

If the design is specified as lib/cell/view, this command netlists the design, if required, and creates the simulator input file. When the `b_recreateAll` argument is set to `t` and the design is specified as lib/cell/view, 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 [design](#) function.

When the `b_display` option is set to `t` (or `nil`) the netlist file is displayed (or undisplayed) to the user. By default, `b_display` is set to `'t` (true).

Note: This command does not work with socket simulators.

Arguments

<code>b_recreateAll</code>	If set and the design is specified as lib/cell/view, the entire netlist is recreated.
----------------------------	---

Value Returned

<code>t_filename</code>	Returns the name of the simulator input file on success.
<code>nil</code>	otherwise <code>nil</code> is returned

Examples

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

Creates simulator input file for the current simulation run.

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

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

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```
design( ?lib "test" ?cell "mytest1" ?view "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 netlist file would be displayed. This can be turned ON/OFF via `?display` set to `t/nil`

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 *Helping a Simulation to Converge* of the *Virtuoso Analog Design Environment User Guide*.

Arguments

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

Value Returned

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

Examples

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

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

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

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

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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 `analysis` command.

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

Note: `t_compParam` is valid only for the spectre, spectreS, spectreVerilog and spectreSVerilog simulators.

Arguments

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

Value Returned

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

Examples

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

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

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

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

definitionFile

```
definitionFile( t_fileName [t_fileName2 ... t_fileNameN ] )  
=> l_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 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

l_fileNames 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.

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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 those particular pieces of information are deleted as opposed to deleting all the information for that command.

Arguments

<i>s_command</i>	Command that was initially used to add the items that are now being deleted. Valid values: analysis, desVar, path, save, ic, forcenode, monteCarlo, monteExpr, nodeset, optimizeGoal, optimizeVar, optimizeAlgoControl, optimizePlotOption Using <code>delete('monteCarlo)</code> turns off the monteCarlo command and sets everything back to the defaults.
<i>g_commandArg1</i>	Argument corresponding to the specified command.
<i>g_commandArg2</i>	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.

Examples

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

Deletes all the saves.

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

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.

design

```
design( t_cktFile )
    => t_cktFile/nil

design( t_lib t_cell t_view )
    => (t_lib t_cell t_view)/nil

design( t_lib t_cell t_view t_mode )
    => (t_lib t_cell t_view)/nil
```

Description

Specifies the name of the design to be simulated.

Note: You can use the *lib*, *cell*, *view* version of the `design` command only if you are running OCEAN within `icms`, `msfb`, or `icfb`. You cannot use this version of the command within the OCEAN process itself.

Arguments

<i>t_cktFile</i>	<p>For the direct simulator, the name of the netlist. The name must end in <code>netlist</code>. Note that the <code>netlistHeader</code> and <code>netlistFooter</code> files are also needed in the same directory.</p> <p>For socket simulators, this is the name of the raw circuit file. If generated in the Virtuoso® Analog Design Environment, the file is named <code>design.c</code> and is found in the <code>netlist</code> directory.</p> <p>Otherwise, <i>cktFile</i> is a pre-existing netlist file from another source. In this case, you might need to remove the <code>.cards</code> from the netlist because the OCEAN commands are converted to <code>.cards</code> and appended to the final netlist. The simulator might give an error or warning if the <code>.cards</code> are read twice.</p>
<i>t_lib</i>	Name of the Virtuoso® Analog Design Environment library that contains the design.
<i>t_cell</i>	Name of the design.
<i>t_view</i>	View of the design (typically <code>schematic</code>).
<i>t_mode</i>	The mode in which the design should be opened. The value can be <code>r</code> , <code>w</code> or <code>a</code> , representing <code>read</code> , <code>write</code> and <code>append</code> ,

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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

<code>t_cktFile</code>	Returns the name of the design if successful.
<code>l_(lib cell view)</code>	Returns the name of the view for an Virtuoso® Analog Design Environment design if successful.
<code>nil</code>	Returns <code>nil</code> and prints an error message if there is a problem using the specified design.

Examples

For the Spectre® circuit simulator,

```
design( "netlist" )  
=> netlist
```

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

For the spectreS simulator,

```
design( "simple.c" )  
=> simple.c
```

specifies that `simple.c`, a raw circuit file, be used in the simulation.

```
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.

```
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.

```
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`.

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 chapters [*Design Variables and Simulation Files for Direct Simulation*](#) and [*Design Variables and Simulation Files for Socket Simulation*](#) of the *Virtuoso Analog Design Environment User Guide*.

Arguments

<i>t_desVar1</i>	Name of the design variable.
<i>f_value1</i>	Value for the design variable.
<i>t_desVarN</i>	Name of an additional design variable.
<i>f_valueN</i>	Value for the additional design variable.

Value Returned

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

Examples

```
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.

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

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Sets `a` to `1k/2` or `500`.

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

envOption

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

Description

Sets environment options.

Use the OCEAN online help command `ocnHelp('envOption)` to get the list of environment options. 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 User Guide*.

Arguments

<code>s_envOption1</code>	Name of the first environment option to set.
<code>g_value1</code>	Value for the option.
<code>s_envOptionN</code>	Name of an additional environment option to set.
<code>g_valueN</code>	Value for the option.

Value Returned

<code>undefined</code>	The return value for this command/function is undefined.
<code>nil</code>	Returns <code>nil</code> if there are problems setting the option.

Examples

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

Sets the `paramRangeCheckFile` environment option.

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

Sets the `initFile` environment option.

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

Sets the `updateFile` environment option.

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 *Helping a Simulation to Converge* of the *Virtuoso Analog Design Environment User Guide*.

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

Arguments

<i>t_netName1</i>	Name of the net.
<i>f_value1</i>	Voltage value for the net.
<i>t_netNameN</i>	Name of an additional net.
<i>f_valueN</i>	Voltage value for the net.

Value Returned

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

Example

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

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

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 *Helping a Simulation to Converge* of the *Virtuoso Analog Design Environment User Guide*.

Arguments

<i>t_netName1</i>	Name of the net.
<i>f_value1</i>	Voltage value for the net.
<i>t_netNameN</i>	Name of an additional net.
<i>f_valueN</i>	Voltage value for the net.

Value Returned

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

Example

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

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

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-SPIICE 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`.

modelFile

```
modelFile( [g_modelFile1 [g_modelFile2 ...]] )  
=> l_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*.

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 *[Helping a Simulation to Converge](#)* of the *Virtuoso Analog Design Environment User Guide*.

Arguments

<i>t_netName1</i>	Name of the net.
<i>f_value1</i>	Voltage value for the net.
<i>t_netNameN</i>	Name of an additional net.
<i>f_valueN</i>	Voltage value for the net.

Value Returned

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

Example

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

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

noise

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

Description

Specifies a noise analysis.

Note: This command cannot be used with the spectre, spectreS, spectreVerilog and spectreSVerilog simulators.

Arguments

<i>t_output</i>	Output node.
<i>t_source</i>	Input source.

Value Returned

<i>undefined</i>	The return value for this command/function is undefined.
<i>nil</i>	Returns <i>nil</i> and prints an error message If there is a problem specifying the analysis.

Example

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

Specifies a noise analysis.

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 *?output* argument, the `ocnDisplay` command opens the file and writes the information to it. If you provide a port (the return value of the `SKILL outfile` command), the `ocnDisplay` command appends the information to the file that is represented by the port.

Arguments

<i>t_filename</i>	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.
<i>p_port</i>	Port (previously opened with <code>outfile</code>) through which to append the information to a file. You are responsible for closing the port. See the outfile command for more information.
<i>s_command</i>	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 <code>design</code> , <code>analysis</code> , <code>tran</code> , <code>ac</code> , <code>dc</code> , <code>noise</code> , <code>netlistDir</code> , <code>resultsDir</code> , <code>temp</code> , <code>option</code> , <code>desVar</code> , <code>path</code> , <code>includeFile</code> , <code>modelFile</code> , <code>stimulusFile</code> , <code>definitionFile</code> , <code>saveOption</code> , <code>envOption</code> , <code>keep</code> , <code>save</code> , <code>converge</code> , <code>ic</code> , <code>forcenode</code> , <code>nodeset</code> , <code>simulator</code> , <code>setup</code> , <code>restore</code> , <code>param</code> , <code>optimizeVar</code> , <code>optimizeAlgoControl</code> , <code>optimizePlotOption</code> , <code>cornerDesVar</code> , <code>monteCarlo</code> , and <code>monteOutputs</code> .
<i>g_commandArg1</i>	Argument corresponding to the specified command.

OCEAN Reference Simulation Commands

g_commandArg2 Additional argument corresponding to the specified command.

Value Returned

t Displays the information and returns *t*.

nil Returns *nil* and prints an error message if there are problems displaying the information.

Examples

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

Displays all the `optimizeGoal` information.

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

Displays only transient analyses.

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

Displays all the keeps.

```
ocnDisplay( ?output myPort 'analysis )  
=> t
```

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

ocnGetWaveformTool

```
ocnGetWaveformTool()  
=> t_toolName
```

Description

Returns the waveform tool name.

Value Returned

t_toolName The waveform tool being used.

Example

```
ocnGetWaveformTool()  
=> awd
```

Specifies that the current waveform tool is AWD.

ocnWaveformTool

```
ocnWaveformTool( s_waveformTool )  
=> t/nil
```

Description

Sets the specified tool as the waveform tool.

Arguments

s_waveformTool Either one of the waveform tools, *awd* or *wavescan*, which you want to set for the current session.

Value Returned

t Indicates that the specified waveform tool has been set.

nil Returns *nil* if there is a problem setting the specified waveform tool.

Example

```
ocnWaveformTool( 'wavescan' )  
=> t
```

Sets WaveScan as the current waveform tool.

OCEAN Reference

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

<i>s_command</i>	Command that was initially used to add the items that are now being turned off. Valid value: <code>restore</code>
<i>g_commandArg1</i>	Argument corresponding to the specified command.
<i>g_commandArg2</i>	Additional argument corresponding to the specified command.

Value Returned

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

Examples

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

Turns off the `restore` command.

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

Turns off the transient `restore` command.

OCEAN Reference 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

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

Value Returned

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

Examples

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

Sets the `abstol` option to `1e-10`.

```
option( 'delmax 50n )
```

Sets the `delmax` option to `50n`.

```
option()
```


OCEAN Reference

Simulation Commands

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.

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

<i>s_analysisType</i>	Type of the analysis. Valid values: <code>dc</code> or <code>tran</code>
<i>t_filename</i>	Name of the file containing the saved state.

Value Returned

<i>undefined</i>	The return value for this command/function is undefined.
<i>nil</i>	Returns <code>nil</code> and prints an error message if there are problems restoring the information.

Examples

```
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.

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 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/spectreS/schematic/psf")=>  
~/simulation/ckt/spectreS/schematic/psf"
```

Specifies the `psf` directory as the directory in which to store the PSF files.

```
resultsDir() => "~/simulation/ckt/spectreS/schematic/psf"
```

Returns the results directory.

OCEAN Reference Simulation Commands

run

```
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] )
    => s_jobName/nil

run( )
    =>t_dirName/nil

run(_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, parameters specific to distributed processing (such as host, job name, and queue) 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 following kinds of analyses. Instead, use the command that is specific to the analysis.

To start	Use this command
parametric analyses	<u>paramRun</u>
corners analyses	<u>cornerRun</u>
Monte Carlo analyses	<u>monteRun</u>
optimizations	<u>optimizeRun</u>

Arguments

analysisList List of analyses to be run with the `run` command.

Note: The following arguments apply only when distributed mode is enabled.

t_jobName If the name given is not unique, an integer is appended to create a unique job name.

OCEAN Reference Simulation Commands

<i>t_hostName</i>	Name of the host on which to run the analysis. If no host is specified, the system assigns the job to an available host.
<i>t_queueName</i>	Name of the queue. If no queue is defined, the analysis is placed in the default queue.
<i>t_startTime</i>	Desired start time for the job. If dependencies are specified, the job does not start until all dependencies are satisfied.
<i>t_termTime</i>	Termination time for job. If the job has not completed by the specified termination time, the job is aborted.
<i>t_dependentOn</i>	List of jobs on which the specified job is dependent. The job is not started until dependent jobs are completed.
<i>t_mailingList</i>	List of users to be notified when the analysis is complete.
<i>s_block</i>	When <i>s_block</i> is not set to <code>nil</code> , the OCEAN script halts until the job is complete. Default value: <code>nil</code>
<i>s_notifyFlag</i>	When not set to <code>nil</code> , the job completion message is echoed to the OCEAN interactive window. Default value: <code>t</code>
<i>s_analysisType1</i>	Name of a prespecified analysis to be simulated.
<i>s_analysisTypeN</i>	Name of another prespecified analysis to be simulated.

Value Returned

<i>s_jobName</i>	Returns the job name of the job submitted. The job name is based on the <code>jobName</code> 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.
<i>t_dirName</i>	Returns the name of the directory in which the results are stored. This value is returned for local and blocking distributed modes.

OCEAN Reference Simulation Commands

`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.)

Examples

```
run( )  
=> t
```

Starts the simulation.

```
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 actual 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.

OCEAN Reference

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 `saveName`, you can include as many outputs as you want. If you want to turn off the default of `save`, 'allv, use the `delete('save)` command.

Arguments

s_categ Type of simulator to be used.
Valid values: analog, digital
Default value: analog
Note: digital is not available.

s_saveType Type of outputs to be saved.
Valid values:

Valid Values	Description
v	Specifies that a list of subsequent net names be kept.
i	Specifies that a list of subsequent currents be kept.
all	Specifies that all nets and all currents are to be saved.
allv	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.

OCEAN Reference Simulation Commands

Value Returned

<i>undefined</i>	The return value for this command/function is undefined.
<i>nil</i>	Returns <i>nil</i> and prints an error message if there is a problem keeping the outputs.

Examples

```
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" "mnl:d" )
```

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(mnl,d)" )
```

For the Cadence-SPICE circuit simulator, saves the current through the same devices.

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 [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 “[save](#)” on page 111.

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.)

OCEAN Reference Simulation Commands

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 ?savehdlvars "all")
```

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.

Examples

```
simulator( 'spectre )  
=> spectre
```

Specifies that the Spectre® circuit simulator be used for the session.

```
simulator( 'spectreVerilog )  
=> spectreVerilog
```

Specifies that spectreVerilog be used for the session.

```
simulator()  
=> spectreVerilog
```

Returns the simulator that you set for the session. If a simulator was not specified, it returns nil.

stimulusFile

```
stimulusFile( t_fileName [t_fileName2 ... t_fileNameN ] [?xlate b_xlate] )  
=> l_fileNames/nil
```

Description

Specifies stimulus files to be used by the simulator.

When the *b_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

<i>t_fileName</i>	The name of the stimulus file to be included.
<i>t_fileName2...t_fileNameN</i>	The names of the additional stimulus files to be included.
<i>b_xlate</i>	If set to <i>t</i> , net and instance expressions are translated to simulator names. The default value of the <i>b_xlate</i> variable is <i>t</i> .

Value Returned

<i>l_fileNames</i>	A list of the stimulus file names is the output if the command is successful.
<i>nil</i>	Otherwise <i>nil</i> 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.

```
stimulusFile()  
=> ( "tran.stimulus" "rf.stimulus" )
```

OCEAN Reference Simulation Commands

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

OCEAN Reference 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

<i>s_analysisType</i>	Type of the analysis. Valid values: <code>dc</code> or <code>tran</code>
<i>t_filename</i>	Name of the file in which to store the simulator's node voltages.

Value Returned

<i>t_filename</i>	Returns the filename.
<code>nil</code>	Returns <code>nil</code> and prints an error message if there are problems storing the information to a file.

Examples

```
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 `tranStoreFile` in the `netlist` (design) directory unless a full path is specified.

temp

```
temp( f_tempValue )  
=> f_tempValue/nil
```

Description

Specifies the circuit temperature.

Arguments

f_tempValue Temperature for the circuit.

Value Returned

f_tempValue Returns the temperature specified.

nil Returns *nil* and prints an error message if there are problems setting the temperature.

Example

```
temp( 125 )  
=> 125
```

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.

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 [analysis](#) 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`, `spectreS`, `spectreVerilog` and `spectreSVerilog` simulators.

Arguments

<i>g_fromValue</i>	Starting time for the analysis.
<i>g_toValue</i>	Ending time.
<i>g_byValue</i>	Increment at which to step through the analysis.

Value Returned

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

Examples

```
tran( 1u )  
=> "1e-06"
```

Specifies a transient analysis to 1u for the Spectre® circuit simulator

```
tran( 0 1u 1n )  
=> "1e-09"
```

Specifies a transient analysis from 0 to 1u by increments of 1n.

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

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Data Access Commands

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dataTypes

```
dataTypes()  
=> l_dataTypes/nil
```

Description

Returns the list of data types that are used in an analysis previously specified with `selectResult`.

Arguments

None.

Value Returned

<i>l_dataTypes</i>	Returns the list of data types.
<i>nil</i>	Returns <i>nil</i> and an error message if the list of datatypes cannot be returned.

Example

```
selectResult( 'dcOp '  
dataTypes() => ( "node" "vs" "resistor" "bjt" )
```

Returns the data types used in the selected file, in this case, `dcOp`.

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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.

Arguments

<i>t_name</i>	Name of the signal.
<i>s_resultName</i>	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.
<i>t_resultsDir</i>	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

<i>x_number</i>	Returns an integer simulation result.
<i>o_waveform</i>	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: drwave:XXXXX.)
<i>nil</i>	Returns <i>nil</i> and an error message if the value cannot be returned.

Examples

```
getData( "/net6" ) => drwave:25178234
```

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Data Access Commands

Returns the number or waveform for `net6`. In this example, the return value is equivalent to `v("/net6")`.

```
getData( "/V1" ?result 'ac )
=> drwave: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" ) ) =>
```

The `getData("net1")` command passes a waveform to the `ocnPrint` command. The `ocnPrint` 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`.

OCEAN Reference

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

<i>s_resultName</i>	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 <code>selectResult</code> command. The default is the current result selected with the <code>selectResult</code> command.
<i>t_resultsDir</i>	Directory containing the PSF files (results). If you supply this argument, you must also supply the <code>resultName</code> argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the <code>openResults</code> command. The default is the current results directory set by the <code>openResults</code> command.

Value Returned

<i>o_results</i>	Returns the object representing the selected results.
<i>nil</i>	Returns <code>nil</code> and an error message if there are problems accessing the analysis.

Example

```
getResult( ?result 'tran )
```

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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

<i>t_component</i>	Name of the component.
<i>s_resultName</i>	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 <code>selectResult</code> command. The default is the current result selected with the <code>selectResult</code> command.
<i>t_resultsDir</i>	Directory containing the PSF files (results). If you supply this argument, you must also supply the <code>resultName</code> argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the <code>openResults</code> command. The default is the current results directory set by the <code>openResults</code> command.

Value Returned

<i>o_waveform</i>	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: <code>drwave:XXXXX</code>).
<i>nil</i>	Returns an error message and <code>nil</code> if there is a problem.

Examples

```
selectResult( 'tran' )  
i( "/R1" )
```

Returns the current through the R1 component.

```
ocnPrint( i( "/R1" ) )
```

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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).

OCEAN Reference

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 `?output` argument, the `ocnHelp` command opens the file and writes the information to it. If you provide a port (the return value of the `SKILL outfile` command), the `ocnHelp` command appends the information to the file that is represented by the port. If you do not specify `?output`, the output goes to standard out (stdout).

Arguments

<i>t_filename</i>	File in which to write the information. The <code>ocnHelp</code> command opens the file, writes to the file, and closes the file. If you specify the filename without a path, the <code>ocnHelp</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.
<i>p_port</i>	Port (previously opened with <code>outfile</code>) through which to append the information to a file. You are responsible for closing the port. See the outfile command for more information.
<i>s_command</i>	Command for which you want help.

Value Returned

<i>t</i>	Displays the online help and returns <i>t</i> .
<i>nil</i>	Returns <i>nil</i> and an error message if help cannot be displayed.

Examples

```
ocnHelp()  
=> t
```

Displays information about using online help.

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Data Access Commands

```
ocnHelp( 'analysis ' )  
=> t
```

Displays help for the `analysis` command.

```
ocnHelp( ?output "helpInfo" )  
=> t
```

Writes information about using online help to a file named `helpInfo`.

OCEAN Reference

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.

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 an ADE data directory; it must have the `psf` directory under it and the `psf` directory must contain `runObjFile`.

OCEAN Reference

Data Access Commands

Value Returned

<code>t_dirName</code>	The directory containing the PSF files.
<code>nil</code>	Returns <code>nil</code> and an error message if there are problems opening the results.

Examples

```
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.

OCEAN Reference

Data Access Commands

outputParams

```
outputParams( t_compType [?result s_resultName [?resultsDir t_resultsDir]] )  
=> l_outputParams/nil
```

Description

Returns the list of output parameters for the specified component.

You can use the [dataTypes](#) 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 [pv](#) command.

Arguments

<i>t_compType</i>	Name of a component.
<i>s_resultName</i>	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 <code>selectResult</code> command. The default is the current result selected with the <code>selectResult</code> command.
<i>t_resultsDir</i>	Directory containing the PSF files (results). If you supply this argument, you must also supply the <code>resultName</code> argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the <code>openResults</code> command. The default is the current results directory set by the <code>openResults</code> command.

Value Returned

<i>l_outputParams</i>	Returns the list of parameters.
<i>nil</i>	Returns <code>nil</code> and an error message if there are no associated parameters or if the specified component (<code>compType</code>) does not exist.

OCEAN Reference

Data Access Commands

Example

```
selectResult( 'dcOp )
dataTypes() => ( "node" "vs" "resistor" "bjt" )
outputParams( "bjt" )
```

Selects the `dcOp` results, returns the list of components for these results, and returns the list of output parameters for the `bjt` component.

```
outputParams("bjt" ?result 'dcOp ?resultsDir "./psf")
```

Returns a list of output parameters for the `bjt` component for `dcOp` (dc analysis with save dc operating point) results stored at the location `./psf`.

outputs

```
outputs( [?result s_resultName [?resultsDir t_resultsDir]]  
        [?type t_signalType])  
=> l_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.

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.

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" )
```

OCEAN Reference

Data Access Commands

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.

OCEAN Reference

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.

Arguments

g_harmonic List of harmonic frequencies.

S_signalResultName Name of the result that stores the signal waveform. Use the `results()` command to obtain the list results.

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.

OCEAN Reference Data Access Commands

Example

```
plot(phaseNoise(0 "pss-fd.pss"))  
phaseNoise(1 "pss_fd" ?result "pnoise" ?resultsDir "./PSF")
```

OCEAN Reference

Data Access Commands

pv

```
pv( t_name t_param [?result s_resultName [?resultsDir t_resultsDir]] )  
    => g_value/nil
```

Description

Returns the value for the specified component parameter. You can use the [outputParams](#) command to get the list of parameters for a particular component.

Arguments

<i>t_name</i>	Name of the node or component.
<i>t_param</i>	Name of the parameter.
<i>s_resultName</i>	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 <code>selectResult</code> command. The default is the current result selected with the <code>selectResult</code> command.
<i>t_resultsDir</i>	Directory containing the PSF files (results). If you supply this argument, you must also supply the <code>resultName</code> argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the <code>openResults</code> command. The default is the current results directory set by the <code>openResults</code> command.

Value Returned

<i>g_value</i>	Returns the requested parameter value.
<i>nil</i>	Returns <code>nil</code> and prints an error message.

Examples

```
selectResult( 'dcOp' )  
pv( "/Q19" "ib" )
```

For the `Q19` component, returns the value of the `ib` parameter.

```
pv( "/Q19" "ib" ?resultsDir "./test2/psf" )
```

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Data Access Commands

For the Q19 component, returns the value of the `ib` parameter for the results from a different run (stored in `test2/psf`).

```
pv( "/Q19" "ib" ?result "dcOp" ?resultDir "./test1/psf" )
```

Returns the value of the `ib` parameter for the Q19 component for the `dcOp` results stored at the location `./test1/psf`.

OCEAN Reference

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

<i>s_propertyName</i>	Name of the parameter
<i>s_resultName</i>	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.
<i>t_resultsDir</i>	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

<i>L_value</i>	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.

Examples

```
resultParam("positive output signal" ?result "pnoise.pss")  
=> "pif"  
resultParam("negative output signal" ?result "pnoise.pss")  
=> "0"
```

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")
```

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Data Access Commands

```
=> 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.

results

```
results( [ ?resultsDir t_resultsDir ] )  
=> l_results/nil
```

Description

Returns a list of the type of results that can be selected.

Arguments

<i>t_resultsDir</i>	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

<i>l_results</i>	Returns the list of result types.
<i>nil</i>	Returns <i>nil</i> and an error message if there are problems returning the results.

Example

```
results()  
=> ( dc tran ac )
```

Returns the list of results available.

```
results("./psf")
```

Returns a list of results stored at the location `./psf`.

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 [results](#) command to get the list of analysis results that you can select.

Arguments

<i>s_resultsName</i>	Results from an analysis.
<i>n_sweepValue</i>	The sweep value you wish to select for an analysis.

Value Returned

<i>o_results</i>	Returns the object representing the selected results.
<i>nil</i>	Returns <code>nil</code> and an error message if there are problems selecting the analysis.

Examples

```
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 )
```

Selects the results for a transient analysis.

```
paramAnalysis("supply" ?start 3 ?stop 5 ?step 1.0/3)
```


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Data Access Commands

```
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.

OCEAN Reference

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

<i>x_iIndex</i>	The <i>i</i> th index of the coefficient in the scattering matrix.
<i>x_jIndex</i>	The <i>j</i> th index of the coefficient in the scattering matrix.
<i>s_resultName</i>	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.
<i>t_resultsDir</i>	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

<i>o_waveform</i>	Waveform object representing the S-parameter.
<i>nil</i>	Returns <i>nil</i> if there is an error.

Examples

```
s21 = sp(2 1)  
s12 = sp(1 2)  
plot(s21 s12)
```

```
s11 = sp(1 1 ?result "sp" ?resultsDir "./simResult/psf")
```

OCEAN Reference

Data Access Commands

Returns the S-parameter s_{11} for results of S-parameter(sp) analysis stored at the location `./simResult/psf`.

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Data Access Commands

sweepNames

```
sweepNames( [o_waveForm] [?result s_resultName [?resultsDir t_resultsDir]] )  
=> l_sweepName/nil
```

Description

Returns the names of all the sweep variables for either a supplied waveform, a currently selected result (via `selectResult()`) or a specified result.

Arguments

<i>o_waveForm</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code>). When this argument is used, the <code>t_resultsDir</code> and <code>s_resultName</code> arguments are ignored.
<i>s_resultName</i>	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 <code>selectResult</code> command. The default is the current result selected with the <code>selectResult</code> command.
<i>t_resultsDir</i>	Directory containing the PSF files (results). If you supply this argument, you must also supply the <code>resultName</code> argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the <code>openResults</code> command. The default is the current results directory set by the <code>openResults</code> command.

Value Returned

<i>l_sweepName</i>	Returns a list of the sweep names.
<code>nil</code>	Returns <code>nil</code> and prints an error message if the sweep names cannot be returned.

Example

```
selectResult('tran)  
sweepNames()  
=> ( "TEMPDC" "time" )
```

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Data Access Commands

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")

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 `./test/psf`.

sweepValues

```
sweepValues( [o_waveForm] )  
=> l_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

<i>o_waveForm</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
-------------------	--

Value Returned

<i>l_sweepValues</i>	Returns the list of sweep values.
<i>nil</i>	Returns <i>nil</i> 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 *w*. In this case, the return values indicate the resistance over which the data was swept.

sweepVarValues

```
sweepVarValues( [t_varName] [?result s_resultName [?resultsDir t_resultsDir]]  
=> l_sweepName/nil
```

Description

Returns the list of sweep values for a particular swept variable name. This command is particularly useful for parametric analyses.

Arguments

<i>t_varName</i>	Name of the specific variable from which the values are retrieved.
<i>s_resultName</i>	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.
<i>t_resultsDir</i>	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

<i>l_sweepValues</i>	Returns the list of sweep values.
<i>nil</i>	Returns nil and an error message if the list of sweep values cannot be returned.

Examples

```
selectResult('tran)  
sweepNames()  
=> ("TEMPDC" "Vsupply" "time")  
sweepVarValues("TEMPDC")  
=> (0 32)
```

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Data Access Commands

```
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)
```


OCEAN Reference

Data Access Commands

V

```
v( t_net [?result s_resultName [?resultsDir t_resultsDir]] )  
    => o_waveform/nil
```

Description

Returns the voltage of the specified net.

Arguments

<i>t_net</i>	Name of the net.
<i>s_resultName</i>	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.
<i>t_resultsDir</i>	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

<i>o_waveform</i>	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: <code>drwave:XXXXX</code>).
<i>nil</i>	Returns an error message and <code>nil</code> if there is a problem.

Example

```
selectResult('tran)  
v( "/net56" )
```

Returns the voltage for net56.

```
ocnPrint( v( "/net56" ) )
```

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Data Access Commands

Prints tabular information representing the voltage for `net56`.

```
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`).

OCEAN Reference

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.

Arguments

<i>x_index</i>	Index of the port.
<i>s_resultName</i>	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.
<i>t_resultsDir</i>	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

<i>o_waveform</i>	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")
```

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Data Access Commands

Returns the voltage standing wave ratio value at port 1 for the results of S-parameter(sp) analysis stored at the location `./simResult/psf`.

OCEAN Reference

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

```
(1 + s( x_index x_index )) / (1 - s( x_index x_index ))  
    * or( zref( x_index ) 50)
```

This command should be run on the results of the Spectre `sp` (S-parameter) analysis.

Arguments

<code>x_index</code>	Index of the port.
<code>s_resultName</code>	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 <code>selectResult</code> command. The default is the current result selected with the <code>selectResult</code> command.
<code>t_resultsDir</code>	Directory containing the PSF files (results). If you supply this argument, you must also supply the <code>resultName</code> argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the <code>openResults</code> command. The default is the current results directory set by the <code>openResults</code> command.

Value Returned

<code>o_waveform</code>	Waveform object representing the port input impedance.
<code>nil</code>	Returns an error message and <code>nil</code> if there is a problem.

Example

```
plot(zm(2))  
zm1 = zm(1 ?result "sp" ?resultsDir "./simResult/psf")
```

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Data Access Commands

Returns input impedance at port 1 for results of S-parameter (sp) analysis stored at the location `./simResult/psf`.

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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

<i>x_portIndex</i>	Index of the port.
<i>s_resultName</i>	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.
<i>t_resultsDir</i>	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

<i>f_impedance</i>	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")
```

Returns the reference impedance at port 1 for the results of S-parameter(sp) analysis stored at the location ./simResult/psf.

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Plotting and Printing Commands

This chapter contains information on the following plotting and printing commands:

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This chapter also includes a topic, [Plotting and Printing SpectreRF Functions in OCEAN](#) on page 212.

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

<i>x_subwindowID</i>	Returns the window ID of the new subwindow.
<i>nil</i>	Returns <i>nil</i> and an error message if there is no current Waveform window.

Example

```
addSubwindow()  
=>3
```

Adds a new subwindow to the Waveform window.

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" )  
=> t
```

Adds the title `waveform 2` to the selected subwindow.

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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.

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Plotting and Printing Commands

addWaveLabel

```
addWaveLabel( x_waveIndex l_location t_label [?textOffset l_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.

Arguments

<i>x_waveIndex</i>	Integer identifying the waveform curve.
<i>l_location</i>	List of two waveform coordinates that describe the location for the label.
<i>t_label</i>	Label for the waveform.
<i>l_textOffset</i>	An offset of the label from <i>l_location</i> , in screen units of the current subwindow. If <i>l_textOffset</i> is not specified, it defaults to 0:0 and the label is displayed at the location. If <i>l_textOffset</i> is specified, the label is offset from the location and a directional arrow is drawn from the label to the location. For example, if the offset is specified as 0:20, the label is drawn 20 units above the location and a directional label is drawn from the label to the location. This feature is useful to label points on a waveform and not obstruct the waveform.
<i>x_color</i>	Label color specified as an index in the technology file. Default value: 10
<i>t_justify</i>	Justification, which is specified as "upperLeft", "centerLeft", "lowerLeft", "upperCenter", "centerCenter", "lowerCenter", "upperRight", "centerRight", or "lowerRight". Default value: "lowerLeft"
<i>t_fontStyle</i>	Font style, which is specified as "euroStyle", "gothic", "math", "roman", "script", "stick", "fixed",

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	"swedish", "raster", or "milSpec". Default value: the font style of the current subwindow
<i>x_height</i>	Height of the font. Default value: the font height of the current subwindow
<i>t_orient</i>	Orientation of the text, specified as either "R0" or "R90". Default value: "R0"
<i>g_drafting</i>	Boolean that specifies whether the label stays backwards or upside-down. If set to <i>t</i> , a backwards or upside-down label is displayed in a readable form. If set to <i>nil</i> , a backwards or upside-down label stays the way it is. Default value: <i>t</i>
<i>g_overbar</i>	Boolean that specifies whether underscores in labels are displayed as overbars. If set to <i>t</i> , underscores in labels are displayed as overbars. If set to <i>nil</i> , underbars are displayed as underbars. Default value: <i>nil</i>

Value Returned

<i>s_labelId</i>	Returns an identification number for the waveform label.
<i>nil</i>	Returns <i>nil</i> if there is an error.

Example

```
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 – 10, and Orientation – *R20*.

The label will be displayed in a readable form. The underscore in the label will be displayed as an overbar.

addWindowLabel

```
addWindowLabel( l_location t_label )  
=> s_labelId/nil
```

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

<i>l_location</i>	List of two waveform coordinates that describe the location for the label. Valid values: 0 through 1
<i>t_label</i>	Label for the waveform.

Value Returned

<i>s_labelId</i>	Returns an identification number for the subwindow label.
<i>nil</i>	Returns <i>nil</i> 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`.

clearAll

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

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.

clearSubwindow

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

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()  
=> t
```

Erases the contents of the current subwindow.

currentSubwindow

```
currentSubwindow( x_subwindow )  
=> t/nil
```

Description

Specifies *x_subwindow* as the current subwindow.

Arguments

<i>x_subwindow</i>	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 <i>x_subwindow</i> .
---	---

nil	Returns nil if there is an error.
-----	-----------------------------------

Example

```
currentSubwindow( 2 )
```

Specifies subwindow 2 as the current subwindow.

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) )
```

Specifies window 2 as the current Waveform window.

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Plotting and Printing Commands

dbCompressionPlot

```
dbCompressionPlot(o_wave x_harmonic x_extrapolationPoint  
  [?compression x_compression] )  
=> t/nil
```

Description

Plots the *n*th 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

<i>o_wave</i>	The waveform for which to plot the compression.
<i>x_harmonic</i>	Harmonic frequency index.
<i>x_extrapolationPoint</i>	The extrapolation point.
<i>x_compression</i>	The amount of dB compression. Default value: 1

Value Returned

<i>t</i>	Returns <i>t</i> if the point is plotted
<i>nil</i>	returns <i>nil</i> if there was an error

Example

```
dbCompressionPlot(v("/Pif") 2 -25)
```

Plots a 1 dB compression point plot for the waveform `v("/Pif")`.

```
dbCompressionPlot(v("/Pif") 2 -25 ?compression 3)
```

Plots a 3 dB compression point plot for the waveform `v("/Pif")`.

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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 `openResults` command. The `dcmatchSummary` command prints the results for that directory and resets the `openResults` command to its previous setting. If you specify a particular result with `resultName`, it is equivalent to temporarily using the `selectResult` command on the specified results. The `dcmatchSummary` command prints the results and resets the `selectResult` command to its previous setting.

This command should be run on the results of the Spectre `dcmatch` analysis.

Arguments

<i>t_resultsDir</i>	The directory containing the <code>dcmatch</code> -analysis results.
<i>S_resultName</i>	Results from an analysis for which you want to print the <code>dcmatchSummary</code> report.
<i>t_fileName</i>	File in which to write the information. The <code>dcmatchSummary</code> command opens the file, writes to the file and closes the file. If you specify the filename without a path, the <code>dcmatchSummary</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.
<i>p_port</i>	Port (previously opened with <code>outfile</code>) through which to append the information to a file. You are responsible for closing the port. See the <code>outfile</code> command for more information.
<i>ln_paramValues</i>	List of values for swept parameters at which the <code>dcmatchSummary</code> is to be printed. In case there is just one swept parameter the value can be specified as is.

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<i>ls_deviceType</i>	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.
<i>ls_variations</i>	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"))
<i>lt_includeInst</i>	List of instance name strings to definitely include in the dcmatchSummary.
<i>lt_excludeInst</i>	List of instance name strings to exclude in the dcmatchSummary.
<i>x_truncateData</i>	Specifies a number that the truncateType argument uses to define the components for which information is to be printed.
<i>s_truncateType</i>	Specifies the method that is used to limit the data being included in the report

Valid Values	Description	Sample Values for truncateData
'top	Saves information for the number of components specified with truncateData. The components with the highest contributions are saved.	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	1.9n
'absolute	Saves information for all the components in the selected set whose contribution are more than truncateData.	0.1

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'none	Saves information for all the components.	Not required
-------	---	--------------

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 message if the summary cannot be printed.

Examples

```
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).

```
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 )
```


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Prints a report for swept DC-Mismatch analysis at swept parameter value of 25 sorted on name.

deleteSubwindow

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

Description

Deletes the current subwindow from the current Waveform window.

Arguments

None.

Value Returned

<code>t</code>	Returns <code>t</code> if the current subwindow is deleted.
<code>nil</code>	Returns <code>nil</code> and an error message if there is no current subwindow.

Example

```
deleteSubwindow()  
=> t
```

Deletes the current subwindow from the Waveform window.

deleteWaveform

```
deleteWaveform( {x_index | all_string } )  
=> t/nil
```

Description

Deletes the specified waveform curve or all the waveform curves from the current subwindow of a Waveform window.

Arguments

<i>x_index</i>	Integer identifying a particular waveform curve.
<i>all_string</i>	The string "all" specifying that all waveform curves are to be deleted.

Value Returned

<i>t</i>	Returns <i>t</i> if the curves are deleted.
<i>nil</i>	Returns <i>nil</i> and an error message if the curves are not deleted.

Examples

```
deleteWaveform( '1' )  
=> t
```

Deletes waveform 1 from the current subwindow.

```
deleteWaveform( "all" )  
=> t
```

Deletes all the curves from the current subwindow.

displayMode

```
displayMode( t_mode )  
=> t/nil
```

Description

Sets the display mode of the current subwindow.

Arguments

<i>t_mode</i>	String representing the display mode for the subwindow. Valid values: <i>strip</i> , <i>smith</i> , or <i>composite</i>
---------------	--

Value Returned

<i>t</i>	Returns <i>t</i> when the display mode of the subwindow is set.
----------	---

<i>nil</i>	Returns <i>nil</i> and an error message if the display mode cannot be set.
------------	--

Example

```
displayMode( "composite" )  
=> t
```

Sets the current subwindow to display in *composite* mode.

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Plotting and Printing Commands

getAsciiWave

```
getAsciiWave( t_filename x_xColumn x_yColumn [x_xskip] [x_yskip])  
=> 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.

Arguments

<i>t_filename</i>	The name of the Ascii file to be read in.
<i>x_xColumn</i>	The column in the data file that contains the X-axis data.
<i>x_yColumn</i>	The column in the data file that contains the Y-axis data.
<i>x_xskip</i>	The number of lines to skip in the X column.
<i>x_yskip</i>	The number of lines to skip in the Y column.

Value Returned

<i>o_wave</i>	The DRL waveform object
<i>nil</i>	Returns <i>nil</i> if the function fails.

Example

```
getAsciiWave("~/mydatafile.txt " 1 2 )  
=> drwave: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 DRL waveform object.

```
getAsciiWave("~/mydatafile.txt " 1 2 ?xskip 1 ?yskip 2)  
=> drwave: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 `xcolumn` and 2 lines in the `ycolumn`, and returns a DRL waveform object.

graphicsOff

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

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

<code>t</code>	Returns <code>t</code> if redrawing is disabled.
<code>nil</code>	Returns <code>nil</code> if there is an error, such as there is no current Waveform window.

Example

```
graphicsOff()  
=> t
```

Disables the redrawing of the Waveform window.

graphicsOn

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

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.

hardCopy

```
hardCopy(w_windowId)  
=> t/nil
```

Description

Sends a Waveform window plot to a printer.

Note: You must first set any plotting options with the [hardCopyOptions](#) command.

Arguments

<i>w_windowId</i>	The window ID of the waveform window whose plot is to be printed. The default value is the window ID of the current window.
-------------------	---

Value Returned

<i>t</i>	Returns <i>t</i> if successful.
<i>nil</i>	Returns <i>nil</i> if there is an error.

Example

```
hardCopy()  
=> t
```

Sends a waveform plot to the printer.

```
w = newWindow()  
plot(v("/vout"))  
hardCopy(w)
```

Sends the waveform plot of *w* to the printer.

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hardCopyOptions

```
hardCopyOptions( [?hcNumCopy x_hcNumCopy] [?hcDisplay t_hcDisplay]  
  [?hcOrientation s_hcOrientation] [?hcOutputFile g_hcOutputFile]  
  [?hcPaperSize t_hcPaperSize] [?hcPlotterName t_hcPlotterName]  
  [?hcTmpDir t_hcTmpDir] )  
=> g_value/nil
```

Description

Sets Waveform window hardcopy plotting options.

The option takes effect for any Waveform window or subwindow that is opened after the option is set.

Arguments

<i>x_hcNumCopy</i>	The number of copies to plot. Valid values: any positive integer Default value: 1
<i>t_hcDisplay</i>	The display name. Valid values: defined in the technology file Default value: "display"
<i>s_hcOrientation</i>	The plot orientation. Valid values: 'portrait, 'landscape, 'automatic Default value: 'automatic
<i>g_hcOutputFile</i>	Name of the output file. Valid values: a string or nil Default value: nil
<i>t_hcPaperSize</i>	The plot paper size. Valid values: specified in .cdsplotinit Default value: specified in .cdsplotinit
<i>t_hcPlotterName</i>	The name of the plotter.

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Valid values: specified in `.cdsplotinit`
Default value: specified in `.cdsplotinit`

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"`

Value Returned

g_value

Returns the new value of the option.

nil

Returns `nil` if there is an error.

Examples

```
hardCopyOptions( ?hcNumCopy 1 )
```

Plots one copy of the window or subwindow.

```
hardCopyOptions(?hcNumCopy 3 ?hcOutputFile "myOutFile")
```

Plots three copies of the window or subwindow and sends them to the file `myOutFile`.

```
hardCopyOptions(?hcNumCopy 2 ?hcOrientation 'portrait ?hcOutputFile "myOutfile")
```

Plots 2 copies of the window in `portrait` orientation and sends them to the file `myOutFile`.

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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 chapter [*Simulating Mixers*](#) of the *SpectreRF User Guide* for more information on ip3Plot.

Arguments

<i>o_wave</i>	Waveform for which to plot the ip3.
<i>x_sigHarmonic</i>	Index of the third order harmonic.
<i>x_refHarmonic</i>	Index of the first order (fundamental) harmonic.
<i>x_extrapolationPoint</i>	Extrapolation point.

Value Returned

<i>t</i>	Returns <i>t</i> if the curves are plotted.
<i>nil</i>	Returns <i>nil</i> if there is an error.

Example

```
ip3Plot(v("/net28") 47 45 -25)
```

newWindow

```
newWindow()  
=> w_windowID/nil
```

Description

Creates a new Waveform window and returns the window ID.

Arguments

None.

Value Returned

<i>w_windowId</i>	Returns the window ID of the new Waveform window.
<i>nil</i>	Returns <i>nil</i> 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.

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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])  
=> t_fileName/p_port/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.

Arguments

<i>s_type</i>	Type of noise-analysis results for which to print the report. Valid values: <i>spot</i> , to specify noise at a particular frequency, or <i>integrated</i> , to specify noise integrated over a frequency range.
<i>s_resultName</i>	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 <i>selectResult</i> command. The default is the current result selected with the <i>selectResult</i> command.
<i>t_resultsDir</i>	Directory containing the PSF files (results). If you supply this argument, you must also supply the <i>resultName</i> argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the <i>openResults</i> command. The default is the current results directory set by the <i>openResults</i> command.
<i>f_frequency</i>	Frequency value of interest.
<i>f_weight</i>	Waveform representing the function with which the integral is weighted. Default value: 1.0
<i>t_fileName</i>	File in which to write the information. The <i>noiseSummary</i> command opens the file, writes to the file, and closes the file. If

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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.

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 `outfile` command for more information.

t_noiseUnit

Specifies the type of noise unit to be saved.
Valid values: "V^2" for V²/Hz or
"V" for V/sqrt(Hz)

x_truncateData

Specifies a number that the `truncateType` argument uses to define the components for which information is to be printed.

s_truncateType

Specifies the method that is used to limit the data being included in the report.

Valid Values	Description	Sample Values
'top	Saves information for the number of components specified with <code>truncateData</code> . The components with the highest contributions are saved.	10
'level	Prints components which have noise contribution higher than that specified by <code>?truncateData</code> .	10u
'relative	Prints components which have noise contribution (percent) higher than that specified by <code>?truncateData</code> .	.1
'none	Saves information for all the components.	

x_digits

Number of significant digits with which the contributors are printed.

x_percentDecimals

Number of decimals printed for any relative contribution.

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<i>f_from</i>	For integrated noise, the start value for frequency.
<i>f_to</i>	For integrated noise, the end value for frequency.
<i>ls_deviceType</i>	List of device type strings to be included. Valid values: a list of strings or 'all
<i>t_weightFile</i>	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, <i>weight</i> and <i>weightFile</i> , the value for <i>weight</i> gets precedence..

Value Returned

<i>t_fileName</i>	Returns the name of the port.
<i>p_port</i>	Returns the name of the file.
nil	Returns nil and an error message if the summary cannot be printed.

Examples

```
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"))
```

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Prints a report for an integrated noise analysis in the frequency range 10-10M for 20 components with deviceType `bjt`, `mos` or `resistor`.

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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] 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 `?output` argument, the `ocnPrint` command opens the file and writes the information to it. If you provide a port (the return value of the `SKILL outfile` command), the `ocnPrint` command appends the information to the file that is represented by the port. There is a limitation of `ocnPrint` for precision. It works upto 30 digits for the Solaris port and 18 digits for HP and AIX.

Arguments

<i>t_filename</i>	File in which to write the information. The <code>ocnPrint</code> 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 <code>getSkillPath()</code> at the OCEAN prompt.
<i>p_port</i>	Port (previously opened with <code>outfile</code>) through which to append the information to a file. You are responsible for closing the port. See the outfile command for more information.
<i>x_precision</i>	The number of significant digits to print. This value overrides any global precision value set with the <code>setup</code> command. Valid values: 1 through 16 Default value: 6
<i>s_numberNotation</i>	The notation for printed information. This value overrides any global format value set with the <code>setup</code> command. Valid values: 'suffix, 'engineering, 'scientific, 'none Default value: 'suffix

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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.

x_numSpaces

The number of spaces between columns.
Valid values: 1 or greater
Default value: 4

x_width

The width of each column.
Valid values: 4 or greater
Default value: 14

x_from

The start value at x axis for the waveform to be printed.

x_to

The end value at x axis for the waveform to be printed.

?step

The step by which text data to be printed is incremented.

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: drwave: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.

Examples

```
ocnPrint( v( "/net56" ) )  
=> t
```

Prints the text for the waveform for the voltage of net56.

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```
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`.

```
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.

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ocnYvsYPlot

```
ocnYvsYPlot([?wavex o_wavex ?wavey o_wavey] [?exprx o_exprx ?expy o_expy]  
            [?title l_titleList] [?color 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 paramset, Corners or MonteCarlo results data.

Arguments

<i>o_wavex</i>	Reference wave against which the wave provided needs to be plotted.
<i>o_wavey</i>	Wave to be plotted against the reference wave.
<i>o_exprx</i>	Reference expression against which the expression provided needs to be plotted.
<i>o_expy</i>	Expression to be plotted against the reference expression.
<i>l_titleList</i>	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.
<i>l_colorList</i>	List of waveform colors. If the waveform is simple, only one color will be required. If the waveform is param, a list of colors needs to be provided.

Value Returned

<i>wave</i>	Returns the waveform specified.
<i>nil</i>	Returns <i>nil</i> if the plot could not be generated.

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Examples

```
wy = VT("/vout")
wx = VT("/vin")
ex = "VT('/vin')"
ey = "VT('/vout')"
ocnYvsYplot(?wavex wx ?wavey wy ?titleList '("simpleWave") ?colorList '(3))
```

Plots wave wy against wave wx with the title being simpleWave and the color being 3.

```
ocnYvsYplot(?exprx ex ?expy ey ?titleList '("simpleWave") ?colorList '(3))
```

Plots expression ey against expression ex with the title being simpleWave and the color being 3.

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Plotting and Printing Commands

plot

```
plot( o_waveform1 [o_waveform2 ...] [?yNumber l_yNumberList] [?expr l_exprList]
      )
      => t/nil
```

Description

Plots waveforms in the current subwindow. If there is no Waveform window, this command opens one.

Arguments

<i>o_waveform1</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>o_waveform2</i>	Additional waveform object.
<i>l_yNumberList</i>	List that specifies the Y axes where the waveforms are to be plotted. The number of Y axes must match the number of waveform objects specified. Valid values: 1, 2, 3, and 4
<i>l_exprList</i>	List of strings used to give names to the waveform objects.

Value Returned

<i>t</i>	Returns <i>t</i> if the waveforms are plotted.
<i>nil</i>	Returns <i>nil</i> and an error message if the waveforms cannot be plotted.

Examples

```
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`.

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```
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" ) ?yNumber list( 1 2 ) )
```

Plots the waveforms `v("OUT")` and `i("VFB")` on the Y axes 1 and 2, respectively.

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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`.

Arguments

S_style Plotting style for the subwindow.
Valid values: `auto`, `scatterplot`, `bar`, `joined`

Argument	Description
<code>auto</code>	The appropriate plotting style is automatically chosen.
<code>scatterplot</code>	Data points are not joined.
<code>bar</code>	Vertical bars are drawn at each data point that extend from the point to the bottom of the graph.
<code>joined</code>	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

```
plotStyle( 'auto' )  
=> t
```

Sets the plot style to `auto`.

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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.

Arguments

<i>t_resultsDir</i>	Directory containing the results. If you specify a directory with <i>resultsDir</i> , the <i>pzPlot</i> command plots the results for that directory.
<i>S_resultName</i>	Pointer to results from the analysis for which you want to plot the report.
<i>S_toPlot</i>	Use this option to plot only poles, only zeros or both poles and zeros information. Valid values: 'poles', 'zeros', 'polesZeros.
<i>f_fval</i>	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.
<i>f_rval</i>	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

<i>t</i>	Returns <i>t</i> if it plots a report.
<i>nil</i>	Returns <i>nil</i> otherwise.

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Examples

```
pzPlot(?resultsDir "/usr/simulation/lowpass/spectre/schematic" ?result 'pz)
```

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).

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Plotting and Printing Commands

pzSummary

```
pzSummary( [?resultsDir t_resultsDir] [?result S_resultName]  
          [?print S_toPrint] [?freqfilter f_fval] [?realfilter f_rval] )  
=>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.

Arguments

<i>t_resultsDir</i>	Directory containing the results. If you specify a directory with <i>resultsDir</i> , the <i>pzSummary</i> command plots the results for that directory.
<i>S_resultName</i>	Pointer to results from the analysis for which you want to print the report.
<i>S_toPlot</i>	Use this option to plot only poles, only zeros or both poles and zeros information. Valid values: 'poles', 'zeros', 'polesZeros.
<i>f_fval</i>	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.
<i>f_rval</i>	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

<i>t</i>	Returns <i>t</i> if it prints a report.
<i>nil</i>	Returns <i>nil</i> otherwise.

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Plotting and Printing Commands

Examples

```
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 greater than 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).

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Plotting and Printing Commands

removeLabel

```
removeLabel( l_id )  
=> t/nil
```

Description

Removes the label, or all the labels identified in a list, from the current subwindow.

Arguments

l_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.

Examples

```
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.

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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 `?output` argument, the `report` command opens the file and writes the information to it. If you provide a port (the return value of the `SKILL outfile` command), the `report` 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 `dcOp` and `opBegin`.

Arguments

<i>t_filename</i>	File in which to write the information. The <code>report</code> 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 <code>getSkillPath()</code> at the OCEAN prompt.
<i>p_port</i>	Port (previously opened with <code>outfile</code>) through which to append the information to a file. You are responsible for closing the port. See the <code>outfile</code> command for more information.
<i>t_type</i>	Type of information to print, such as all bjts.
<i>t_name</i>	Name of the node or component.
<i>t_param</i>	Name of the parameter to print. It is also a list.
<i>s_reportStyle</i>	Specifies the format of the output. Valid values: <code>spice</code> and <code>paramValPair</code> Default value: <code>paramValPair</code>

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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

charsPerLine

Number of characters to be printed per line.

Value Returned

`t` Returns `t` if the information is printed.

`nil` Returns `nil` and an error message if the information cannot be printed.

Examples

```
selectResult( dcOp )  
= > t  
report()
```

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 `bjts`.

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```
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.

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xLimit

```
xLimit( l_minMax )  
=> t/nil
```

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.

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yLimit

```
yLimit( l_minMax [?yNumber x_yNumber] [?stripNumber x_stripNumber])  
=> t/nil
```

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

<i>l_minMax</i>	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 <i>nil</i> , the limit is set to <i>auto</i> .
<i>x_yNumber</i>	Specifies the Y axis that will have limited display with the range specified by <i>l_minMax</i> . Valid values: 1 through 4
<i>x_stripNumber</i>	Specifies the strip in which the y display is to be limited as specified by <i>x_yNumber</i> . Valid values: 1 through 20

Value Returned

<i>t</i>	Returns <i>t</i> if the Y axis display limits are set.
<i>nil</i>	Returns <i>nil</i> and an error message if the Y axis display limits cannot be set.

Example

```
yLimit( list( 4.5 7.5 ) ?yNumber 1 )  
=> t
```

Sets Y axis 1 to display from 4.5 to 7.5.

```
yLimit( list( 4.5 7.5 ) ?yNumber 1 ?stripNumber 3)
```

OCEAN Reference

Plotting and Printing Commands

Sets Y axis 1 to display from 4.5 to 7.5 in `stripNumber` 3.

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 Artist environment, click *Results – Direct Plot – S-param*. 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 Artist environment. When all the desired expressions are created in the *Outputs* pane, use the *ADE – Session – Save 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 very similar to plotting RN.

```
; start ocean with SpectreS as the simulator.
simulator( 'spectreS )
; if you wanted to use Spectre as the simulator, then
; simulator( 'spectre )
;specify design and model path
design(  "/usr1/mnt4/myhome/simulation/myckt/spectreS/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
desVar(  "r2" 37)
desVar(  "r1" 150)
;set temperature
temp( 25 )
```

OCEAN Reference

Plotting and Printing Commands

```
;run sp noise analysis with the above desVar list.
run()
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 *SpectreRF User Guide*.

For more information on these functions, click these links: [getData](#), [sp](#), [ocnPrint](#), and [plot](#).

OCEAN Reference
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 `selectResult`. 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
<code>vm</code>	<code>vm(t_net [?resultsDir t_resultsDir][?result s_resultname]) => o_waveform/nil</code>	Aliased to <code>mag(v())</code> . Gets the magnitude of the voltage of a net.
<code>vdb</code>	<code>vdb(t_net [?resultsDir t_resultsDir][?result s_resultname]) => o_waveform/nil</code>	Aliased to <code>db20(v())</code> . Gets the power gain in decibels from net in to net out.
<code>vp</code>	<code>vp(t_net [?resultsDir t_resultsDir][?result s_resultname]) => o_waveform/nil</code>	Aliased to <code>phase(v())</code> . Gets the phase of the voltage of a net.

OCEAN Reference

OCEAN Aliases

List of Aliases, *continued*

vr	<code>vr(t_net [?resultsDir t_resultsDir][?result s_resultName]) => o_waveform/ nil</code>	Aliased to <code>real(v())</code> . Gets the real part of a complex number representing the voltage of a net.
vim	<code>vim(t_net [?resultsDir t_resultsDir][?result s_resultName]) => o_waveform/ nil</code>	Aliased to <code>imag(v())</code> . Gets the imaginary part of a complex number representing the voltage of a net.
im	<code>im(t_component [?resultsDir t_resultsDir][?result s_resultName]) => o_waveform/ nil</code>	Aliased to <code>mag(i())</code> . Gets the magnitude of the AC current through a component.
ip	<code>ip(t_component [?resultsDir t_resultsDir][?result s_resultName]) => o_waveform/ nil</code>	Aliased to <code>phase(i())</code> . Gets the phase of the AC current through a component.
ir	<code>ir(t_component [?resultsDir t_resultsDir][?result s_resultName]) => o_waveform/ nil</code>	Aliased to <code>real(i())</code> . Gets the real part of a complex number representing the AC current through a component.
iim	<code>iim(t_component [?resultsDir t_resultsDir][?result s_resultName]) => o_waveform/ nil</code>	Aliased to <code>imag(i())</code> . Gets the imaginary part of a complex number representing the AC current through a component.

Predefined Functions and Waveform (Calculator) Functions

This chapter contains information about the following functions. Some additional predefined data access commands are described in the Virtuoso® Analog Design Environment SKILL Language Reference manual.

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Predefined Arithmetic Functions

Several functions are predefined in the Virtuoso[®] SKILL language. The full syntax and brief definitions for these functions follows the table.

Predefined Arithmetic Functions

Synopsis	Result
General Functions	
<code>add1(n)</code>	$n + 1$
<code>abs</code>	$ n $
<code>sub1(n)</code>	$n - 1$
<code>exp(n)</code>	e raised to the power n
<code>linRg(<i>n_from</i>, <i>n_to</i>, <i>n_by</i>)</code>	Returns list of numbers in linear range from <i>n_from</i> to <i>n_to</i> in <i>n_by</i> steps
<code>log(n)</code>	Natural logarithm of n
<code>logRg(<i>n_from</i>, <i>n_to</i>, <i>n_by</i>)</code>	Returns list of numbers in log10 range from <i>n_from</i> to <i>n_to</i> in <i>n_by</i> steps
<code>max(<i>n1</i> <i>n2</i> ...)</code>	Maximum of the given arguments
<code>min(<i>n1</i> <i>n2</i> ...)</code>	Minimum of the given arguments
<code>mod(<i>x1</i> <i>x2</i>)</code>	<i>x1</i> modulo <i>x2</i> , that is, the integer remainder of dividing <i>x1</i> by <i>x2</i>
<code>round(n)</code>	Integer whose value is closest to n
<code>sqrt(n)</code>	Square root of n

OCEAN Reference
Predefined Functions and Waveform (Calculator) Functions

Predefined Arithmetic Functions

Synopsis	Result
-----------------	---------------

Trigonometric Functions

<code>sin(<i>n</i>)</code>	sine, argument <i>n</i> is in radians
<code>cos(<i>n</i>)</code>	cosine
<code>tan(<i>n</i>)</code>	tangent
<code>asin(<i>n</i>)</code>	arc sine, result is in radians
<code>acos(<i>n</i>)</code>	arc cosine
<code>atan(<i>n</i>)</code>	arc tangent

Random Number Generator

<code>random(<i>x</i>)</code>	Returns a random integer between 0 and <i>x</i> -1. If <code>random</code> is called with no arguments, it returns an integer that has all of its bits randomly set.
<code>srandom(<i>x</i>)</code>	Sets the initial state of the random number generator to <i>x</i> .

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

abs

```
abs( n_number )  
=> n_result
```

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*.

Example

```
abs( -209.625 )  
=> 209.625  
abs( -23 )  
=> 23
```

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

acos

```
acos( n_number )  
    => f_result
```

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
```


OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

add1

```
add1( n_number )  
    => n_result
```

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

```
add1( 59 )  
=> 60
```

Adds 1 to 59.

OCEAN Reference

Predefined Functions 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*.

Example

```
asin(0.3)  
=> 0.3046927
```

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

atan

```
atan( n_number )  
    => f_result
```

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
```

OCEAN Reference

Predefined Functions 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*.

Examples

```
cos(0.3)  
=> 0.9553365  
cos(3.14/2)  
=> 0.0007963
```

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

exp

```
exp( n_number )  
    => f_result
```

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*.

Examples

```
exp( 1 )  
=> 2.718282  
exp( 3.0 )  
=> 20.08554
```

OCEAN Reference

Predefined Functions 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

<i>n_from</i>	Smaller number in the linear range.
<i>n_to</i>	Larger number in the linear range.
<i>n_by</i>	Increment value when stepping through the range.

Value Returned

<i>l_range</i>	List of numbers in the linear range.
nil	Returned if error.

Example

```
range = linRg(-30 30 5)  
(-30 -25 -20 -15 -10 -5 0 5 10 15 20 25 30)
```

OCEAN Reference

Predefined Functions 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*.

Example

```
log( 3.0 )  
=> 1.098612
```

OCEAN Reference

Predefined Functions 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

<i>n_from</i>	Smaller number in the linear range.
<i>n_to</i>	Larger number in the linear range.
<i>n_by</i>	Increment value when stepping through the range.

Value Returned

<i>l_range</i>	List of numbers in the linear range.
nil	Returned if error.

Example

```
logRg(1 1M 0.5)  
(1.0 100.0 10000.0 1000000.0)
```


OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

max

```
max( n_num1 n_num2 [n_num3 ...] )  
=> n_result
```

Description

Returns the maximum of the values passed in. Requires a minimum of two arguments.

Arguments

<i>n_num1</i>	First value to check.
<i>n_num2</i>	Next value to check.
[<i>n_num3</i> ...]	Additional values to check.

Value Returned

<i>n_result</i>	The maximum of the values passed in.
-----------------	--------------------------------------

Examples

```
max(3 2 1)  
=> 3  
max(-3 -2 -1)  
=> -1
```

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

min

```
min( n_num1 n_num2 [n_num3 ...] )  
    => n_result
```

Description

Returns the minimum of the values passed in. Requires a minimum of two arguments.

Arguments

<i>n_num1</i>	First value to check.
<i>n_num2</i>	Next value to check.
[<i>n_num3</i> ...]	Additional values to check.

Value Returned

<i>n_result</i>	The minimum of the values passed in.
-----------------	--------------------------------------

Examples

```
min(1 2 3)  
=> 1  
min(-1 -2.0 -3)  
=> -3.0
```

OCEAN Reference

Predefined Functions 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.

Example

```
mod(4 3)  
=> 1
```

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

random

```
random( [x_number] )  
=> x_result
```

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.

Example

```
random( 93 )  
=> 26
```

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

round

```
round( n_arg )  
    => x_result
```

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*.

Examples

```
round(1.5)  
=> 2  
round(-1.49)  
=> -1  
round(1.49)  
=> 1
```

OCEAN Reference

Predefined Functions 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*.

Examples

```
sin(3.14/2)  
=> 0.9999997  
sin(3.14159/2)  
=> 1.0
```

Floating-point results from evaluating the same expressions might be machine-dependent.

OCEAN Reference

Predefined Functions 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*.

Examples

```
sqrt( 49 )  
=> 7.0  
sqrt( 43942 )  
=> 209.6235
```

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

srandom

```
srandom( x_number )  
=> t
```

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.

Example

```
srandom( 89 )  
=> t
```


OCEAN Reference

Predefined Functions 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

```
sub1( 59 )  
=> 58
```

Subtracts 1 from 59.

OCEAN Reference

Predefined Functions 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*.

Example

```
tan( 3.0 )  
=> -0.1425465
```

Waveform (Calculator) Functions

The calculator commands are described in this section.

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

average

```
average( o_waveform )  
=> 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 $y=f(x)$, $average(y)=$

$$\frac{\int_{from}^{to} f(x)dx}{to - from}$$

where *from* is the initial value for x and *to* is the final value.

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <i>drwave:XXXXX</i> .)
-------------------	--

Value Returned

<i>n_average</i>	Returns a number representing the average value of the input waveform.
------------------	--

<i>o_waveformAverage</i>	Returns a waveform (or family of waveforms) representing the average value if the input is a family of waveforms.
--------------------------	---

<i>nil</i>	Returns <i>nil</i> and an error message otherwise.
------------	--

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

Example

```
average( v( "/net9" ) )
```

Gets the average voltage (Y-axis value) of `/net9` over the entire time range specified in the simulation analysis.

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

awvPlaceXMarker

```
awvPlaceXMarker(w_windowId n_xLoc [?subwindow x_subwindowId])  
=> t_xLoc/t/nil
```

Description

Places a vertical marker at a specific x-coordinate in the optionally specified subwindow of the specified window.

Arguments

<i>w_windowId</i>	Waveform window ID.
<i>n_xLoc</i>	The x-coordinate at which to place the marker.
<i>x_subwindowId</i>	Waveform subwindow ID.

Value Returned

<i>t_xLoc</i>	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.

Examples

```
awvPlaceXMarker( window 5)  
=> "5"
```

Vertical marker info form is opened when the command is executed.

```
awvPlaceXMarker( window 6 ?subwindow 2)  
=> t
```

Vertical marker info form is not opened.

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

awvPlaceYMarker

```
awvPlaceYMarker(w_windowId n_yLoc [?subwindow x_subwindowId])  
=> t_yLoc/t/nil
```

Description

Places a horizontal marker at a specific y-coordinate in the optionally specified subwindow of the specified window.

Arguments

<i>w_windowId</i>	Waveform window ID.
<i>n_yLoc</i>	The y-coordinate at which to place the marker.
<i>x_subwindowId</i>	Waveform subwindow ID.

Value Returned

<i>t_yLoc</i>	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 <i>nil</i> or an error message.

Examples

```
awvPlaceYMarker( window 5)  
=> "5"
```

Horizontal marker info form is opened when the command is executed.

```
awvPlaceYMarker( window 6 ?subwindow 2)  
=> t
```

Horizontal marker info form is not opened.

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

b1f

```
b1f( o_s11 o_s12 o_s21 o_s22 )  
    => o_waveform/nil
```

Description

Returns the alternative stability factor in terms of the supplied parameters.

Arguments

<i>o_s11</i>	Waveform object representing s11.
<i>o_s12</i>	Waveform object representing s12.
<i>o_s21</i>	Waveform object representing s21.
<i>o_s22</i>	Waveform object representing s22.

Value Returned

<i>o_waveform</i>	Waveform object representing the alternative stability factor.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

Examples

```
s11 = sp(1 1)  
s12 = sp(1 2)  
s21 = sp(2 1)  
s22 = sp(2 2)  
plot(b1f(s11 s12 s21 s22))
```


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Predefined Functions 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

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>n_db</i>	Positive number that defines the bandwidth.
<i>t_type</i>	Type of input filter. Valid values: "low", "high" or "band".

Value Returned

<i>n_value</i>	Returns a number representing the value of the bandwidth if the input argument is a single waveform.
<i>o_waveform</i>	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.

Examples

```
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.

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Predefined Functions 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

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <i>drwave:XXXXX</i> .)
<i>n_from</i>	Starting value for the range on the X axis.
<i>n_to</i>	Ending value for the range on the X axis.

Value Returned

<i>o_waveform</i>	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.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

Examples

```
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 ) )
```

Plots the original waveform.

```
plot( clip( v( "/net9" ) nil 3m ) )
```

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Predefined Functions and Waveform (Calculator) Functions

Plots the portion of a waveform that ranges from 0 to 3 ms.

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Predefined Functions and Waveform (Calculator) Functions

compression

```
compression( o_waveform [ ?x f_x ] [ ?y f_y ] [ ?compression f_compression ]  
            [ ?io s_measure ] )  
            => 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.

Arguments

<i>o_waveform</i>	Waveform object representing output power (in dBm) versus input power (in dBm).
<i>f_x</i>	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 <i>f_y</i> is specified, defaults to the X coordinate of the first point of the <i>o_waveform</i> wave.
<i>f_y</i>	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 <i>f_x</i> is specified, defaults to the Y coordinate of the first point of the <i>o_waveform</i> wave.
<i>f_compression</i>	The delta (in dB) between the power waveform and the ideal gain line that marks the compression point Default value: 1
<i>s_measure</i>	Symbol indicating whether the measurement is to be input referred ('input') or output referred ('output') Default value: 'input'

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Predefined Functions and Waveform (Calculator) Functions

Value Returned

<i>f_compPoint</i>	Depending on the setting of <i>s_measure</i> , returns either input referred or output referred compression point.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

Examples

```
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)))
compression(dBm(harmonic(spectralPower(v("/Pif"
?result "pss_fd")/ 50.0
v("/Pif" ?result "pss_fd")) 2)))
compression(dBm(harmonic(spectralPower(v("/Pif"
?result "pss_fd")/resultParam("rif:r"
?result "pss_td") v("/Pif"
?result "pss_fd")) 2)))
compression(dBm(harmonic(spectralPower(i("/rif/PLUS"
?result "pss_fd") v("/Pif" ?result "pss_fd")) 2))
?x -25 ?compress 0.1 ?measure "Output")
```

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Predefined Functions 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] )
=> 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 `dBm(spectralPower((i or v/r),v))` to calculate a power waveform. The function passes this power curve and the remaining arguments to the `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.

Arguments

<i>o_vport</i>	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.
<i>x_harm</i>	Harmonic index of the voltage wave contained in <i>o_vport</i> . When <i>o_iport</i> is specified, also applies to a current waveform contained in <i>o_iport</i> .
<i>o_iport</i>	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: <code>nil</code>
<i>f_rport</i>	Resistance into the output port. When specified and <i>o_iport</i> is <code>nil</code> , the output power is calculated using voltage and resistance. Default value: <code>50</code>

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Predefined Functions and Waveform (Calculator) Functions

<i>f_epoint</i>	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 <i>o_waveform wave</i>
<i>f_gcomp</i>	The delta (in dB) between the power waveform and the ideal gain line that marks the compression point. Default value: 1
<i>s_measure</i>	Symbol indicating if measurement is to be input referred ('input) or output referred ('output). Default value: 'input
Value Returned	
<i>o_waveform</i>	Returns a waveform when <i>o_waveform1</i> is a family of waveforms.
<i>f_number</i>	Returns a number when <i>o_waveform1</i> is a waveform.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

Examples

Each of the following returns a compression measurement:

```
compressionVRI(v("/Pif" ?result "pss_fd") 2)
compressionVRI(v("/Pif" ?result "pss_fd") 2
  ?rport resultParam("rif:r" ?result "pss_td"))
compressionVRI(v("/Pif" ?result "pss_fd") 2
  ?iport i("/rif/PLUS" ?result "pss_fd") ?epoint -25
  ?gcomp 0.1 ?measure "Output")
```

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Predefined Functions and Waveform (Calculator) Functions

compressionVRICurves

```
compressionVRICurves( o_vport x_harm [?iport o_iport] [?rport f_rport]
    [?epoint f_epoint] [?gcomp f_gcomp] )
    => 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 `compressionVRI` function for making measurements.

Arguments

<i>o_vport</i>	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.
<i>x_harm</i>	Harmonic index of the wave contained in <i>o_vport</i> . When <i>o_iport</i> is specified, also applies to a current waveform contained in <i>o_iport</i> .
<i>o_iport</i>	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: <code>nil</code>
<i>f_rport</i>	Resistance into the output port. When specified and <i>o_iport</i> is <code>nil</code> , the output power is calculated using voltage and

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	resistance. Default value: 50
<i>f_epoint</i>	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 <i>o_waveform</i> wave
<i>f_gcomp</i>	The delta (in dB) between the power waveform and the ideal gain line that marks the compression point. Default value: 1
Value Returned	
<i>o_waveform</i>	Returns a family of waveforms containing the output power and tangent line.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

Examples

Each of following examples returns curves related to a compression measurement:

```
compressionVRICurves(v("/Pif" ?result "pss_fd") 2)
compressionVRICurves(v("/Pif" ?result "pss_fd") 2
  ?rport resultParam("rif:r" ?result "pss_td"))
compressionVRICurves(v("/Pif" ?result "pss_fd") 2
  ?iport i("/rif/PLUS" ?result "pss_fd") ?epoint -25
  ?gcomp 0.1)
```

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Predefined Functions 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

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>n_x</i>	Complex or imaginary number.

Value Returned

<i>o_waveform</i>	Returns the conjugate of a waveform if the input argument is a waveform.
<i>n_y</i>	Returns the result of <i>n_x</i> being mirrored against the real axis (X axis) if the input argument is a number.
<i>nil</i>	Returns <i>nil</i> 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.

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Predefined Functions 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

<i>f_real</i>	The real part of the complex number.
<i>f_imaginary</i>	The imaginary part of the complex number.

Value Returned

<i>o_complex</i>	Returns the complex number.
------------------	-----------------------------

Example

```
complex( 1.0 2.0 )  
=> complex( 1, 2 )
```

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Predefined Functions and Waveform (Calculator) Functions

complexp

```
complexp( g_value )  
=> t/nil
```

Description

Checks if an object is a complex number. The suffix *p* is added to the name of a function to indicate that it is a predicate function.

Arguements

g_value A skill object.

Values Returned

t Returns *t* when *g_value* is a complex number.

nil Returns *nil* if there is an error.

Example

```
complexp( (complex 0 1) )  
=> t  
complexp( 1.0 )  
=> nil
```

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Predefined Functions 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

$$\int_{\text{from}}^{\text{to}} f1(s)f2(t-s)ds$$

$f1$ and $f2$ 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.

Arguments

<i>o_waveform1</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>o_waveform2</i>	Additional waveform object.
<i>n_from</i>	Starting point (X-axis value) of the integration range.
<i>n_to</i>	Ending point (X-axis value) of the integration range.
<i>t_type</i>	Type of interpolation. Valid values: "linear" or "log".
<i>n_by</i>	Increment.

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Predefined Functions and Waveform (Calculator) Functions

Value Returned

<i>o_waveform</i>	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.
<i>n_number</i>	Returns a value representing the convolution if both of the input arguments are numbers.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

Example

```
sinWave = expr( n sin( n ) linRg( 0 20 0.01 ) )
triWave = artListToWaveform( '( ( -4, 0 ) ( -3, 1 ) ( -2, 0 ) ( -1, -1 ) ( 0, 0 )
( 1, 1 ) ( 2, 0 ) ( 3, -1 ) ( 4, 0 ) )' )
plot( convolve( sinWave triWave 0 10 "linear" 1 ) )
```

Gets the waveform from the convolution of the sine waveform and triangle waveform within the range of 0 to 10.

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Predefined Functions 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 b_closeCont]
             [?modifier s_modifier] )
=> 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 `spectralPower` function. The reference reflection coefficients are calculated using voltage, current, and a reference resistance.

Arguments

<i>o_iwave</i>	Current used to calculate power, expected to be a two-dimensional family of harmonic waveforms.
<i>o_vwave</i>	Voltage used to calculate power, expected to be a two-dimensional family of harmonic waveforms.
<i>x_harm</i>	Harmonic index of the waves contained in <i>o_iwave</i> and <i>o_vwave</i> .
<i>o_iwaveLoad</i>	Current used to calculate reflection coefficient, expected to be a two-dimensional family of harmonic waveforms. Default value: <i>o_iwave</i>
<i>o_vwaveLoad</i>	Voltage used to calculate reflection coefficient, expected to be a two-dimensional family of harmonic waveforms. Default value: <i>o_vwave</i>
<i>f_maxPower</i>	Maximum power magnitude value for contours. Default value: automatic
<i>f_minPower</i>	Minimum power magnitude value for contours. Default value: automatic

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Predefined Functions and Waveform (Calculator) Functions

<i>x_numCont</i>	Total number of contours returned. Default value: 8
<i>f_refImp</i>	Reference resistance used to calculate reflection coefficients. Default value: 50
<i>b_closeCont</i>	Boolean indicating when to close the contours. When <i>nil</i> , largest segment of each contour is left open. Default value: <i>nil</i>
<i>s_modifier</i>	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: 'mag
Value Returned	
<i>o_waveform</i>	Returns a family of waveforms (contours) for Z-Smith plotting.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

Examples

The following example plots constant output power contours according to output:

```
cPwrContour(i("/I8/out" ?result "pss_fd") v("/net28"
?result "pss_fd")1)
```

The following example plots constant output power contours according to output reflection coefficients:

```
cPwrContour(i("/I8/out" ?result "pss_fd") v("/net28"
?result "pss_fd") 1 ?maxPower 0.002 ?minPower 0.001 ?numCont 9)
```

The following example plots constant input power contours according to output reflection coefficients:

```
cPwrContour(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 ?modifier "mag")
```


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Predefined Functions 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 b_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.

Arguments

<i>o_iwave</i>	Current used to calculate reflection coefficient magnitude, expected to be a two-dimensional family of spectrum waveforms.
<i>o_vwave</i>	Voltage used to calculate reflection coefficient magnitude, expected to be a two-dimensional family of spectrum waveforms.
<i>x_harm</i>	Harmonic index of the waves contained in <i>o_iwave</i> and <i>o_vwave</i> .
<i>o_iwaveLoad</i>	Current used to calculate reference reflection coefficient, expected to be a two-dimensional family of harmonic waveforms. Default value: <i>o_iwave</i>
<i>o_vwaveLoad</i>	Voltage used to calculate reference reflection coefficient, expected to be a two-dimensional family of spectrum waveforms. Default value: <i>o_vwave</i>
<i>f_maxRefl</i>	Maximum reflection coefficient magnitude value for contours. Default value: automatic
<i>f_minRefl</i>	Minimum reflection coefficient magnitude value for contours. Default value: automatic

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Predefined Functions and Waveform (Calculator) Functions

<i>x_numCont</i>	Total number of contours returned. Default value: 8
<i>f_refImp</i>	Reference resistance used to calculate reflection coefficients. Default value: 50
<i>b_closeCont</i>	Boolean indicating when to close the contours. When <i>nil</i> , the largest segment of each contour is left open. Default value: <i>nil</i>
Value Returned	
<i>o_waveform</i>	Returns a family of waveforms (contours) for Z-Smith plotting.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

Examples

The following example plots constant output reflection coefficient contours according to output reflection coefficients:

```
cReflContour(i("/I8/out" ?result "pss_fd") v("/net28"
?result "pss_fd") 1)
```

The following example plots constant output reflection coefficient contours according to output reflection coefficients:

```
cReflContour(i("/I8/out" ?result "pss_fd") v("/net28"
?result "pss_fd") 1 ?maxRefl 0.7 ?minRefl 0.1 ?numCont 7)
```

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)
```

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Predefined Functions and Waveform (Calculator) Functions

CROSS

```
cross( o_waveform n_crossVal x_n s_crossType )  
      => 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.

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>n_crossVal</i>	Y-axis value at which the corresponding values of X are calculated.
<i>x_n</i>	Number that specifies which X value to return. If <i>x_n</i> equals 1, the first X value with a crossing is returned. If <i>x_n</i> equals 2, the second X value with a crossing is returned, and so on. If you specify a negative integer for <i>x_n</i> , the X values with crossings are counted from right to left (from maximum to minimum).
<i>s_crossType</i>	Type of the crossing. Valid values: 'rising', 'falling', 'either.'

Value Returned

<i>o_waveform</i>	Returns a waveform if the input argument is a family of waveforms.
<i>g_value</i>	Returns the X-axis value of the crossing point if the input argument is a single waveform.
<i>nil</i>	Returns <code>nil</code> and an error message otherwise.

Examples

```
cross( v( "/net9" ) 2.5 2 'rising )
```

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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.

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Predefined Functions and Waveform (Calculator) Functions

db10

```
db10( {o_waveform | n_number} )  
=> o_waveform/n_number/nil
```

Description

Returns 10 times the \log_{10} of the specified waveform object or number. This function can also be written as dB10.

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>n_number</i>	Number.

Value Returned

<i>o_waveform</i>	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.
<i>n_number</i>	Returns a number if the input argument is a number.
<i>nil</i>	Returns <code>nil</code> and an error message otherwise.

Examples

```
db10( ymax( v( "/net9" ) ) )
```

Returns a waveform representing $\log_{10}(\text{ymax}(v("/net9"))) multiplied by 10.$

```
db10( 1000 )  
=> 30.0
```

Gets the value $\log_{10}(1000)$ multiplied by 10, or 30.

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Predefined Functions 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

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>n_number</i>	Number.

Value Returned

<i>o_waveform</i>	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.
<i>n_number</i>	Returns a number if the input argument is a number.
<i>nil</i>	Returns <code>nil</code> and an error message otherwise.

Examples

```
db20( ymax( v( "/net9" ) ) )
```

Returns a waveform representing 20 times $\log_{10}(\text{ymax}(v("/net9")))$.

```
db20( 1000 )  
=> 60.0
```

Returns the value of 20 times $\log_{10}(1000)$, or 60.

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Predefined Functions and Waveform (Calculator) Functions

dbm

```
dbm( {o_waveform | n_number} )  
=> o_waveform/n_number/nil
```

Description

Returns 10 times the log₁₀ of the specified waveform object plus 30. This function can also be written as dBm.

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <i>drwave:XXXXX</i> .)
<i>n_number</i>	Number.

Value Returned

<i>o_waveform</i>	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.
<i>n_number</i>	Returns a number if the input argument is a number.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

Example

```
dbm( ymax( v( "/net9" ) ) ) )
```

Returns a waveform representing 10 times log₁₀(ymax(v("/net9"))) plus 30.

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Predefined Functions and Waveform (Calculator) Functions

delay

```
delay( ?wf1 o_wf1 ?value1 n_value1 ?edge1 s_edge1 ?nth1 x_nth1 ?td1 n_td1
      ?wf2 o_wf2 ?value2 n_value2 ?edge2 s_edge2 ?nth2 x_nth2 {[?td2 n_td2] |
      [?td2r0 n_td2r0]} ?stop n_stop )
=> 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.

Arguments

<i>o_wf1</i>	First waveform object.
<i>n_value1</i>	Value at which the crossing is significant for the first waveform object.
<i>s_edge1</i>	Type of the edge that must cross <i>n_value1</i> . Valid values: 'rising', 'falling', 'either'
<i>x_nth1</i>	Number that specifies which crossing is to be the trigger event. For example, if <i>x_nth1</i> is 2, the trigger event is the second edge of the first waveform with the specified type that crosses <i>n_value1</i> .
<i>n_td1</i>	Time at which to start the delay measurement. The simulator begins looking for the trigger event, as defined by <i>o_wf1</i> , <i>n_value1</i> , <i>s_edge1</i> , and <i>x_nth1</i> , only after the <i>n_td1</i> time is reached.
<i>o_wf2</i>	Second waveform object.
<i>n_value2</i>	Value at which the crossing is significant for the second waveform.
<i>s_edge2</i>	Type of the edge for the second waveform. Valid values: 'rising', 'falling', 'either'
<i>x_nth2</i>	Number that specifies which crossing is to be the target event. For example, if <i>x_nth2</i> is 2, the target event is the second edge

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Predefined Functions and Waveform (Calculator) Functions

of the second waveform with the specified type that crosses *n_value2*.

n_td2

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_wf2*, *n_value2*, *t_edge2*, and *x_nth2*, only after the *n_td2* time is reached.

If you specify neither *n_td2* nor *n_td2r0*, the simulator begins looking for the target event at $t = 0$.

n_td2r0

Time to start observing the target event, relative to $t = 0$. Only applicable if both *o_wf1* and *o_wf2* 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_wf2*, *n_value2*, *t_edge2*, and *x_nth2*, only after the *n_tdr0* time is reached.

If 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.

n_stop

Time to stop observing the target event.

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.

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Predefined Functions and Waveform (Calculator) Functions

Examples

```
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 wf2 ?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.

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Predefined Functions 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: After the second derivative, the results become inaccurate because the derivative is obtained numerically.

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
-------------------	--

Value Returned

<i>o_waveform</i>	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.
-------------------	---

<i>nil</i>	Returns <code>nil</code> and an error message otherwise.
------------	--

Example

```
plot( deriv( v( "/net8" ) ) )
```

Plots the waveform representing the derivative of the voltage of `"/net8"`.

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Predefined Functions and Waveform (Calculator) Functions

dft

```
dft( o_waveform n_from n_to x_num [t_windowName [n_param1]] )  
=> o_waveform/nil
```

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 $w(to)$.

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>n_from</i>	Starting value for the <code>dft</code> computation.
<i>n_to</i>	Ending value for the <code>dft</code> computation.
<i>x_num</i>	Number of timepoints. If <i>x_num</i> is not a power of 2, it is forced to be the next higher power of 2.
<i>t_windowName</i>	Variable representing different methods for taking a <code>dft</code> computation. Valid values: <code>Rectangular</code> , <code>ExtCosBell</code> , <code>HalfCycleSine</code> , <code>Hanning</code> or <code>Cosine2</code> , <code>Triangle</code> or <code>Triangular</code> , <code>Half3CycleSine</code> or <code>HalfCycleSine3</code> , <code>Hamming</code> , <code>Cosine4</code> , <code>Parzen</code> , <code>Half6CycleSine</code> or <code>HalfCycleSine6</code> , <code>Blackman</code> , or <code>Kaiser</code> .

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Predefined Functions and Waveform (Calculator) Functions

For more information about *windowName*, see the information about Discrete Fourier Transform (dft) in the *Virtuoso® Analog Design Environment User Guide*.

n_param1

Smoothing parameter.

Applies only if the *t_windowName* argument is set to `Kaiser`.

Value Returned

o_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

```
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.

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Predefined Functions 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.

Arguments

<i>o_waveform1</i>	Time domain waveform object with units of volts or amps.
<i>o_waveform2</i>	Time domain waveform object with units of volts or amps.
<i>f_timeStart</i>	Start time for the spectral analysis interval. Use this parameter and <i>f_timeEnd</i> to exclude part of the interval. For example, you might set these values to discard initial transient data.
<i>f_timeEnd</i>	End time for the spectral analysis interval.
<i>x_num</i>	The number of time domain points to use. The maximum frequency in the Fourier analysis is directly proportionate to <i>x_num</i> and inversely proportional to the difference between <i>f_timeStart</i> and <i>f_timeEnd</i> .
<i>t_windowName</i>	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, or Kaiser. Default value: Hanning.
<i>x_smooth</i>	The Kaiser window smoothing parameter. If there are no requests, there is no smoothing. Valid values: $0 \leq x_smooth \leq 15$ Default value: 1
<i>f_cohGain</i>	A scaling parameter. A non-zero value scales the power spectral density by $1/(f_cohGain)$. Valid values: $0 \leq f_cohGain \leq 1$. You can use 1 if you do

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Predefined Functions and Waveform (Calculator) Functions

not want the scaling parameter to be used.
Default value: 1

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.

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")
```

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Predefined Functions and Waveform (Calculator) Functions

eyeDiagram

```
eyeDiagram ( o_waveform n_start n_stop n_period )  
=> o_waveform/nil
```

Description

This function gives an eye-diagram plot of the input waveform signal. It returns the waveform object of the eye-diagram plot.

Arguments

<i>n_start</i>	x-axis start value from where the eye-diagram plot is to commence
<i>n_stop</i>	x-axis stop value where the eye-diagram plot will terminate
<i>n_period</i>	period value

Value Returned

<i>o_waveform</i>	Returns a waveform object representing the eye-diagram plot of the input waveform
<i>nil</i>	Returns <i>nil</i> and an error message otherwise

Example

```
eyeDiagram( v("/out" ) 0n 500n 12.5n)
```


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Predefined Functions and Waveform (Calculator) Functions

flip

```
flip( o_waveform )  
    => o_waveform/nil
```

Description

Returns a waveform with the X vector values negated.

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX.</code>)
-------------------	--

Value Returned

<i>o_waveform</i>	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.
-------------------	--

<i>nil</i>	Returns <i>nil</i> and an error message otherwise.
------------	--

Example

```
plot( flip( v("/net4") ) ) )
```

Plots the waveform for the voltage of `"/net4"` with the X vector values negated.

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Predefined Functions and Waveform (Calculator) Functions

fourEval

```
fourEval( o_waveform n_from n_to n_by )  
=> 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 `dft` command. The `fourEval` function transforms the expression from the frequency domain to the time domain.

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>n_from</i>	Starting point on the X axis at which to start the evaluation.
<i>n_to</i>	Increment.
<i>n_by</i>	Ending point on the X axis.

Value Returned

<i>o_waveform</i>	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.
<i>nil</i>	Returns <code>nil</code> and an error message otherwise.

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.

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Predefined Functions 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

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
-------------------	--

Value Returned

<i>o_waveform</i>	Returns a waveform representing the frequency of a family of waveforms if the input argument is a family of waveforms.
<i>n_value</i>	Returns a number representing the frequency of the specified waveform.
<i>nil</i>	Returns <code>nil</code> and an error message otherwise.

Example

```
frequency( v( "/net12" ) )
```

Returns the frequency of `"/net12"`.

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Predefined Functions 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

<i>o_s11</i>	Waveform object representing s11.
<i>o_s12</i>	Waveform object representing s12.
<i>o_s21</i>	Waveform object representing s21.
<i>o_s22</i>	Waveform object representing s22.
<i>n_gs</i>	Source reflection coefficient. Default value: 0

Value Returned

<i>o_waveform</i>	Waveform object representing the available gain.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

Examples

```
s11 = sp(1 1)  
s12 = sp(1 2)  
s21 = sp(2 1)  
s22 = sp(2 2)  
plot(ga(s11 s12 s21 s22))
```

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Predefined Functions 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.

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_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, `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, an available gain circle is calculated at each one of the 10 frequencies.

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Predefined Functions and Waveform (Calculator) Functions

Value Returned

o_waveform Waveform object representing the available gain circles.

nil Returns *nil* and an error message otherwise.

Examples

```
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))
```

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Predefined Functions 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.

.

$$\text{gainBwProd}(\text{gain}) = A_o * f_2$$

The gain-bandwidth product is calculated as the product of the DC gain A_o and the critical frequency f_2 . The critical frequency f_2 is the smallest frequency for which the gain equals $1/\sqrt{2}$ times the DC gain A_o .

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
-------------------	--

Value Returned

<i>o_waveform</i>	Returns a waveform representing the gain-bandwidth product for a family of waveforms if the input argument is a family of waveforms.
-------------------	--

<i>n_value</i>	Returns a value for the gain-bandwidth product for the specified waveform.
----------------	--

<i>nil</i>	Returns <code>nil</code> and an error message otherwise.
------------	--

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Predefined Functions and Waveform (Calculator) Functions

Example

```
gainBwProd( v( "/OUT" ) )
```

Returns the gain-bandwidth product for the waveform representing the voltage of the "/OUT" net.

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Predefined Functions and Waveform (Calculator) Functions

gainMargin

```
gainMargin( o_waveform [b_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 b_stable is set to nil. If b_stable value is set to t, then a stable design will have a positive value.

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>b_stable</i>	Boolean optional value that takes the value <code>nil</code> by default.

Value Returned

<i>o_waveform</i>	Returns a waveform representing the gain margin for a family of waveforms if the input argument is a family of waveforms.
<i>n_value</i>	Returns the value for the gain margin of the specified waveform.
<code>nil</code>	Returns <code>nil</code> and an error message otherwise.

Example

```
gainMargin( v( "/OUT" ) ) = -9.234  
gainMargin( v( "/OUT" ) nil ) = -9.234  
gainMargin( v( "/OUT" ) t ) = 9.234
```

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Predefined Functions 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

<i>o_s11</i>	Waveform object representing s11.
<i>o_s12</i>	Waveform object representing s12.
<i>o_s21</i>	Waveform object representing s21.
<i>o_s22</i>	Waveform object representing s22.

Value Returned

<i>o_waveform</i>	Load reflection coefficient.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

Examples

```
s11 = sp(1 1)  
s12 = sp(1 2)  
s21 = sp(2 1)  
s22 = sp(2 2)  
plot(gmax(s11 s12 s21 s22))
```

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Predefined Functions 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:

```
yOpt = o_Gopt + (complex 0 1) * o_Bopt  
return ( 1 / f_zref(1) - yOpt ) / ( 1 / f_zref(1) + yOpt )
```

Arguments

<i>o_Gopt</i>	Waveform object representing the optimum source conductance.
<i>o_Bopt</i>	Waveform object representing the optimum source susceptance.
<i>f_zref</i>	Reference impedance.

Value Returned

<i>o_gminWave</i>	Waveform object representing the optimum noise reflection coefficient.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

Examples

```
Gopt = getData("Gopt")  
Bopt = getData("Bopt")  
Zref = zref(1 ?result "sp")  
plot(gmin(Gopt Bopt Zref))
```

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Predefined Functions 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

<i>o_s11</i>	Waveform object representing s11.
<i>o_s12</i>	Waveform object representing s12.
<i>o_s21</i>	Waveform object representing s21.
<i>o_s22</i>	Waveform object representing s22.

Value Returned

<i>o_waveform</i>	Waveform object representing the maximum stable power gain.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

Example

```
s11 = sp(1 1)  
s12 = sp(1 2)  
s21 = sp(2 1)  
s22 = sp(2 2)  
plot(gmsg(s11 s12 s21 s22))
```

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Predefined Functions and Waveform (Calculator) Functions

gmux

```
gmux( o_s11 o_s12 o_s21 o_s22 )  
      => o_waveform/nil
```

Description

Returns the maximum unilateral power gain in terms of the supplied parameters.

Arguments

<i>o_s11</i>	Waveform object representing s11.
<i>o_s12</i>	Waveform object representing s12.
<i>o_s21</i>	Waveform object representing s21.
<i>o_s22</i>	Waveform object representing s22.

Value Returned

<i>o_waveform</i>	Waveform object representing the maximum unilateral power gain.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

Examples

```
s11 = sp(1 1)  
s12 = sp(1 2)  
s21 = sp(2 1)  
s22 = sp(2 2)  
plot(gmux(s11 s12 s21 s22))
```

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Predefined Functions 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

<i>o_s11</i>	Waveform object representing s11.
<i>o_s12</i>	Waveform object representing s12.
<i>o_s21</i>	Waveform object representing s21.
<i>o_s22</i>	Waveform object representing s22
<i>n_gl</i>	Load reflection coefficient. Default value: 0

Value Returned

<i>o_waveform</i>	Waveform object representing the power gain.
<i>nil</i>	Returns <i>nil</i> 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>))`

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Predefined Functions 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.

Arguments

<i>o_s11</i>	Waveform object representing s11.
<i>o_s12</i>	Waveform object representing s12.
<i>o_s21</i>	Waveform object representing s21.
<i>o_s22</i>	Waveform object representing s22.
<i>g_level</i>	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 <code>linRg</code> function can be called to generate a linear range. For example, <code>linRg(-30 30 5)</code> is the same as <code>list(-30 -25 -20 -15 -10 -5 0 5 10 15 20 25 30)</code> and the <i>g_level</i> argument can be specified as either. In that case, a power gain circle is calculated at each one of the 13 levels.
<i>g_frequency</i>	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, <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, a power gain circle is calculated at each one of the 10 frequencies.

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Predefined Functions and Waveform (Calculator) Functions

Value Returned

<i>o_waveform</i>	Waveform object representing the power gain circles.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

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:

$$\text{Group Delay} = \frac{d\phi}{d\omega} = \frac{d}{df} \left[\frac{\text{phase}(/netX)}{360} \right]$$

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: *drwave: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*.

OCEAN Reference

Predefined Functions 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

<i>o_s11</i>	Waveform object representing s11.
<i>o_s12</i>	Waveform object representing s12.
<i>o_s21</i>	Waveform object representing s21.
<i>o_s22</i>	Waveform object representing s22.
<i>n_gs</i>	Source reflection coefficient. Default value: 0
<i>n_gl</i>	Input reflection coefficient. Default value: 0

Value Returned

<i>o_waveform</i>	Waveform object representing the transducer gain.
<i>nil</i>	Returns <i>nil</i> and displays a message if there is an error.

Examples

```
s11 = sp(1 1)  
s12 = sp(1 2)  
s21 = sp(2 1)  
s22 = sp(2 2)  
plot(gt(s11 s12 s21 s22))
```

Note: *gl* is an imaginary number which should be input in the following format:
`gt(s11 s12 s21 s22 ?gl complex(<realPart> <imagPart>))`

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Predefined Functions 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

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>h_index</i>	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 <code>funds</code> analysis parameter in the netlist. If more than one <i>h_index</i> is desired at one time, a list can be specified.

Value Returned

<i>o_waveform</i>	Returns a waveform (when a single <i>h_index</i> is specified) or family of waveforms (when more than one <i>h_index</i> is specified) if the input argument is a family of waveforms.
<i>g_value</i>	Returns the harmonic value if the input is a single waveform with the X values being harmonics
<i>nil</i>	Returns <code>nil</code> and displays a message if there is an error.

Examples

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.

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Predefined Functions 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:

```
harmonic(v("/net54" ?result "pac-pac") list(1 5))
harmonic(v("/net51" ?result "sweeppss_pss_fd-sweep") '(1 8))
harmonic(v("/Pif" ?result "sweeppss_pac-sweep") list(-8 0))
harmonic(v("/net36" ?result "sweeppdisto_pdisto_fi-sweep") '((1 -1) (2 -2) (-1 2)))
```

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.

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Predefined Functions 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

<i>t_resultsDir</i>	Directory containing the PSF files (results). If you supply this argument, you must also supply the <i>resultName</i> argument.
<i>S_resultName</i>	Results from an analysis.

Value Returned

<i>n_list</i>	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 <i>funds</i> 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.
<i>nil</i>	Returns <i>nil</i> if no harmonics are found in the data.

Examples

For each of the following commands:

```
harmonicFreqList( ?result "pss-fd.pss" )  
harmonicFreqList( ?result "pac-pac" )  
harmonicFreqList( ?result "sweppss_pss_fd-sweep" )  
harmonicFreqList( ?result "sweppss_pac-sweep" )
```

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Predefined Functions and Waveform (Calculator) Functions

Each result is a list of integers.

For each of the following commands:

```
harmonicFreqList( ?result "pdisto-fi.pdisto" )  
harmonicFreqList( ?result "sweep_pdisto_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.

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Predefined Functions 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

<i>t_resultsDir</i>	Directory containing the PSF files (results). If you supply this argument, you must also supply the <i>resultName</i> argument.
<i>S_resultName</i>	Results from an analysis.

Value Returned

<i>n_list</i>	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.

Examples

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" )
```

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Predefined Functions 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.

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Predefined Functions 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: `drwave: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`.

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Predefined Functions 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

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>n_input</i>	Complex number.

Value Returned

<i>o_waveformImag</i>	Returns a waveform when the input argument is a waveform.
<i>n_numberImag</i>	Returns a number when the input argument is a number.
<code>nil</code>	Returns <code>nil</code> and an error message otherwise.

Examples

```
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 `x` representing a complex number, and returns the real portion of that complex number.

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Predefined Functions and Waveform (Calculator) Functions

integ

```
integ(o_waveform,[n_initial_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

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>initial_limit_n</i>	Initial limit for definite integration.
<i>final_limit_n</i>	Final limit for definite integration.

Value Returned

<i>o_waveform</i>	Returns a waveform representing the definite integral for a family of waveforms if the input argument is a family of waveforms.
<i>n_value</i>	Returns a numerical value representing the definite integral of the input waveform if the input argument is a single waveform.
<i>nil</i>	Returns <code>nil</code> and an error message otherwise.

Example

```
integ( v( "/out" ) )
```

Returns the definite integral of the waveform representing the voltage of `"/out"` over its entire range.

```
integ( VT( "/out" ),12.5n,18n)
```

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Predefined Functions and Waveform (Calculator) Functions

Returns the definite integral of the waveform representing the voltage of `"/out"` within a specified range.

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Predefined Functions and Waveform (Calculator) Functions

ipn

```
ipn( o_spurious o_reference [ f_ordspur f_ordref f_epspur f_epref b_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 `ipn` 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 `ipn` function calculates the intersection of these two lines and returns the value of either the X coordinate (input referred) or Y coordinate.

Arguments

<i>o_spurious</i>	Waveform or number representing the spurious output power (in dBm).
<i>o_reference</i>	Waveform or number representing the reference output power (in dBm).
<i>f_ordspur</i>	Order or slope of the spurious constant-slope power line. Default value: 3
<i>f_ordref</i>	Order or slope of the reference constant-slope power line. Default value: 1
<i>f_epspur</i>	Value (in dBm) used to indicate the point where the spurious constant-slope power line begins. If <i>b_pswEEP</i> is <code>t</code> , this value is the input power value of the point on the <i>o_spurious</i> waveform, otherwise this value is paired with the <i>o_spurious</i> value to define the point. This point should be in the linear region of operation. (If <i>b_pswEEP</i> is <code>t</code> , <i>f_spspur</i> defaults to the X coordinate of the first point of the <i>o_spurious</i> wave; if <i>s_measure</i> is <code>'input</code> , a number must be specified.)

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Predefined Functions and Waveform (Calculator) Functions

<i>f_epref</i>	Value (in dBm) used to indicate the point where the reference constant-slope power line begins. If <i>b_pswEEP</i> is <i>t</i> , this value is the input power value of the point on the <i>o_reference</i> waveform, otherwise this value is paired with the <i>o_reference</i> value to define the point. This point should be in the linear region of operation. (If <i>b_pswEEP</i> is <i>t</i> , <i>f_epref</i> defaults to the X coordinate of the first point of the <i>o_reference</i> wave; if <i>s_measure</i> is 'input', a number must be specified.)
<i>b_pswEEP</i>	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: <i>t</i>
<i>s_measure</i>	Name indicating if measurement is to be input referred ('input') or output referred ('output'). Default value: 'input'

Value Returned

<i>o_waveform</i>	Depending on setting of <i>b_pswEEP</i> and the dimension of the input waveforms, returns either a waveform or a family of waveforms.
<i>f_number</i>	If <i>o_spurious</i> and <i>o_reference</i> are numbers or they are waveforms when <i>b_pswEEP</i> is <i>t</i> , returns a number.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

Examples

```
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))
```

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Predefined Functions and Waveform (Calculator) Functions

```
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))
dbm(harmonic(spectralPower(v("/Pif" ?result "pss_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)))
```

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Predefined Functions 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 b_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 `dBm(spectralPower((i or v/r),v))` to 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 `ipn` 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 `ipn` function calculates the intersection of these two lines and returns the value of either the X coordinate (input referred) or the Y coordinate.

Arguments

<i>o_vport</i>	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).
<i>x_harmspur</i>	Harmonic number of the spurious voltage contained in <i>o_vport</i> . When <i>o_iport</i> is specified, also applies to a current waveform contained in <i>o_iport</i> .
<i>x_harmref</i>	Harmonic index of the reference voltage contained in <i>o_vport</i> . When <i>o_iport</i> is specified, also applies to a current waveform contained in <i>o_iport</i> .
<i>o_iport</i>	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.
<i>f_rport</i>	Resistance into the output port. When specified and <i>o_iport</i> is <code>nil</code> , the output power is calculated using voltage and

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Predefined Functions and Waveform (Calculator) Functions

	resistance. Default value: 50
<i>f_ordspur</i>	Order or slope of the spurious constant-slope power line. Default value: 3
<i>f_epoint</i>	Value (in dBm) used to indicate the point where the spurious constant-slope power line begins. If <i>b_pswEEP</i> is <i>t</i> , this value is the input power value of the point on the <i>o_spurious</i> waveform, otherwise this value is paired with the <i>o_spurious</i> value to define the point. This point should be in the linear region of operation. Default value: If <i>b_pswEEP</i> is <i>t</i> , the lowest input power value; if <i>s_measure</i> is 'input, a number must be specified.
<i>b_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: <i>t</i>
<i>f_epref</i>	Value (in dBm) used to indicate the point where the reference constant-slope power line begins. If <i>b_pswEEP</i> is <i>t</i> , this value is the input power value of the point on the <i>o_reference</i> waveform, otherwise this value is paired with the <i>o_reference</i> value to define the point. This point should be in the linear region of operation. Default value: If <i>f_epoint</i> is not <i>nil</i> , <i>f_epoint</i> ; else if <i>b_pswEEP</i> is <i>t</i> , the X coordinate of the first point of the <i>o_reference</i> wave; else if <i>s_measure</i> is 'input, a number must be specified.
<i>f_ordref</i>	Order or slope of the reference constant-slope power line. Default value: 1
<i>s_measure</i>	Symbol indicating if measurement is to be input referred ('input) or output referred ('output). Default value: 'input

Value Returned

<i>o_waveform</i>	Depending on the setting of <i>b_pswEEP</i> and the dimension of input waveform(s), the <i>ipnVRI</i> function returns either a waveform or a family of waveforms.
-------------------	--

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Predefined Functions and Waveform (Calculator) Functions

<i>f_number</i>	Depending on the setting of <i>b_pswEEP</i> and the dimension of input waveform(s), the <i>ipnVRI</i> function returns a number.
<i>nil</i>	Returns <i>nil</i> 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
  ?rport resultParam("rif:r" ?result "pss_td"))
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"))
```

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Predefined Functions and Waveform (Calculator) Functions

ipnVRICurves

```
ipnVRICurves( o_vport x_harmspurs x_harmref [?iport o_iport] [?rport f_rport]
             [?ordspur f_ordspur] [?epoint f_epoint] [?psweep b_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 `ipnVRICurves` 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.

Arguments

<code>o_vport</code>	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).
<code>x_harmspurs</code>	Harmonic index 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> .
<code>x_harmref</code>	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> .
<code>o_iport</code>	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.

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<i>f_rport</i>	Resistance into the output port. When specified and <i>o_iport</i> is <i>nil</i> , the output power is calculated using voltage and resistance. Default value: 50
<i>f_ordspur</i>	Order or slope of the spurious constant-slope power line. Default value: 3
<i>f_epoint</i>	Value (in dBm) used to indicate the point where the spurious constant-slope power line begins. If <i>b_pswEEP</i> is <i>t</i> , this value is the input power value of the point on the <i>o_spurious</i> waveform, otherwise this value is paired with the <i>o_spurious</i> value to define the point. This point should be in the linear region of operation. Default value: If <i>b_pswEEP</i> is <i>t</i> , the X coordinate of the first point of the <i>o_spurious</i> wave; otherwise a number must be specified.
<i>b_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: <i>t</i>
<i>f_epref</i>	Value (in dBm) used to indicate the point where the reference constant-slope power line begins. If <i>b_pswEEP</i> is <i>t</i> , this value is the input power value of the point on the <i>o_reference</i> waveform, otherwise this value is paired with the <i>o_reference</i> value to define the point. This point should be in the linear region of operation. Default value: If <i>f_epoint</i> is not <i>nil</i> , <i>f_epoint</i> ; else if <i>b_pswEEP</i> is <i>t</i> , the X coordinate of the first point of the <i>o_reference</i> wave; else a number must be specified.
<i>f_ordref</i>	Order or slope of the reference constant-slope power line. Default value: 1

Value Returned

<i>o_waveform</i>	A family of waveforms that contains the spurious and reference tangent lines, and when <i>b_pswEEP</i> is <i>t</i> , contains the spurious and reference waveforms.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

Examples

Each of following examples returns curves related to an ip3 measurement:

```
ipnVRICurves(v("/Pif" ?result "pss_fd") 9 8)
ipnVRICurves(v("/Pif" ?result "pss_fd") 9 8
  ?rport resultParam("rif:r" ?result "pss_td"))
ipnVRICurves(v("/Pif" ?result "pss_fd") 9 8
  ?iport i("/rif/PLUS" ?result "pss_fd") ?epoint -25)
ipnVRICurves(v("/Pif" ?result "pac") -21 -25
  ?rport resultParam("rif:r" ?result "pss_td"))
```

OCEAN Reference

Predefined Functions 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

<i>o_s11</i>	Waveform object representing s11.
<i>o_s12</i>	Waveform object representing s12.
<i>o_s21</i>	Waveform object representing s21.
<i>o_s22</i>	Waveform object representing s22.

Value Returned

<i>o_waveform</i>	Waveform object representing the stability factor.
<i>nil</i>	Returns <i>nil</i> if there is an error.

Examples

```
s11 = sp(1 1)  
s12 = sp(1 2)  
s21 = sp(2 1)  
s22 = sp(2 2)  
plot(kf(s11 s12 s21 s22))
```

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

In

```
ln( {o_waveform | n_number} )  
=> o_waveform/f_number/nil
```

Description

Gets the base-e (natural) logarithm of a waveform or number.

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>n_number</i>	Number.

Value Returned

<i>o_waveform</i>	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
<i>f_number</i>	Returns a number if the input argument is a number.
<i>nil</i>	Returns <code>nil</code> and an error message otherwise.

Examples

```
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("/net9"))`.

```
ln(100)  
=> 4.60517
```

Gets the natural logarithm of 100.

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

log10

```
log10( {o_waveform | n_number} )  
=> o_waveform/n_number/nil
```

Description

Gets the base-10 logarithm of a waveform or a number.

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>n_number</i>	Number.

Value Returned

<i>o_waveform</i>	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.
<i>n_number</i>	Returns a number that is calculated as the base-10 logarithm of the input number.
<i>nil</i>	Returns <code>nil</code> and an error message otherwise.

Examples

```
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.

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

lsb

```
lsb( o_s11 o_s12 o_s21 o_s22 g_frequency )  
    => o_waveform/nil
```

Description

Computes the load stability circles.

Arguments

<i>o_s11</i>	Waveform object representing s11.
<i>o_s12</i>	Waveform object representing s12.
<i>o_s21</i>	Waveform object representing s21.
<i>o_s22</i>	Waveform object representing s22.
<i>g_frequency</i>	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, <code>list(100M 1G 100M)</code> specifies a linear range with the following values: <pre>{ 100M, 200M, 300M, 400M, 500M, 600M, 700M, 800M, 900M, 1G }</pre> In that case, a load stability circle is calculated at each one of the 10 frequencies

Value Returned

<i>o_waveform</i>	Waveform object representing the load stability circles.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

Examples

```
plot(lsb(s11 s12 s21 s22 list(800M 1G 100M)))
```

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

lshift

```
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 rshift command.

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>n_delta</i>	Value by which the waveform is to be shifted.

Value Returned

<i>o_waveform</i>	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.
<i>nil</i>	Returns <code>nil</code> 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.

OCEAN Reference

Predefined Functions 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

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>n_number</i>	Number.

Value Returned

<i>o_waveform</i>	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.
<i>n_number</i>	Returns a number if the input argument is a number.
<code>nil</code>	Returns <code>nil</code> and an error message otherwise.

Examples

```
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`.

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Predefined Functions and Waveform (Calculator) Functions

nc

```
nc( o_Fmin o_Gmin o_rn g_level g_frequency )  
    => o_waveform/nil
```

Description

Computes the noise circles.

Arguments

<i>o_Fmin</i>	Waveform object representing the minimum noise factor.
<i>o_Gmin</i>	Waveform object representing the optimum noise reflection.
<i>o_rn</i>	Waveform object representing the normalized equivalent noise resistance.
<i>g_level</i>	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 <code>linRg</code> function can be called to generate a linear range. For example, <code>linRg(-30 30 5)</code> is the same as <code>list(-30 -25 -20 -15 -10 -5 0 5 10 15 20 25 30)</code> and the <i>g_level</i> argument can be specified as either of the above. In that case, a noise circle is calculated at each one of the 13 levels.
<i>g_frequency</i>	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, <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, a noise circle is calculated at each one of the 10 frequencies.

Value Returned

o_waveform Waveform object representing the noise circles.

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Predefined Functions and Waveform (Calculator) Functions

`nil` Returns `nil` and an error message otherwise.

Examples

```
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))
displayMode("smith")
smithType("impedance")
plot(NC)
```

overshoot

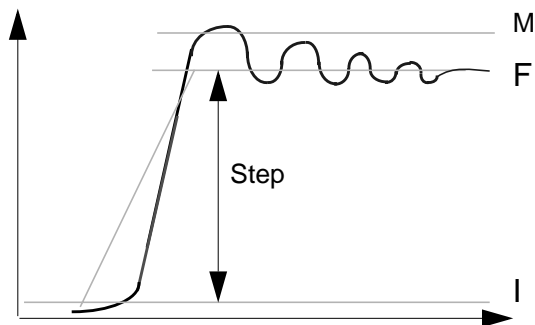
```
overshoot( o_waveform n_initVal g_initType n_finalVal g_finalType )  
=> o_waveform/n_value/nil
```

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.



$$\text{Overshoot} = \frac{(M - F) \times 100}{F - I}$$

Arguments

- | | |
|-------------------|--|
| <i>o_waveform</i> | Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .) |
| <i>n_initVal</i> | Initial X value at which to start the computation. |
| <i>g_initType</i> | Specifies how <i>initVal</i> functions.
Valid values: a non-nil value specifies that the initial value is taken to be the value of the waveform, interpolated at <i>initVal</i> , and the waveform is clipped from below, as follows:
<code>o_waveform = clip(o_waveform initVal nil)</code> |

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Predefined Functions and Waveform (Calculator) Functions

`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 `finalVal` functions.

Valid values: a non-`nil` value specifies that the final value is taken to be the value of the waveform, interpolated at `finalVal`, and the waveform is clipped from above, as follows:

```
o_waveform = clip( o_waveform nil finalVal )
```

`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`.)

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.

`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"`.

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Predefined Functions and Waveform (Calculator) Functions

peakToPeak

```
peakToPeak( o_waveform )  
=> o_waveform/n_value/nil
```

Description

Returns the difference between the maximum and minimum values of a waveform.

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
-------------------	--

Value Returned

<i>o_waveform</i>	Returns a waveform or a family of waveforms if the input argument is a family of waveforms.
<i>n_value</i>	Returns the difference between the maximum and minimum values of a waveform if the input argument is a single waveform.
<i>nil</i>	Returns <code>nil</code> 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.

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Predefined Functions and Waveform (Calculator) Functions

phase

```
phase( {o_waveform | n_number} )  
=> o_waveform/n_number/nil
```

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: `drwave: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.

Examples

```
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
```

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Predefined Functions and Waveform (Calculator) Functions

Gets the phase of -2 .

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Predefined Functions and Waveform (Calculator) Functions

phaseDeg

```
phaseDeg( {o_waveform | n_number} )  
=> o_waveform/n_number/nil
```

Description

Calculates the wrapped phase in degrees of a waveform and returns a waveform.

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>n_number</i>	Number.

Value Returned

<i>o_waveform</i>	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.
<i>n_number</i>	Returns a number if the input argument is a number.
<code>nil</code>	Returns <code>nil</code> 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.

phaseDegUnwrapped

```
phaseDegUnwrapped( {o_waveform | n_number} )  
=> o_waveform/n_number/nil
```

Description

Calculates the unwrapped phase in degrees of a waveform and returns a waveform.

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>n_number</i>	Number.

Value Returned

<i>o_waveform</i>	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.
<i>n_number</i>	Returns a number if the input argument is a number.
<i>nil</i>	Returns <code>nil</code> 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.

phaseMargin

```
phaseMargin( o_waveform )  
=> o_waveform/n_value/nil
```

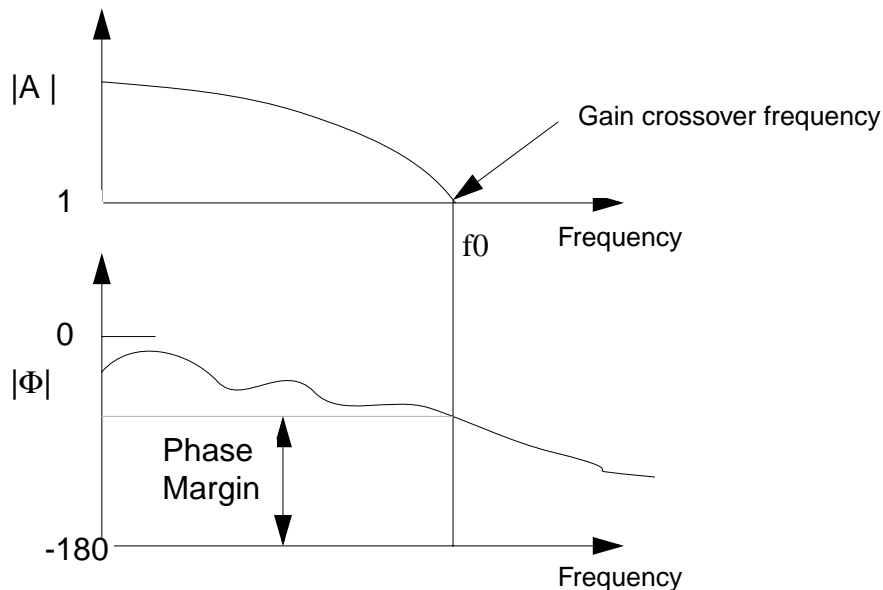
Description

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.

$\text{phaseMargin}(\text{gain}) = 180 + \text{phase}(\text{value}(\text{gain } f_0))$

The phase margin is calculated as the difference between the phase of the gain in degrees at f_0 and at -180 degrees. The frequency f_0 is the lowest frequency where the gain is 1. For stability, the phase margin must be positive.



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: `drwave:XXXXX`.)

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Predefined Functions and Waveform (Calculator) Functions

Value Returned

<i>o_waveform</i>	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.
<i>n_value</i>	Returns the value (in degrees) equivalent to the phase margin of the input waveform.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

Example

```
phaseMargin( v( "/OUT" ) )
```

Returns the phase margin for the waveform representing the voltage of the "/OUT" net.

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Predefined Functions and Waveform (Calculator) Functions

phaseRad

```
phaseRad( {o_waveform | n_number} )  
=> o_waveform/n_number/nil
```

Description

Calculates the wrapped (discontinuous) phase in radians of a waveform.

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX.</code>)
<i>n_number</i>	Number.

Value Returned

<i>o_waveform</i>	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.
<i>n_number</i>	Returns a number when the input argument is a number.
<i>nil</i>	Returns <code>nil</code> and an error message otherwise.

Example

```
plot( phaseRad( v( "/OUT" ) ) ) )
```

Returns the wrapped phase of the waveform representing the voltage of the `"/OUT"` net.

phaseRadUnwrapped

```
phaseRadUnwrapped( o_waveform )  
=> o_waveform/nil
```

Description

Calculates the unwrapped (continuous) phase in radians of a waveform and returns a waveform.

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
-------------------	--

Value Returned

<i>o_waveform</i>	Returns a waveform representing 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.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

Example

```
plot( phaseRadUnwrapped( v( "/OUT" ) ) )
```

Returns the unwrapped phase of the waveform representing the voltage of the "/OUT" net.

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Predefined Functions and Waveform (Calculator) Functions

pow

```
pow( {o_waveformBase | n_numberBas} {o_waveformExpn | n_numberExpn} )  
=> o_waveform/n_result/nil
```

Description

Takes the exponent of a given waveform or number.

Arguments

<i>o_waveformBase</i>	Waveform object to be used as the base for the expression.
<i>o_waveformExpn</i>	Waveform object to be used as the exponent for the expression.
<i>n_numberBase</i>	Number to be used as the base for the expression.
<i>n_numberExpn</i>	Number to be used as the exponent for the expression.

Value Returned

<i>o_waveform</i>	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).
<i>n_result</i>	Returns a number if both the input arguments are numbers.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

Examples

```
pow( average( v( "/net9" ) ) 0.5 )
```

Gets the square root of the average value of the voltage at "/net9".

```
pow( 2 3 )  
=> 8
```

Gets the value of 2 to the third power, or 8.

```
pow( -2 2 )  
=> 4
```

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Predefined Functions and Waveform (Calculator) Functions

Gets the value of -2 to the second power.

```
pow( 2.5 -1.2 )  
=> 0.3330213
```

Gets the value of 2.5 to the power of -1.2.

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Predefined Functions 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*.

Arguments

<i>o_waveform</i>	Time domain waveform object with units of volts or amps.
<i>f_timeStart</i>	Starting time for the spectral analysis interval. Use this parameter and <i>f_timeEnd</i> to exclude part of the interval. For example, you might set these values to discard initial transient data.
<i>f_timeEnd</i>	Ending time for the spectral analysis interval.
<i>x_num</i>	The number of time domain points to use. The maximum frequency in the Fourier analysis is proportional to <i>x_num</i> and inversely proportional to the difference between <i>f_timeStart</i> and <i>f_timeEnd</i> . Default value: 512
<i>t_windowName</i>	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 or 'Triangular. Default value: 'Hanning
<i>x_smooth</i>	The Kaiser window smoothing parameter. The 0 value requests no smoothing. Valid values: 0 <= <i>x_smooth</i> <= 15. Default value: 1

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Predefined Functions and Waveform (Calculator) Functions

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

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

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.

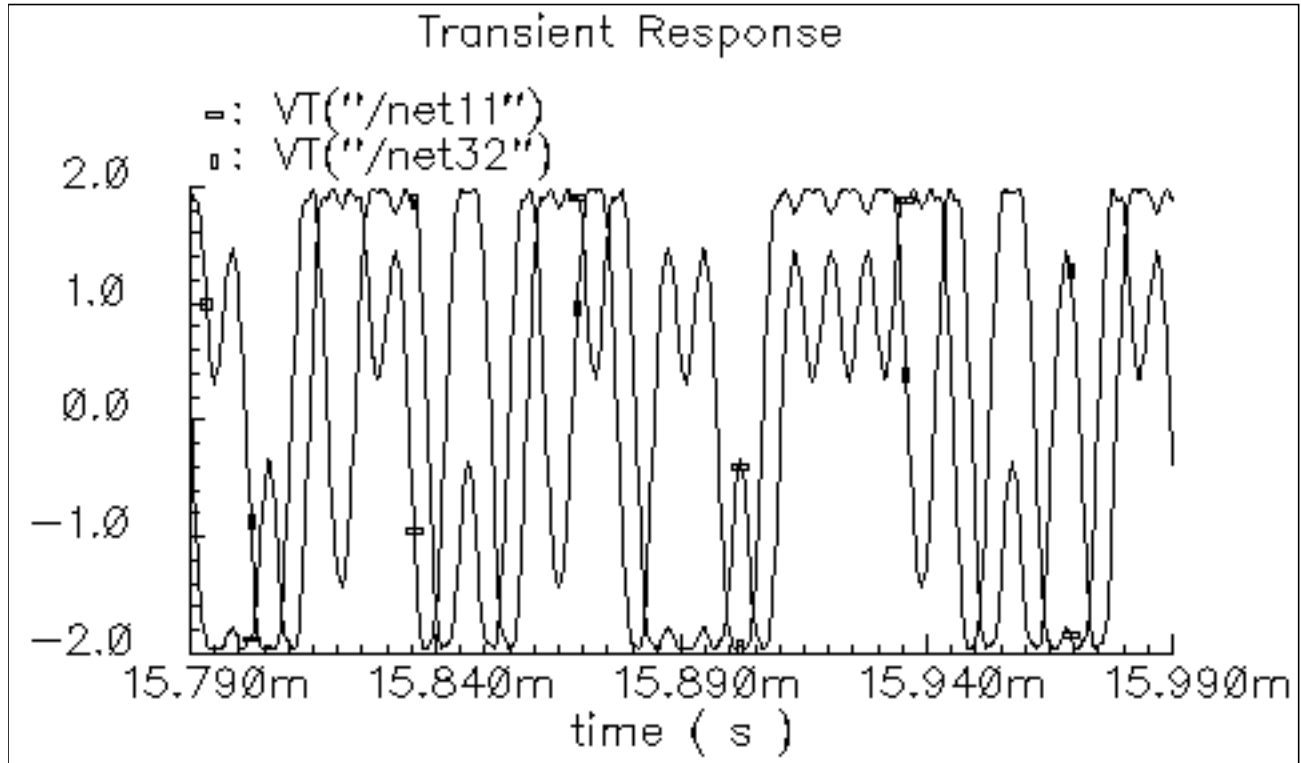
nil Returns `nil` when the command fails.

Example

```
psd(VT("/net32" "/hm/test_bench/spectre/schematic"), 0, 16m, 12000,  
    ?windowName 'Hanning,?smooth 1, ?windowSize 256,  
    ?detrending 'None, ?cohGain 1)
```

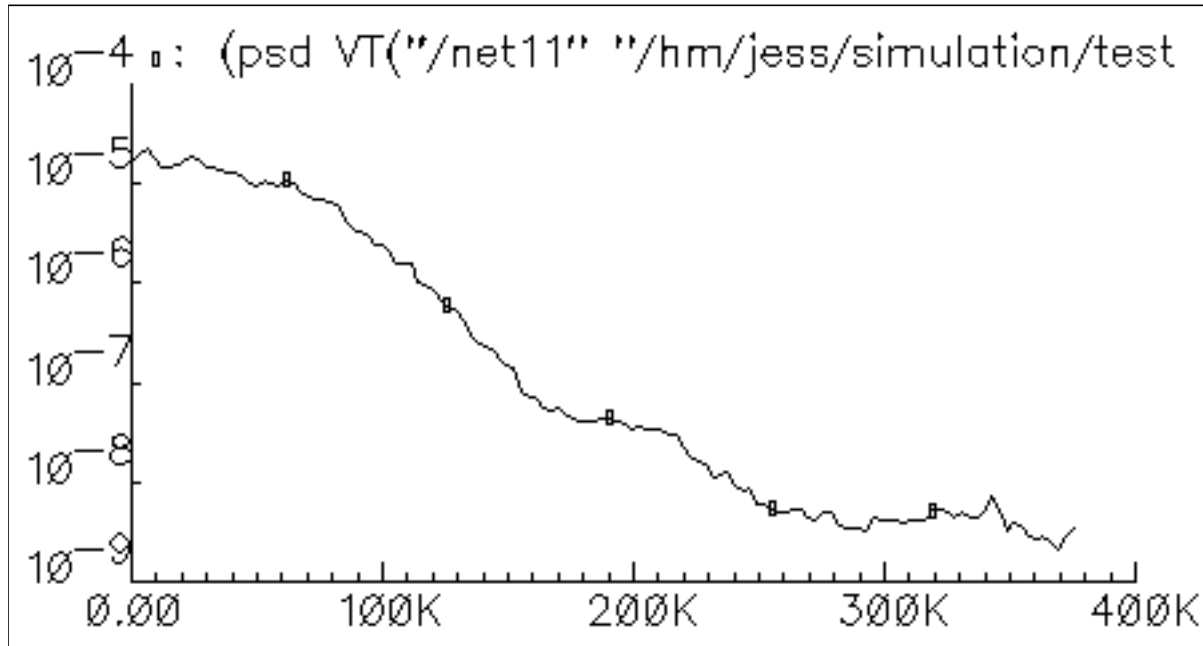
OCEAN Reference
Predefined Functions and Waveform (Calculator) Functions

Consider applying this command to one of the waveforms in the following illustration.



OCEAN Reference
Predefined Functions and Waveform (Calculator) Functions

The result is the following spectrum, which is displayed with a logarithmic vertical scale.



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Predefined Functions 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 $x_windowSize$.

Arguments

<i>o_waveform1</i>	Time domain waveform object with units of volts or amps.
<i>o_waveform2</i>	Time domain waveform object with units of volts or amps.
<i>f_timeStart</i>	Starting time for the spectral analysis interval. Use this parameter and <i>f_timeEnd</i> to exclude part of the interval. For example, you might set these values to discard initial transient data.
<i>f_timeEnd</i>	Ending time for the spectral analysis interval.
<i>x_num</i>	The number of time domain points to use. The maximum frequency in the Fourier analysis is proportional to <i>x_num</i> and inversely proportional to the difference between <i>f_timeStart</i> and <i>f_timeEnd</i> .
<i>t_windowName</i>	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 or 'Triangular. Default value: 'Hanning
<i>x_smooth</i>	The Kaiser window smoothing parameter. 0 requests no smoothing. Valid values: $0 \leq x_smooth \leq 15$. Default value: 1

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Predefined Functions and Waveform (Calculator) Functions

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

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.

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.

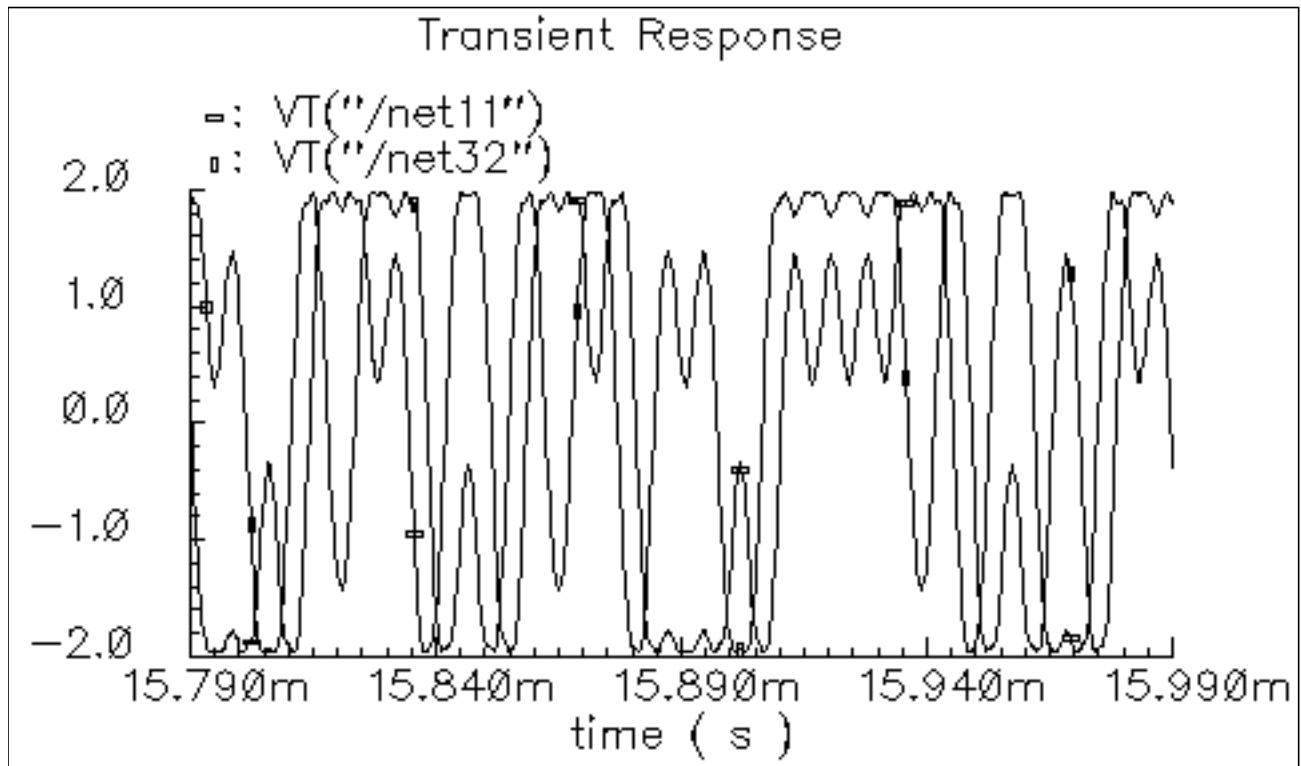
nil Returns `nil` when the command fails.

Example

```
psdbb(VT("/net32" "/hm/test_bench/spectre/schematic"),  
      VT("/net11" "/hm/test_bench/spectre/schematic"), 0, 16m, 12000,  
      ?windowName 'Hanning,?smooth 1, ?windowSize 256,  
      ?detrending 'None, ?cohGain 1)
```

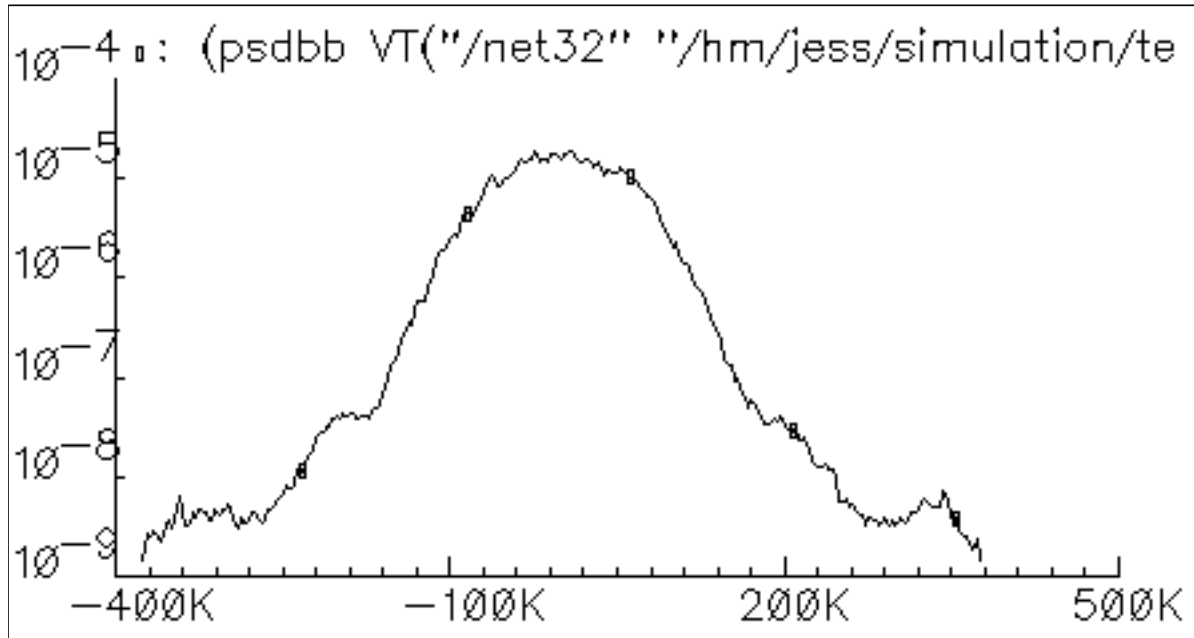

OCEAN Reference
Predefined Functions and Waveform (Calculator) Functions

Consider applying this command to both of the waveforms in the following illustration.



OCEAN Reference
Predefined Functions and Waveform (Calculator) Functions

The result is the following spectrum, which is displayed with a logarithmic vertical scale.



OCEAN Reference

Predefined Functions 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

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>n_input</i>	Complex number.

Value Returned

<i>o_waveformReal</i>	Returns a waveform when the input argument is a waveform.
<i>n_numberReal</i>	Returns a number when the input argument is a number.
<i>nil</i>	Returns <code>nil</code> 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 `x` representing a complex number, and returns the real portion of that complex number.

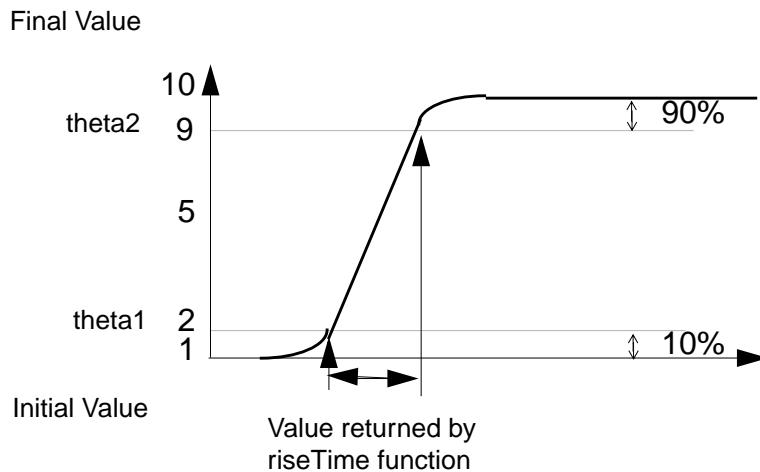
riseTime

```
riseTime( o_waveform n_initVal g_initType n_finalVal g_finalType n_theta1
          n_theta2 )
=> 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*.



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: *drwave: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:

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Predefined Functions and Waveform (Calculator) Functions

```
o_waveform = clip( o_waveform g_initVal nil )
```

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:

```
o_waveform = clip( o_waveform nil n_finalVal )
```

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.

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.

n_value Returns a value for the rise time if the input is a single waveform.

nil Returns *nil* and an error message otherwise.

Examples

```
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.

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Predefined Functions 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: `drwave: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.

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Predefined Functions 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.

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Predefined Functions 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 (*rootVal*).

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <i>drwave:XXXXX</i> .)
<i>n_rootVal</i>	Y value of interest.
<i>x_n</i>	Number that specifies which X value to return. If <i>n</i> equals 1, the first X value that crosses over the Y <i>rootVal</i> is returned. If <i>n</i> equals 2, the second X value that crosses over the Y <i>rootVal</i> is returned, and so on. If you specify a negative integer for <i>n</i> , the X values that cross the <i>rootVal</i> are counted from right to left (from maximum to minimum). If you specify <i>n</i> as 0, the list of root values is returned.

Value Returned

<i>o_waveform</i>	Returns a waveform if the input argument is a family of waveforms.
<i>n_value</i>	Returns an X value when the input argument is a single waveform.
<i>l_value</i>	Returns a list of all the root values when <i>n</i> is 0.
<i>nil</i>	Returns <i>nil</i> and an error message otherwise.

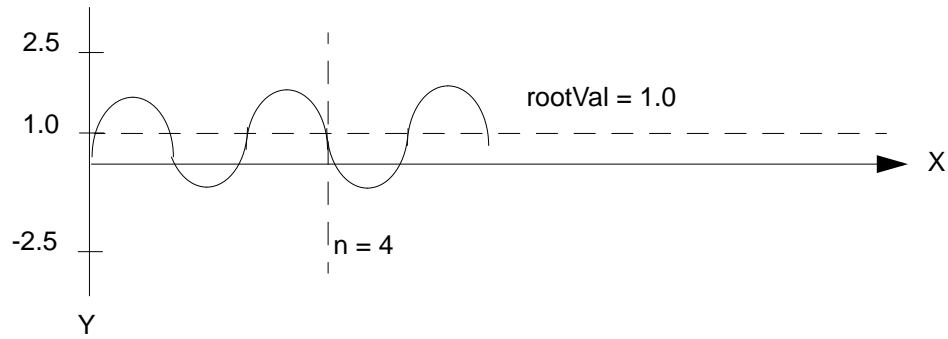
Example

```
root( v( "vout" ), 1.0, 4 )
```


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Predefined Functions 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.



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Predefined Functions 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: *drwave:XXXXX*.)

n_delta Value by which the waveform is to be shifted.

Value Returned

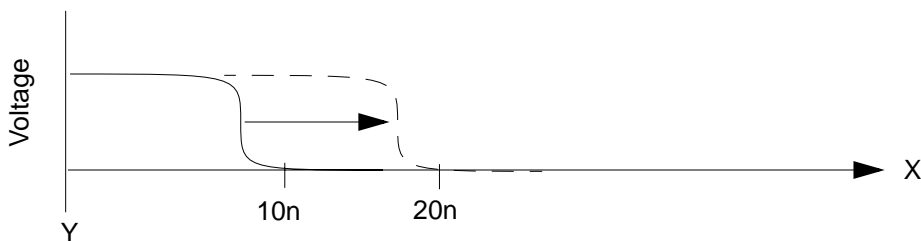
o_waveform Returns a waveform object. Returns a family of waveforms if the input argument is a family of waveforms.

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.



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Predefined Functions 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).

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>n_from</i>	Starting value for the sampling.
<i>n_to</i>	Ending value for the sampling.
<i>t_type</i>	Type of the sampling. Valid values: "linear" or "log"
<i>n_by</i>	Interval at which to sample.

Value Returned

<i>o_waveform</i>	Returns a waveform representing the sampling you specified.
<i>n_number</i>	Returns a number if the output contains only one point.
<i>nil</i>	Returns <code>nil</code> and an error message otherwise.

Examples

```
sample( v( "vout" ) 0 50n "linear" 0.1n )
```

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Predefined Functions and Waveform (Calculator) Functions

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.

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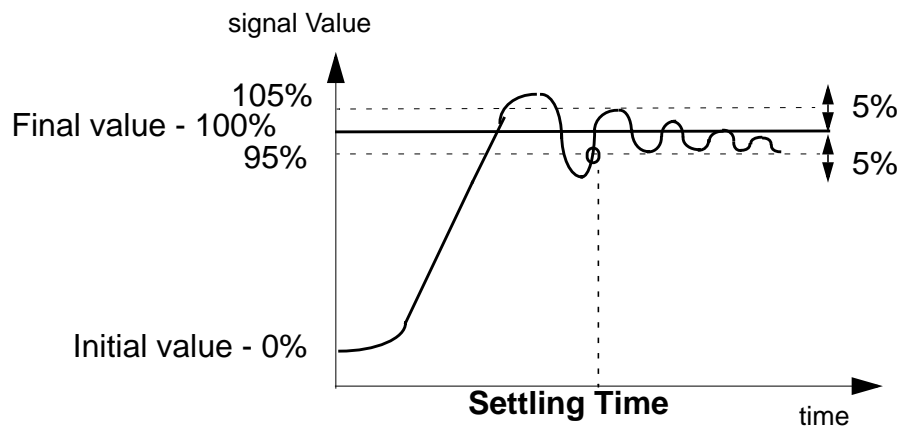
Predefined Functions and Waveform (Calculator) Functions

settlingTime

```
settlingTime( o_waveform n_initVal g_initType n_finalVal g_finalType n_theta )  
=> 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%.

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>n_initVal</i>	Initial value at which to start the computation.
<i>g_initType</i>	Specifies whether the values entered are X values or Y values. Valid values: <code>t</code> specifies that <i>initVal</i> is defined by the X value entered; <code>nil</code> specifies that <i>initVal</i> is defined by the Y value entered
<i>n_finalVal</i>	Final value at which to start the computation.

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Predefined Functions and Waveform (Calculator) Functions

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.

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.

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 )
```

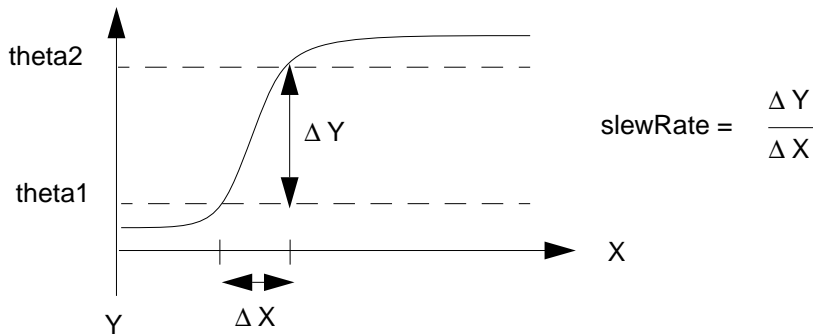
Computes the time required for the waveform representing the voltage of the `/out` net to settle within 90 percent of the step from 0 to 2.

slewRate

```
slewRate( o_waveform n_initVal g_initType n_finalVal g_finalType n_theta1  
          n_theta2 )  
=> 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.



Arguments

- | | |
|--------------------|---|
| <i>o_waveform</i> | Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <i>drwave:XXXXX</i> .) |
| <i>n_initVal</i> | Initial X-axis value at which to start the computation. |
| <i>g_initType</i> | Specifies whether the values entered are X values or Y values. Valid values: <i>t</i> specifies that <i>initVal</i> is defined by the X value entered; <i>nil</i> specifies that <i>initVal</i> is defined by the Y value entered |
| <i>n_finalVal</i> | Final value at which to end the computation. |
| <i>g_finalType</i> | Specifies whether the values entered are X values or Y values. Valid values: <i>t</i> specifies that <i>finalVal</i> is defined by the X value entered; <i>nil</i> specifies that <i>finalVal</i> is defined by the Y value entered |
| <i>n_theta1</i> | Percent low (percentage of the total step). |

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Predefined Functions and Waveform (Calculator) Functions

n_theta2 Percent high (percentage of the total step).

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.

n_value Returns a value for the slew rate for the specified waveform if the input is a single waveform.

nil Returns *nil* and 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.

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Predefined Functions 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

<i>o_current</i>	Waveform representing the current. The current can be obtained by calling the <code>i</code> data access function for the desired terminal.
<i>o_voltage</i>	Waveform representing the voltage. The voltage can be obtained by calling the <code>v</code> data access function for the desired net. To obtain meaningful results, the terminal used to obtain the current must be a member of the net used to obtain the voltage.

Value Returned

<i>o_power</i>	Waveform representing the power of the net.
<code>nil</code>	Returns <code>nil</code> if there is an error.

Example

```
plot(dbl10(spectralPower(i("/PORT0/PLUS") v("/net28"))))
```

Plots power of the output `"/net28"`. `"/PORT0/PLUS"` is a member of `"/net28"`.

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Predefined Functions 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.

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)))
```

stddev

```
stddev( o_waveform )  
=> 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 $y=f(x)$

$$stddev(y) = \sqrt{\frac{\int_{from}^{to} (y - average(y))^2}{to - from}}$$

Arguments

<i>o_waveform</i>	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: drwave:XXXXX)
-------------------	--

Value Returned

<i>n_stddev</i>	Returns a number representing the standard deviation value of the input waveform.
<i>o_waveformStddev</i>	Returns a waveform representing the average value if the input is a family of waveforms.
<i>nil</i>	Returns <i>nil</i> 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.

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Predefined Functions and Waveform (Calculator) Functions

tangent

```
tangent( o_waveform [ ?x n_x ] [ ?y n_y ] [ ?slope n_slope ] )  
=> o_waveform/nil
```

Description

Returns the tangent to a waveform through the point (n_x , n_y) with the given slope.

Arguments

<i>o_waveform</i>	Waveform object representing the wave.
<i>n_x</i>	X coordinate of the point. The default value is the X coordinate of the first point on the wave.
<i>n_y</i>	Y coordinate of the point. The default value is the Y coordinate at the given or default X coordinate.
<i>n_slope</i>	Slope of the line. Default value: 1.0

Value Returned

<i>o_waveform</i>	Wave object representing the line.
<i>nil</i>	Returns <i>nil</i> if there is an error.

Example

```
refLine  
=> tangent(refWave ?x -25 ?slope 1.0)
```

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Predefined Functions 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.

The computation uses the `dft` function. Assume that the `dft` function returns complex coefficients $A_0, A_1, \dots, A_f, \dots$. 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 i th harmonic where $i \neq 0, f$. Then, total harmonic distortion is computed as:

$$\frac{\sqrt{\sum_{i=1, i \neq 0, f} |A_i|^2}}{|A_f|} \times 100\%$$

Arguments

<code>o_waveform</code>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<code>n_from</code>	Starting value for the computation.
<code>n_to</code>	Ending value for the computation.
<code>x_num</code>	Number of timepoints. If <code>x_num</code> is not a power of 2, it is forced to be the next higher power of 2.
<code>n_fund</code>	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.

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Predefined Functions and Waveform (Calculator) Functions

Value Returned

<i>o_waveform</i>	Returns a waveform representing the absolute value of the total harmonic distortion if the input argument is a family of waveforms.
<i>n_thdValue</i>	Returns the absolute value of the total harmonic distortion of the input waveform.
<i>nil</i>	Returns <i>nil</i> 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.

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Predefined Functions and Waveform (Calculator) Functions

value

```
value( o_waveform [s_name] g_value )  
      => o_waveform/g_value/nil
```

Description

Returns the Y value of a waveform for a given X value.

Arguments

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
<i>s_name</i>	The name of the innermost or outermost sweep variable. If the sweep variable name is not supplied, the innermost sweep variable is used.
<i>g_value</i>	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.

Value Returned

<i>o_waveform</i>	Returns a waveform or a family of waveforms if the input argument is a family of waveforms.
<i>g_value</i>	Returns the Y value if the input argument is a single waveform. Note: For parametric sweeps, the value might be a waveform that can be printed with the <code>ocnPrint</code> command.
<i>nil</i>	Returns <code>nil</code> and an error message if the value cannot be printed.

Examples

```
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.

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

```
value( v( "/OUT" )'TEMPDC 20.0 )
```

Returns `drwave: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`.

OCEAN Reference

Predefined Functions 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: `drwave:XXXXX`.)

x_numberOfPeaks Specifies the *n*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.

Examples

```
xmax( v( "/net9" ) 1 )
```

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

Gets the time value (X-axis value) at which the voltage of `"/net9"` attains its first peak value.

```
xmax( v( "/net9" ) 0 )
```

Gets the list of time values (X-axis values) at which the voltage of `"/net9"` attains each of its peak values.

OCEAN Reference

Predefined Functions 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: `drwave: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.

Examples

```
xmin( v( "/net9" ) 1 )
```

OCEAN Reference

Predefined Functions 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.

OCEAN Reference

Predefined Functions 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

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX.</code>)
-------------------	--

Value Returned

<i>o_waveform</i>	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.
-------------------	--

<i>nil</i>	Returns <i>nil</i> 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.

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

y_{max}

```
ymax( owaveform )  
=> nmax/owaveformMax/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: `drwave:XXXXX`.)

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"`.

OCEAN Reference

Predefined Functions and Waveform (Calculator) Functions

ymin

```
ymin( o_waveform )  
=> 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

<i>o_waveform</i>	Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: <code>drwave:XXXXX</code> .)
-------------------	--

Value Returned

<i>n_min</i>	Returns a number representing the minimum value of Y if the input argument is a single waveform.
<i>o_waveformMin</i>	Returns a waveform (or family of waveforms) representing the minimum value of Y if the input argument is a family of waveforms.
<i>nil</i>	Returns <code>nil</code> and an error message otherwise.

Example

```
ymin( v( "/net9" ) )
```

Gets the minimum voltage (Y value) of `"/net9"`.

OCEAN Reference
Predefined Functions and Waveform (Calculator) Functions

Advanced Analysis

The OCEAN commands for advanced analyses let you run parametric analysis, corners analysis, Monte Carlo analysis, and Optimization. This chapter includes setup commands for these analyses and the special data-access or plot commands that are used for these analyses.

The following sections contain the commands and other information relating to advanced analyses.

- [Parametric Analysis Commands](#) on page 377
- [Corners Analysis Commands](#) on page 383
- [Monte Carlo Analysis Commands](#) on page 391
- [Optimization Commands](#) on page 413

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.

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]  
              [?values l_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.

<i>n_start, n_stop</i>	highest precedence
	↓
<i>n_center, n_span</i>	
<i>n_center, n_spanPercent</i>	lowest precedence

Similarly, if you specify more than one step control, the OCEAN environment uses the following precedence.

<i>f_step</i>	highest precedence
	↓
<i>n_lin</i>	
<i>n_dec</i>	
<i>n_log</i>	
<i>n_oct</i>	
<i>n_times</i>	lowest precedence

To run the analysis, use the `paramRun` command described in “[paramRun](#)” on page 382.

Arguments

<i>t_desVar</i>	Name of the design variable to be swept.
<i>n_start</i>	Beginning value for the design variable.

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<i>n_stop</i>	Final value for the design variable.
<i>n_center</i>	Center point for a range of values that you want to sweep.
<i>n_span</i>	Range of values that you want to sweep around the center point. For example, if <i>n_center</i> is 100 and <i>n_span</i> is 20 then the sweep range extends from 90 to 110.
<i>f_step</i>	Increment by which the value of the design variable changes. For example, if <i>n_start</i> is 1.0, <i>n_stop</i> is 2.1, and <i>f_step</i> is 0.2, the parametric analyzer simulates at values 1.0, 1.2, 1.4, 1.6, 1.8, and 2.0.
<i>n_lin</i>	<p>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 <i>n_start</i> and <i>n_stop</i>. The value for the <i>n_lin</i> argument must be an integer greater than 0.</p> <p>For example, if <i>n_start</i> is 0.5, <i>n_stop</i> is 2.0, and <i>n_lin</i> is 4, the parametric analyzer simulates at values 0.5, 1.0, 1.5, and 2.0.</p>
<i>n_log</i>	<p>The number of steps between the starting and stopping points at equal-ratio intervals using the following formula:</p> $\log \text{ multiplier} = (n_{\text{stop}}/n_{\text{start}})^{(n_{\text{log}} - 1)}$ <p>The number of steps can be any positive number, such as 0.5, 2, or 6.25. Default value: 5</p> <p>For example, if <i>n_start</i> is 3, <i>n_stop</i> is 15, and <i>n_log</i> is 5, the parametric analyzer simulates at values 3, 4.48605, 6.7082, 10.0311, and 15.</p> <p>The ratios of consecutive values are equal, as shown below.</p> $3/4.48605 = 4.48605/6.7082 = 6.7082/10.0311 = 10.0311/15 = .67$
<i>n_dec</i>	The number of steps between the starting and stopping points calculated using the following formula:

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$$\text{decade multiplier} = 10^{1/n_dec}$$

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 .

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 .

$$\text{octave multiplier} = 2^{1/(n_oct)}$$

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.

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.

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l_values List of values to be swept. You can use *l_values* by itself or in conjunction with *n_start*, *n_stop*, and *f_step* to specify the set of values to sweep.

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.

Examples

```
paramAnalysis( "rs" ?start 200 ?stop 1000 ?step 200
  ?values '(1030 1050 1090) )
```

Sets up a parametric analysis for the `rs` design variable. The swept values are 200, 400, 600, 800, 1000, 1030, 1050, and 1090.

```
paramAnalysis( "r1" ?start 200 ?stop 600 ?step 200
  paramAnalysis( "rs" ?start 300 ?stop 700 ?step 200
  )
)
```

Sets up a nested parametric analysis for the `r1` design variable.

```
paramAnalysis("temp" ?start -50 ?stop 100 ?step 50)
```

Sets up a parametric analysis for temperature.

paramRun

```
paramRun( [s_paramAnalysis] )  
=> t/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 [“paramAnalysis”](#) on page 378.

Arguments

s_paramAnalysis Parametric analysis.

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.

Corners Analysis Commands

The corners analysis commands let you set up and run analyses to measure circuit performance with respect to variations in a semiconductor manufacturing process. This section lists the commands that you can use to configure and run corners analyses in the OCEAN environment. The following manuals provide more information on corners analysis.

- [Advanced Analysis Tools User Guide](#)
- [Virtuoso® Analog Design Environment SKILL Language Reference](#)

The corners analysis commands follow.

cornerDesVar

```
cornerDesVar(t_cornerName t_desVarName t_value)  
=> t/nil
```

Description

Sets the design variable value for the specified corner.

Arguments

<i>t_cornerName</i>	Name of the corner.
<i>t_desVarName</i>	Name of the design variable.
<i>t_value</i>	Value of the design variable.

Value Returned

t	Returned if successful.
nil	Returns nil and prints an error message.

Example

```
cornerDesVar("slow" "vcc" "5")
```

Sets the value of `vcc` to 5 for corner `slow`.

cornerMeas

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

Description

Displays all the predefined enabled measurements from a Design Customization file, either graphically (plot) or textually (print), according to your choices.

Each measurement is plotted or printed in a separate subwindow.

Arguments

None.

Value Returned

t	Returned if successful.
nil	Returns nil and prints an error message.

Example

```
cornerMeas()
```

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Advanced Analysis

cornerRun

```
cornerRun( [t_cornerName1 t_cornerName2 ...] [?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] )
=> t/s_jobName/nil
```

Description

Runs the corner analysis that has been predefined in the .pcf and .dcf files and selected via the `selectProcess` command. If specific corners are specified, only those corners run; otherwise all the corners run.

You can load your .pcf and .dcf files with the `loadPcf` and `loadDcf` commands. See the Virtuoso® Analog Design Environment SKILL Language Reference for information on these commands.

Arguments

t_cornerName A specific corner to be run. If you do not specify one or more corners, then all the enabled corners run.

Note: The following arguments are valid only when running in distributed processing mode.

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.

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.

t_queueName Name of the queue. If no queue is defined, the analysis is placed in the default queue (your home machine).

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 is not completed by *t_termTime*, the job is terminated.

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t_dependentOn

List of jobs on which the specified analysis is dependent. The analysis is not started until after dependent jobs are complete.

t_mailingList

List of users to be notified by e-mail when the analysis is complete.

s_block

When *s_block* is not `nil`, the OCEAN script halts until the job is complete.

Default value: `nil`

s_notifyFlag

When *notifyFlag* is not `nil`, a job completion message is echoed to the OCEAN interactive window.

Default value: `t`

Value Returned

`t`

Returned if successful.

s_jobName

For a distributed process, the job name specified or assigned by the system to the analysis.

`nil`

Returns `nil` and prints an error message.

Examples

```
cornerRun( )
```

Runs all corners analysis defined in the `.pcf` and `.dcf` files and selected by the `selectProcess` command.

```
cornerRun( ?startTime 10 ?host "mach14" ?mail "preampGroup" )
```

Runs all corners analysis defined in the `.pcf` and `.dcf` files and selected by the `selectProcess` command in distributed mode with a *starttime* of 10, using `mach14` as *host*, and notifying the mail group `preampGroup` when the analysis is complete.

cornerRunTemp

```
cornerRunTemp(t_cornerName t_value)  
=> t/nil
```

Description

Sets the analysis temperature (in degrees Celsius) to be used for a corner.

Arguments

<i>t_cornerName</i>	Name of the corner.
<i>t_value</i>	Temperature value in degrees Celsius.

Value Returned

t	Returned if successful.
nil	Returns nil and prints an error message.

Example

```
cornerRunTemp("slow" "50")
```

Sets the temperature to 50 for corner `slow`.

residual

```
residual(x_scalarExpression ?upper x_upperValue ?target x_targetValue  
        ?lower x_lowerValue)  
=> t/nil
```

Description

Creates a residual plot of the given scalar expression given the upper and lower performance bounds and target.

Arguments

<i>x_scalarExpression</i>	Scalar expression from a corners analysis.
<i>x_upperValue</i>	Upper performance bound.
<i>x_targetValue</i>	Target value.
<i>x_lowerValue</i>	Lower performance bound.

Value Returned

t	Returned if successful.
nil	Returns nil and prints an error message.

Example

```
residual( bandwidth(v("net1"), 3, "low") ?upper 5 ?target 2.5 ?lower 0)
```

Creates a residual plot of `v("net1")` with an upper boundary of 5, a target of 2.5, and a lower boundary of 0.

selectProcess

```
selectProcess(t_processName)  
=> t/nil
```

Description

Selects one of the processes already loaded with a `loadPcf` or `loadDcf` command.

Arguments

<i>t_processName</i>	Name of the process, as specified in the <code>.pcf</code> or <code>.dcf</code> file with the <code>corAddProcess</code> function.
----------------------	--

Value Returned

<code>t</code>	Returned if successful.
<code>nil</code>	Returns <code>nil</code> and prints an error message.

Example

```
selectProcess("fab6")
```

Selects the process `fab6`.

Monte Carlo Analysis Commands

The commands for running Monte Carlo in the OCEAN environment are as follows.

correlationTable

```
correlationTable(?suppress x_suppress)  
=> t/nil
```

Description

Prints the correlation between all pairs of declared `monteExpr` expressions.

Pairs of the same expression, which have a correlation value of 1.0, are excluded. This exclusion means that the `correlationTable` command prints only the off-diagonal terms in the correlation matrix.

Arguments

<i>x_suppress</i>	Suppresses the printing for correlations less than this value. Default value: .5
-------------------	---

Value Returned

t	Returned if successful.
nil	Returned otherwise.

Example

```
correlationTable()
```


dataFilter

```
dataFilter(t_monteExprName ( {?sigma x_sigma | ?upper x_upper ?lower x_lower}  
    ?filterBy s_filterBy ) )  
=> t/nil
```

Description

Eliminates bad data points (*outliers*) from a Monte Carlo data set.

Arguments

t_monteExprName

The `monteExpr` name with the appended swept parameter.

x_sigma

Filters data lying outside an established sigma point from the mean. For instance, you might filter data lying outside 3 standard deviations (sigma) from the mean. You can specify *x_sigma* or you can specify *x_upper* and *x_lower*, but you cannot specify both.

Default value: 3

x_upper

Filters data that is greater than an upper numerical limit.

Default value: `inf`

x_lower

Filters data that is less than a lower numerical limit.

Default value: `-inf`

s_filterBy

Type of filter to be used. This setting affects all of your data so you only need to specify the type of filter once.

Valid values: `'dataSet', 'point'`

Default value: `'dataSet'`

`'dataSet'` ignores all measurements for a point if the value of any of the measurements for that point is outside the filter limits.

`'point'` filters an outlying point only from the specific measurement that recorded the outlying point.

OCEAN Reference

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Value Returned

t	Returned if successful.
nil	Returned otherwise.

Example

```
dataFilter('bandwidth ?upper 10Mhz  
          ?lower 0.1Mhz ) ;For nominal 1Mhz  
dataFilter('bandwidth ?sigma 3 )
```

The second example sets the upper limit to
 $\text{mean}(\text{bandwidth}) + 3 * \text{sigma}(\text{bandwidth})$
and sets the lower limit to
 $\text{mean}(\text{bandwidth}) - 3 * \text{sigma}(\text{bandwidth})$

histogram

```
histogram( t_monteExprName ?type s_type ?numBins x_numBins ?density b_density )  
=> t/nil
```

Description

Plots a histogram of Monte Carlo data.

This command plots to an individual subwindow. The value of the *s_type* argument determines the style of the line. Setting *b_density* to `t` causes the `histogram` command to plot a smooth distribution curve for the data.

Arguments

t_monteExprName

The `monteExpr` name with the appended swept parameter.

s_type

Style of line to be used.

Valid values: 'standard, 'passFail, 'cumulativeLine,
'cumulativeBox

Default value: 'standard (if you do not specify *s_type*)

'standard prints a bar graph of the output versus parameter.

'passFail requires that specification limits be specified. This option plots a bar graph where the runs that pass are shown in green and the runs that fail are shown in red.

'cumulativeLine uses a *joined* line style to plot the cumulative distribution function. The cumulative distribution function is the area under the standard histogram bars.

'cumulativeBox plots the same information as the
'cumulativeLine option but uses a *bar* plotting style.

x_numBins

Number of bins to be used for the histogram.

Default value: 10

b_density

If set to `t`, plots the probability density function for the data. Valid values: `t` or `nil`.

OCEAN Reference

Advanced Analysis

Value Returned

t	Returned if successful.
nil	Returned otherwise.

Example

```
monteExpr( "bw" 'bandwidth( v("vout"),3,"low" ) )
monteExpr( "DCgain" 'ymax( vdb("vout") ) )
histogram( "bw_27" )
histogram( "bw_27" ?numBins 12 ?density t )
```

iterVsValue

```
iterVsValue( t_monteExprName ?outputFormat s_outputFormat )  
=> t/nil
```

Description

Prints the value of every scalar measurement for each Monte Carlo iteration.

Arguments

t_monteExprName

The `monteExpr` name with the appended swept parameter.

s_outputFormat

The output format for the printout.

Valid values: 'sorted, 'unsorted

Default value: 'sorted

'sorted sorts the output from highest to lowest value.

'unsorted prints the values without sorting.

Value Returned

t

Returned if successful.

nil

Returned otherwise.

Example

```
iterVsValue( "bw_27" )
```

monteCarlo

```
monteCarlo(  
  [?numIters x_numIters] [?startIter x_startIter]  
  [?analysisVariation s_analysisVariation] [?sweptParam t_sweptParam]  
  [?sweptParamVals l_sweptParamVals] [?saveData saveData] [?append b_append]  
)  
=> t/nil
```

Description

Sets up a Monte Carlo analysis.

To run the analysis, use the `monteRun` command described in [“monteRun”](#) on page 405.

Arguments

x_numIters Number of iterations (runs).
 Default value: 100

x_startIter Starting iteration.
 Default value: 1

Note: *x_startIter* must not be 1 when

- You want to append to existing data. For example, you run 100 Monte Carlo analyses and then want to run 100 more in addition to the previous 100. In this case, *x_startIter* must be 101. If *x_startIter* is 1, the same results are recalculated as before.
- You want to rerun a particular run. In this case, *startIter* must be the number of that particular run.

s_analysisVariation Analysis variations.
 Valid values: 'process, 'mismatch,
 'processAndMismatch
 Default value: 'process

t_sweptParam Design variable (or temperature) that can be swept with Monte Carlo.
 Default value: none (this is the inner loop)

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Advanced Analysis

t_sweptParamVals

List of values of *sweptParam*.

saveData

Indicates when to save data to allow family plots.
Default value: `nil`

b_append

Appends the new results to data from a previous Monte Carlo run.
Default value: `nil`

Value Returned

`t`

Returned if successful.

`nil`

Returned otherwise.

Example

```
monteCarlo()
```

Sets up a Monte Carlo analysis using some of the defaults.

```
monteCarlo(?numRuns 300 ?analysisType 'processAndMismatch ?sweptParam temp  
?sweptParamVals list(-50, 0, 50) ?nomRun n)
```

monteCorrelate

```
monteCorrelate( {t_param1 ... t_paramN | t_deviceName1 ... t_deviceNameN}  
               f_correlationValue)  
=> t/nil
```

Description

Specifies a correlation coefficient for a list of process parameters or a list of devices specified in individual subcircuits.

Use this command to specify matched pairs of devices or to specify mismatch of devices in excess of that specified for the process. You must not mix devices and parameters on the same command line.

Arguments

f_correlationValue

Value of the correlation coefficient that describes the correlation among the listed parameters or devices.

t_param1

Name of the first process parameter to be correlated.

t_paramN

Name of another process parameter to be correlated.

t_deviceName1

Name of the first device to be correlated.

t_deviceName2

Name of another device to be correlated.

Value Returned

t

Returned if successful.

nil

Returned otherwise.

Example

```
monteCorrelate("cje_27 bw_27" 0.8)
```


monteDisplay

```
monteDisplay()  
=> undefined/nil
```

Description

Displays the currently defined Monte Carlo analysis, including all expressions that are defined.

Arguments

None.

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

```
monteDisplay()
```

monteExpr

```
monteExpr( t_monteExprName s_expression )  
=> t/nil
```

Description

Sets up the Monte Carlo scalar expressions that are used to create the histogram file.

Arguments

<i>t_monteExprName</i>	Name of the expression.
<i>s_expression</i>	Expression.

Value Returned

t	Returned if successful.
nil	Returned otherwise.

Example

```
monteExpr( "bw" 'bandwidth( v(\ "net7\ " ) 3 \ "low\ " ) )
```

monteOutputs

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

Description

Returns the names of the `monteExpr` expressions, concatenating the `monteExprName` set in the `monteExpr` command with the value of the swept variable.

If no variable is swept, the `monteOutputs` command concatenates the default temperature to the `monteExprName`. For example, the returned name might be `bw_27`.

Arguments

None.

Value Returned

<code>t</code>	Returned if successful.
<code>nil</code>	Returned otherwise.

Example

```
monteOutputs()
```

monteResults

```
monteResults(?dataFileName t_scalarDataFile ?paramFileName t_parameterFile)  
=> t/nil
```

Description

Initializes the Monte Carlo data analysis tools.

The `monteResults` command reads in the specified data and parameter files, opens a new Waveform window, and adds a statistical analysis menu to the Waveform window. The menu items are equivalent to those found on the *Monte Carlo Results* menu in the Virtuoso® Analog Design Environment.

Arguments

t_scalarDataFile

Name of scalar data file to be read in.
Default value: `mcdData`

t_parameterFile

Name of parameter file associated with scalar data.
Default value: `mcParam`

Value Returned

`t` Returned if successful.

`nil` Returned otherwise.

Examples

```
monteResults()  
monteResults( ?dataFileName "myData" ?paramFileName "myParams" )
```

monteRun

```
monteRun(  
  [?jobName t_jobName] [?host t_hostName] [?tasks x_tasks]  
  [?queue t_queueName] [?startTime t_startTime] [?termTime t_termTime]  
  [?dependentOn t_dependentOn] [?mail t_mailingList] [?block s_block]  
  [?notify s_notifyFlag] )  
=> s_jobName/nil/t
```

Description

Runs a Monte Carlo analysis previously set up with the `monteCarlo` and `monteExpr` commands.

The `monteRun` command runs all the Monte Carlo processes defined in the `.pcf` and `.dcf` files. You can load your `.pcf` and `.dcf` files with the `loadPcf` and `loadDcf` commands. See the *Artist SKILL Language Reference Manual* for information on these commands.

Arguments

Note: Arguments to the `monteRun` command are valid only when running in distributed (processing) mode.

<code>t_jobName</code>	Used as the basis of the job name. The value entered for <code>t_jobName</code> is used as the job name and return value if the <code>run</code> command is successful. If the name given is not unique, a value is appended to create a unique job name.
<code>t_hostName</code>	Name of the host on which to run the analysis. If no host is specified, the system assigns the analysis to an available host.
<code>x_tasks</code>	Number of tasks in which to divide the Monte Carlo job. Default value: calculated from your setup
<code>t_queueName</code>	Name of the queue. If no queue is defined, the analysis is placed in the default queue (your home machine).
<code>t_startTime</code>	Desired start time for the job. If dependencies are specified, the job does not start until all dependencies are satisfied.
<code>t_termTime</code>	Termination time for job. If the job has not completed by <code>t_termTime</code> , the job is terminated.

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<i>t_dependentOn</i>	List of jobs on which the specified analysis is dependent. The analysis is not started until after dependent jobs are complete.
<i>t_mailingList</i>	List of users to be notified when the analysis is complete.
<i>s_block</i>	When <i>s_block</i> is not <code>nil</code> , the OCEAN script halts until the job is complete. Default value: <code>nil</code>
<i>s_notifyFlag</i>	When <i>s_notifyFlag</i> is not <code>nil</code> , a job completion message is echoed to the OCEAN interactive window. Default value: <code>t</code>

Value Returned

<code>t</code>	Returned if successful.
<code>nil</code>	Returned otherwise.
<i>s_jobName</i>	For a distributed process, the job name that the system specified or assigned to the analysis.

Example

```
monteRun( )
```

Runs all the Monte Carlo analyses defined in the `.pcf` and `.dcf` files.

monteSelectResults

```
monteSelectResults( ?mcdDataFileName t_mcdDataFileName  
  ?paramFileName t_paramFileName )  
=> t/nil
```

Description

Selects the specified mcdData file, which is the file that contains the scalar data.

Before you use this command, you must have access to mcdData and param files, either produced by an earlier successful Monte Carlo simulation or pointed to by a previous `openResults()` command.

Arguments

t_mcdDataFileName The name of the mcdData file.
Default value: `mcdData`

t_paramFileName The name of the param file.
Default value: `param`

Value Returned

`t` Returned if successful.

`nil` Returned otherwise.

Example

```
monteSelectResults()  
monteSelectResults(?mcdDataFileName mcdDataRun2  
  ?paramFileName paramRun2)
```

scatterplot

```
scatterplot(t_monteExprName_X t_monteExprName_Y ?bestFit b_bestFit )  
=> t/nil
```

Description

Plots different statistical measurements against each other so you can determine whether there is a relationship between two parameters.

Tightly correlated parameters show linear relationships.

Arguments

t_monteExprName_X

The `monteExpr` name with the appended swept parameter for the X-axis variable.

t_monteExprName_Y

The `monteExpr` name with the appended swept parameter for the Y-axis variable.

b_bestFit

If `t`, the `scatterplot` command computes and draws on the plot the best fitting straight line through the data. The best line is defined as the line that minimizes the sum of squares of the distances between the data points and the line.

Value Returned

`t` Returned if successful.

`nil` Returned otherwise.

Example

```
monteExpr( "bw" 'bandwidth( v("vout"), 3, "low" ) )  
monteExpr( "DCgain" 'ymax( vdb("vout") ) )  
scatterplot( "bw_27" "DCgain_27" ?bestFit t )
```


specLimits

```
specLimits(t_monteExprName ( {?sigma x_sigma | ?upper x_upper ?lower x_lower}  
  ) )  
=> t/nil
```

Description

Sets specification limits for yield analysis and histograms.

You can set specification limits for each of your measured values and then analyze how many runs are outside those limits (pass/fail) or you can analyze the spec sensitivity of measured quantities to changing input parameters.

You can specify limits using *x_upper* and *x_lower* options, or you can use the *x_sigma* option to have limits calculated for you based on a specified number of standard deviations of the actual data.

Note: You can specify *x_sigma* or you can specify *x_upper* and *x_lower*, but you cannot specify both.

Arguments

<i>t_monteExprName</i>	The <i>monteExpr</i> name with the appended swept parameter.
<i>x_sigma</i>	Identifies data lying outside an established sigma point from the mean. For instance, you might identify data lying outside 3 standard deviations (sigma) from the mean.
<i>x_upper</i>	Identifies data that is greater than this value.
<i>x_lower</i>	Identifies data that is less than this value.

Value Returned

<i>t</i>	Returned if successful.
<i>nil</i>	Returned otherwise.

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Example

```
specLimits("bw_27" ?upper 15E+06 ?lower 5+06)
```

yield

```
yield( s_type ?exprList l_monteExprName ?suppress x_suppress )  
=> t/nil
```

Description

Prints simple, conditional, or multiconditional yield statistics for the Monte Carlo data set.

Arguments

s_type

The type of statistics to print.

Valid values: 'simple, 'conditional, or
'multiconditional

'simple prints the yields for each measurement. Based upon the specification limits you set, the 'simple option prints the percentage of *pass* runs compared to the total number of Monte Carlo runs. For example, you set your specification limits for bandwidth, run 100 runs, and find that 60 of the runs pass the specification limits. For this example, the `yield` command calculates and displays a yield of 60% for bandwidth. The command also displays the total yield number, which is used when you have multiple measurements, each with its own limits. Total yield is the total percentage of *pass* runs where every parameter is within its specification limits for a Monte Carlo run.

'conditional prints conditional yields. To use a conditional yield, you specify a single measurement against which all other measurements are compared. The 'conditional option first sorts all of the Monte Carlo runs and picks out only the runs where the specified measurement passes. These passing runs are the starting point for the conditional yield calculation. So, in the bandwidth example above, instead of using 100 runs, the tool uses 60 runs as the base. Next, all of the other measurements are analyzed. For example, you have a second measurement called `maximum_25`. Out of the base 60 runs, `maximum_25` passes 30 times. It has a conditional yield of 50%. In addition to the conditional yield, the tool prints the total yield (based on all Monte Carlo runs) and the difference between the conditional and total yield numbers.

'multiconditional prints multiconditional yields. As in

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calculating the conditional yield, multiconditional yields are calculated from a base set of passing runs. However, instead of using one parameter to build the base set, for multiconditional yields you use two. Only runs where both measurements pass become part of the base set. All other measurements are then compared against that base.

l_monteExprName

The `monteExpr` name with the appended swept parameter.

x_suppress

If *s_type* is 'simple, suppresses the printing for yields greater than this percentage of the value.

Default value: 98

If *s_type* is 'conditional or 'multiconditional, suppresses the printing for delta yields less than this percentage of the value.

Default value: 98

Value Returned

t Returned if successful.

nil Returned otherwise.

Example

```
yield('simple ?exprList '("bw_27" "slew_27") ?suppress 70)
yield('conditional ?exprList '("max_27" "slew_27") )
yield('multiconditional ?exprList '("max_27" "slew_27") )
```

Optimization Commands

The commands for running optimization in the OCEAN environment are as follows.

optimizeAlgoControl

```
optimizeAlgoControl( ?relDelta x_relDelta ?relFunTol x_relFunTol ?relVarTol  
                    x_relVarTol )  
=> undefined/nil
```

Description

Changes the internal algorithm controls.

Arguments

<i>x_relDelta</i>	Finite difference relative perturbation. Default value: .005
<i>x_relFunTol</i>	Relative function convergence tolerance. Default value: .0001
<i>x_relVarTol</i>	Relative variable convergence tolerance. Default value: .0001

Value Returned

<i>undefined</i>	The return value for this function is not defined.
<i>nil</i>	Returns <i>nil</i> and an error message if there was a problem.

Example

```
optimizeAlgoControl(?relDelta .05)
```

optimizeGoal

```
optimizeGoal( t_name t_expr s_direction x_target x_acceptable [?percent  
             b_percent] )  
=> undefined/nil
```

Description

Sets up the goals for optimization.

Arguments

<i>t_name</i>	Name of the goal.
<i>t_expr</i>	Expression defining the goal.
<i>s_direction</i>	Valid values: 'max, 'min, 'match, 'le or 'ge Default value: 'match
<i>x_target</i>	The value to be matched or the lower or upper bound (depending on <i>s_direction</i>).
<i>x_acceptable</i>	Number or a waveform specifying the acceptable value. When a waveform is entered, each target point has its own acceptable value. Both <i>x_target</i> and <i>x_acceptable</i> must be expressions. The expression returns a number or a waveform.
<i>b_percent</i>	Specifies whether the <i>x_acceptable</i> field is a percentage of the target. When this is specified, <i>x_acceptable</i> is ignored.

Value Returned

<i>undefined</i>	The return value for this command is not defined.
<i>nil</i>	Returns <i>nil</i> and an error message if there was a problem.

Example

```
optimizeGoal( "bandwidth" 'bandwidth(v("/out") 3 "low") 'le 18M 15M )
```

optimizePlotOption

```
optimizePlotOption(  
    ?auto b_auto ?varHist b_varHist ?scalHist b_scalHist  
    ?funcObjHist b_funcObjHist ?numIter x_numIter ?fontSize x_fontSize  
    ?width x_width ?height x_height ?xloc xloc ?yloc yloc )  
=> undefined/nil
```

Description

Sets the plot options used to view the optimization iterations.

Arguments

<i>b_auto</i>	If set to <code>t</code> , auto plots after each iteration. Default value: <code>t</code>
<i>b_varHist</i>	If set to <code>t</code> , displays the history of the variables. Default value: <code>t</code>
<i>b_scalHist</i>	If set to <code>t</code> , displays the history of the scalars. Default value: <code>t</code>
<i>b_funcObjHist</i>	If set to <code>t</code> , displays the history of the functional objectives. Default value: <code>t</code>
<i>x_numIter</i>	Number of waveforms to display. There is one waveform stored available per functional iteration. Default value: 5
<i>x_fontSize</i>	Font size used in the Waveform window. Default value: 9
<i>x_width</i>	Width of the Waveform window. Default value: 630
<i>x_height</i>	Height of the Waveform window. Default value: 376
<i>xloc</i>	Specifies the top boundary of the optimize window when it is opened in the windowing environment. Default value: 511

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yloc Specifies the left boundary of the optimize window when it is opened in the windowing environment.
Default value: 377

Value Returned

undefined The return value is for this value is not defined.

nil Returns *nil* and an error message if there was a problem setting plot options.

Example

```
optimizePlotOption(?delta .05)
```

optimizeRun

```
optimizeRun(?goals l_goalNames ?vars l_varNames ?numIter x_numIter  
            ?algo s_algoName ?continue b_continue)  
=> t/nil
```

Description

Runs the optimizer using the goals specified with the `optimizeGoal` command.

Arguments

<i>l_goalNames</i>	Names of the goals to be used with this run of the optimizer. If none are specified, all declared goals are used.
<i>l_varNames</i>	Names of the variables to be used with this run of the optimizer. If none are specified, all declared variables are used.
<i>x_numIter</i>	Number of iterations that you want the optimizer to perform.
<i>s_algoName</i>	Algorithm that you want to use. Valid values: 'lsq, 'cfsqp, 'auto
<i>b_continue</i>	t indicates that this <code>optimizeRun</code> needs to continue from the previous <code>optimizeRun</code> (using the last design variables calculated from the last <code>optimizeRun</code>).

Value Returned

t	If the command was successful.
nil	Returns nil and an error message if there was a problem.

Example

```
optimizeRun()  
  optimizeRun(?goals '("bandwidth" "slewrates" )  
              ?vars '("rs" "vs") ?numIter 5)  
  optimizeRun( ?numIter 5 ?continue t )
```

Continues the previous `optimizeRun` for another 5 iterations.

optimizeVar

```
optimizeVar( t_name x_initVal x_minVal x_maxVal )  
=> undefined/nil
```

Description

Specifies the design variables to be used with optimization.

Arguments

<i>t_name</i>	Name of the design variable.
<i>x_initVal</i>	Initial value of the variable.
<i>x_minVal</i>	Lower bound of the variable.
<i>x_maxVal</i>	Upper bound of the variable.

Value Returned

<i>undefined</i>	The return value for this function is not defined.
<i>nil</i>	Returns <i>nil</i> and an error message if there was a problem.

Example

```
optimizeVar( "res" 100 ?minVal 1 ?maxVal 1000 )
```

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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](#) on page 422
- [digitalHostMode](#) on page 423
- [digitalHostName](#) on page 424
- [hostMode](#) on page 425
- [hostName](#) on page 426
- [killJob](#) on page 427
- [monitor](#) on page 428
- [remoteDir](#) on page 429
- [resumeJob](#) on page 430
- [suspendJob](#) on page 431
- [wait](#) on page 432

This chapter also provides sample OCEAN scripts that optimally use these commands. See the section [Sample Scripts](#) on page 433.

For detailed information on distributed processing, refer to [Virtuoso® Analog Distributed Processing Option User Guide](#).

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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_jobname2...t_jobnameN
 Additional jobs that you want to delete.

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.

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

<code>'local</code>	Sets the simulation to run locally on the user's machine.
<code>'remote</code>	Sets the simulation to run on a remote host. If you use this argument, you must specify the host name by using the <code>digitalHostName</code> command.

Value Returned

<code>t</code>	Returns <code>t</code> if successful.
<code>nil</code>	Returns <code>nil</code> and prints an error message if unsuccessful.

Example

```
digitalHostMode( 'local )
```

Sets the digital simulator to run locally on the user's host.

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`.

hostMode

```
hostMode( { 'local | 'remote | 'distributed } )  
=> t/nil
```

Description

Sets the simulation host mode.

The default value for `hostMode` is specified in the `asimenv.startup` file with the `hostMode` environment variable.

Arguments

<code>'local</code>	Sets the simulation to run locally on the user's machine.
<code>'remote</code>	Sets the simulation to run on a remote host queue. For this release, the remote host is specified in the <code>.cdsenv</code> file.
<code>'distributed</code>	Sets the simulation to run using the distributed processing software.

Value Returned

<code>t</code>	Returns <code>t</code> if successful.
<code>nil</code>	Returns <code>nil</code> and prints an error message if unsuccessful.

Example

```
hostMode( 'distributed )  
=> t
```

Enables distributed processing on the current host.

hostName

```
hostName( t_name )  
=> t/nil
```

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

<i>t_name</i>	Name used to identify the remote host.
---------------	--

Value Returned

<i>t</i>	Returns <i>t</i> if successful.
----------	---------------------------------

<i>nil</i>	Returns <i>nil</i> and prints an error message if unsuccessful.
------------	---

Example

```
hostName( "remotehost" )
```

Specifies that the host called `remotehost` is to be used for remote simulation.

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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.

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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.

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

<i>t_path</i>	Specifies the path to the project directory on the remote host to be used for remote simulation.
---------------	--

Value Returned

<i>t</i>	Returns <i>t</i> if successful.
<i>nil</i>	Returns <i>nil</i> and prints an error message if unsuccessful.

Example

```
remoteDir( "~/simulation" )
```

Specifies that the project directory is `~/simulation`.

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resumeJob

```
resumeJob( t_jobName [t_jobName2 t_jobName3 ... t_jobNameN] )  
=> t/nil
```

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.

t_jobName2...*t_jobNameN*
 Additional jobs that you want to resume

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.

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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`.

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OCEAN Distributed Processing Commands

wait

```
wait( 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 very 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 [Virtuoso® Analog Distributed Processing Option User Guide](#).

Arguments

<i>t_jobName</i>	Name used to identify the job. The job name is user defined or system generated, depending on how the user submitted the job.
<i>t_jobName2...t_jobnameN</i>	Additional jobs that you want to postpone.

Value Returned

<i>t</i>	Returns <i>t</i> if successful.
<i>nil</i>	Returns <i>nil</i> and prints an error message if unsuccessful.

Example

```
wait( 'mycktl' )  
=> t
```

Postpones execution of all subsequent OCEAN commands until the job `mycktl` completes.

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 set up and run a Corners analysis in blocking mode and select results
- To set up and run a montecarlo 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 `dependentOn` 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 = appendl( jobList run( ?queue "test" ?host "menaka" ) )

analysis('tran ?stop "50u" )

; starting second job
jobList = appendl( jobList run(?jobName "job_2" ?queue "test" ?host "menaka" ))

analysis('tran ?stop "10u" )

; starting third job, which is dependent on job_2
```

OCEAN Reference

OCEAN Distributed Processing Commands

```
jobList= appendl(jobList run(?jobName "disable" ?queue "test" ?dependentOn
                    symbolToString(car(last(jobList))))))

; wait for all the jobs to complete
wait((appendl 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"))
```

OCEAN Reference

OCEAN Distributed Processing Commands

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(
"/home/amit/Artist446/simulation/ampTest/spectre/schematic/netlist/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 set up and run a Corners analysis in blocking mode and select results

```
; set up the environment for corners analysis
simulator('spectre)
design("./netlist/netlist")
hostMode( 'distributed )
analysis('tran ?stop 50n)
keep('allv)
definitionFile("model")
loadPcf("./singleNumeric.pcf")
loadDcf("./singleNumeric.dcf")

; submit the job in blocking mode, to the queue test and machine menaka
cornerRun( ?block t ?queue "fast" ?host "menaka" )

; select and print/plot the results
selectResults('tran)
plot v("2")
```

OCEAN Reference

OCEAN Distributed Processing Commands

```
ocnPrint v("2")
```

To set up and run a montecarlo analysis in blocking mode and select results

```
; set up the environment for montecarlo analysis
simulator( 'spectre )
hostMode( "distributed" )
design("./spectre/netlist/netlist")
resultsDir( "./spectre" )
path("./spectre/netlist" )
modelFile( '( "spectreLib.scs" "statistics" ) )
definitionFile( "update" "init" "lowpassStats" )
analysis('ac ?start "1" ?stop "100M" )
desVar( "rout2" 3K )
desVar( "rout1" 1K )
desVar( "rin2" 5K )
desVar( "rin1" 1K )
desVar( "cloop" .001u )
desVar( "cin" .017u )
temp( 27 )
monteCarlo( ?numIters "100" ?startIter "1"
  ?analysisVariation "Process Only" ?sweptParam "None"
  ?sweptParamVals "27" ?saveData t
  ?nomRun nil ?append nil)
monteExpr( "bw" "bandwidth(VF('OUT') 3 'low')" )
monteExpr( "phase" "value(phase(VF('OUT'))) 100000)" )
monteExpr( "db20" "value(dB20(VF('OUT'))) 100000)" )

; submit the job in blocking mode, to the queue test and machine menaka
monteRun( ?block t ?queue "fast" ?host "menaka" )

; Initializes the Monte Carlo data analysis tools
monteResults()
```

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")
```

OCEAN Reference

OCEAN Distributed Processing Commands

```
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 )
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" )
```

OCEAN Reference

OCEAN Distributed Processing Commands

```
analysis('tran ?stop "3u" )
desVar( "CAP" 0.8p )
temp( 27 )
jobList = appendl( jobList run(?queue "SUN5_5032" ?host "menaka"))

; wait for both the jobs to finish
wait( (appendl 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 )
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` on page 449

`if` on page 440

`cond` on page 451

`unless` on page 442

`for` on page 444

`when` on page 443

`foreach` on page 446

`while` on page 448

if

```
if( g_condition g_thenExpression [g_elseExpression] )  
    => g_result
```

Description

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.

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.

Examples

```
x = 2  
if( x > 5 1 0 )  
=> 0
```

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"
```


OCEAN Reference Language Constructs

Returns "non-nil" because x was not nil. If x was nil, "nil" would be returned.

```
x = 7  
if( x > 5 1 0 )  
=> 1
```

Returns 1 because x is greater than 5.

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`.

Examples

```
x = -123  
unless( x >= 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`.

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

<i>g_condition</i>	Any SKILL expression.
<i>g_expr1</i> ...	Any SKILL expression.

Value Returned

<i>g_result</i>	Returns the value of the last expression of the sequence <i>g_expr1</i> ... if <i>g_condition</i> evaluates to non-nil.
<code>nil</code>	returns <code>nil</code> if the <i>g_condition</i> expression evaluates to <code>nil</code> .

Examples

```
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`.

for

```
for( s_loopVar x_initialValue x_finalValue g_expr1 [g_expr2 ...] )  
    => t
```

Description

Evaluates the sequence *g_expr1 g_expr2 ...* 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_expr1 g_expr2 ...* 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.

Arguments

<i>s_loopVar</i>	Name of the local loop variable that must not be changed inside the loop.
<i>x_initialValue</i>	Integer expression setting the initial value for the local loop variable.
<i>x_finalValue</i>	Integer expression giving final limit value for the loop.
<i>g_expr1</i>	Expression to evaluate inside loop.
<i>g_expr2</i> ...	Additional expressions to evaluate inside loop.

Value Returned

t This construct always returns *t*.

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Examples

```
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

```
foreach( s_formalVar g_exprList g_expr1 [g_expr2 ...] )  
    => l_valueList  
  
foreach( (s_formalVar1...s_formalVarN) g_exprList1... g_exprListN g_expr1  
    [g_expr2 ...] )  
    => l_valueList  
  
foreach( s_formalVar g_exprTable g_expr1 [g_expr2 ...])  
    => o_valueTable
```

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,

```
foreach( (s_formalVar1...s_formalVarN) g_exprList1... g_exprListN g_expr1  
[g_expr2 ...] )=> l_valueList
```

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
```

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

<i>s_formalVar</i>	Name of the variable.
<i>g_exprList</i>	Expression whose value is a list of elements to assign to the formal variable <i>s_formalVar</i> .
<i>g_expr1 g_expr2 ...</i>	Expressions to execute.
<i>g_exprTable</i>	Association table whose elements are to be processed.

Value Returned

<i>l_valueList</i>	Returns the value of the second argument, <i>g_exprList</i> .
<i>o_valueTable</i>	Returns the value of <i>g_exprTable</i> .

Examples

```
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.

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.

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*). Note that because 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_expr1 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.

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 *l_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 *l_clause1* *l_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.

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.

l_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.

l_clause2...

Zero or more clauses of the same form as *l_clause1*.

Value Returned

g_result

Returns the value of the last expression evaluated in the matched clause.

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`nil` Returns `nil` if there is no match.

Example

```
cornersType = "min"
type = case( cornersType
  ("min" path("./min"))
  ("typ" path("./typ"))
  ("max" path("./max"))
  (t println("you have not chosen an appropriate
    corner")))
=> path is set to "./min"
```

Sets `path` to `./min`.

cond

```
cond( l_clause1 ... )  
    => g_result/nil
```

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

l_clause1 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 Returns the value of the last expression of the satisfied clause.

nil Returns nil if no clause is satisfied.

Example

```
procedure( test(x)  
    cond(((null x) (println "Arg is null"))  
        ((numberp x) (println "Arg is a number"))  
        ((stringp x) (println "Arg is a string"))  
        (t (println "Arg is an unknown type")))  
    )  
test( nil )  
=> nil; Prints "Arg is null".  
test( 5 )  
=> nil; Prints "Arg is a number".  
test( 'sym )  
=> nil; Prints "Arg is an unknown type".
```

OCEAN Reference Language Constructs

Tests each of the arguments according to the conditions specified with `cond`.

File Commands and Functions

This chapter contains information on the following commands:

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fscanf on page 455

gets on page 457

infile on page 458

load on page 459

newline on page 461

outfile on page 462

printf on page 464

println on page 465

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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.

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

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.

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File Commands and Functions

Value Returned

<code>x_items</code>	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.
<code>nil</code>	Returns <code>nil</code> 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")
```


OCEAN Reference

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

<i>s_variableName</i>	Variable in which to store the input string.
<i>p_inputPort</i>	Name of input port. Default value: <code>piport</code>

Value Returned

<i>t_string</i>	Returns the input string when successful.
<code>nil</code>	Returns <code>nil</code> when the end of file is reached. (<i>s_variableName</i> 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.

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_inport Returns the port opened for reading the named file.

`nil` Returns `nil` if the file does not exist or cannot be opened for reading.

Examples

```
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.

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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.

Arguments

<i>t_fileName</i>	File to be loaded. Uses the file name extension to determine the language mode to use. Valid values: <ul style="list-style-type: none"><code>.ils</code> Means the file contains SKILL++ code.<code>.il</code> Means the file contains SKILL code.<code>.ocn</code> Means the file contains OCEAN commands (with SKILL or SKILL++ commands)
<i>t_password</i>	Password, if <i>t_fileName</i> is an encrypted file.

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File Commands and Functions

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.

newline

```
newline( [p_outputPort] )  
=> nil
```

Description

Prints a newline (backslash `\n`) character and then flushes the output port.

Arguments

<i>p_outputPort</i>	Output port. Defaults value: <code>poport</code>
---------------------	---

Value Returned

<code>nil</code>	Prints a newline and then returns <code>nil</code> .
------------------	--

Example

```
print( "Hello" ) newline() print( "World!" )  
"Hello"  
"World!"  
=> nil
```

Prints a newline character after `Hello`.

OCEAN Reference

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

<i>S_fileName</i>	Name of the file to open or create.
<i>t_mode</i>	Mode in which to open the file. If <code>a</code> , the file is opened in append mode; if <code>w</code> , a new file is created for writing (any existing file is overwritten). Default value: <code>w</code>

Value Returned

<i>p_outport</i>	An output port ready to write to the specified file.
<code>nil</code>	returns <code>nil</code> if the named file cannot be opened for writing. An error is signaled if an illegal mode string is supplied.

Examples

```
p = outfile( "/tmp/out.il" "w" )  
=> port:"/tmp/out.il"
```

Opens the `/tmp/out.il` port.

```
outfile( "/bin/ls" )  
=> nil
```

OCEAN Reference

File Commands and Functions

Returns `nil`, indicating that the specified port could not be opened.

printf

```
printf( t_formatString [g_arg1 ...] )  
=> t
```

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 *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 => 197.9687  
printf( "The test measures %10.2f." x )
```

Prints the following line to *poport* and returns *t*.

```
The test measures 197.97. => t
```


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

<i>g_value</i>	Any SKILL value.
<i>p_outputPort</i>	Port to be used for output. Default value: <code>poport</code>

Value Returned

<code>nil</code>	Prints the given object and returns <code>nil</code> .
------------------	--

Example

```
for( i 1 3 println( "hello" ))  
"hello"  
"hello"  
"hello"  
=> t
```

Prints hello three times. `for` always returns `t`.

OCEAN Reference
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® 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 `path` command.

Mixed-Signal in OCEAN 4.4.6

All of the analog OCEAN features are available in mixed-signal. This means you can set up analyses, change options, change the path, and so forth.

There are limitations in the area of mixed-signal simulation.

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OCEAN 4.4.6 Issues

- If mixed-signal simulation is run using a standalone OCEAN tool, then the complete netlist must be created before running the simulation. The netlist can be created using Affirma Analog Design Environment or by specifying the design as lib-cell-view using the ocean command `design` in CIW of the workbench followed by the OCEAN commands `createNetlist` and `run`.

For example:

```
design("mylib" "ampTest" "schematic")

; design using lib-cell-view can only be specified in CIW of
workbench

createNetlist()

run()
```

- If mixed-signal simulation is run using OCEAN commands in the CIW of the workbench, then the design should be specified as lib-cell-view.

Otherwise, if the design is specified as the path to the netlist, for example as `design("./simulation/ampTest/specter/netlist")`, then the complete netlist should be created before running the simulation using the procedure specified above.

In the 4.4.6 release, there are no commands that operate on Verilog-XL final netlists. If you need to change anything in the final netlist, you must make those changes by hand.

However, you can change any of the command line arguments that are sent to the Verilog-XL simulator. This means you can change any of the digital options or any of the mixed-signal options. To see what these options are, choose *Simulation – Options – Digital* in the Virtuoso® Analog Design Environment window.

For example, you can change acceleration, keep nodes, and library files.

For detailed information, please refer to the [Virtuoso® Mixed-Signal Circuit Design Environment User Guide](#).

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