Contents

Preface ................................................................. 17
Licensing in OCEAN .................................................. 18
Related Documents for OCEAN ........................................ 18
  Installation, Environment, and Infrastructure .................. 18
  Virtuoso Tools ...................................................... 19
Typographic and Syntax Conventions ................................. 19
  SKILL Syntax Examples .............................................. 20
Identifiers Used to Denote Data Types .............................. 22

1
Introduction to OCEAN ................................................. 25
Types of OCEAN Commands .......................................... 26
OCEAN Online Help ................................................... 26
OCEAN Syntax Overview .............................................. 27
  Common SKILL Syntax Characters Used in OCEAN ............. 27
  Parentheses ......................................................... 27
  Quotation Marks .................................................... 28
  Single Quotation Marks .......................................... 29
  Question Mark ....................................................... 29
  Data Types Used in OCEAN ......................................... 30
  OCEAN Return Values .............................................. 31
  Design Variables in OCEAN ........................................ 31
  outputs() in OCEAN ............................................... 32
Parametric Analysis .................................................. 33
  Data Access Without Running a Simulation ...................... 34
Distributed Processing ............................................... 34
  Blocking and Nonblocking Modes ................................ 35
Waveform Tool ......................................................... 36
2
Using OCEAN ................................................................. 37
OCEAN Use Models ......................................................... 37
Using OCEAN Interactively .............................................. 38
    Using OCEAN from a UNIX Shell ................................ 38
    Using OCEAN from the CIW ...................................... 39
    Interactive Session Demonstrating the OCEAN Use Model  40
License Requirements ..................................................... 41
Creating OCEAN Scripts ................................................ 42
    Creating Scripts Using Sample Script Files .................. 42
    Creating Scripts from the Analog Design Environment ...... 42
    Selectively Creating Scripts .................................... 42
    Loading OCEAN Scripts ........................................... 45
Selecting Results .......................................................... 46
    Selecting Results Run from Worst Case Scripts for Cross-Probing or Back Annotating Operating Points ..................... 46
    Selecting Results Run from Spectre Stand Alone .......... 47
Running Multiple Simulators .......................................... 48
OCEAN Tips ................................................................. 48

3
Introduction to SKILL .................................................. 51
The Advantages of SKILL ............................................... 51
Naming Conventions ..................................................... 52
Arithmetic Operators ................................................... 52
Scaling Factors .......................................................... 52
Relational and Logical Operators .................................... 54
    Relational Operators ............................................. 54
    Logical Operators ................................................ 55
SKILL Syntax ............................................................ 56
    Special Characters ................................................ 56
    White Space ......................................................... 57
    Comments .......................................................... 57
    Role of Parentheses .............................................. 58
    Line Continuation ................................................ 59
4
Working with SKILL .............................................. 63
Skill Functions .................................................. 63
Data Types ....................................................... 63
   Numbers ....................................................... 64
   Atoms ......................................................... 65
   Constants and Variables .................................. 65
   Strings ....................................................... 65
Arrays ............................................................ 66
   Allocating an Array of a Given Size ..................... 66
Concatenating Strings (Lists) ............................... 66
Comparing Strings ............................................. 67
Declaring a SKILL Function ................................. 68
   Defining Function Parameters ................................ 69
   Defining Local Variables (let) ............................. 69
Skill Function Return Values ............................... 70
Syntax Functions for Defining Functions .................. 70
   procedure .................................................... 70
   Terms and Definitions ..................................... 70

5
OCEAN Environment Commands ............................... 73
   appendPath .................................................. 74
   path .......................................................... 75
   prependPath ................................................ 76
   setup ........................................................ 77
   history ....................................................... 79
   ocnSetSilentMode ........................................... 80
# 6 Simulation Commands

<table>
<thead>
<tr>
<th>Command</th>
<th>Page</th>
</tr>
</thead>
<tbody>
<tr>
<td>ac</td>
<td>83</td>
</tr>
<tr>
<td>analysis</td>
<td>85</td>
</tr>
<tr>
<td>converge</td>
<td>88</td>
</tr>
<tr>
<td>connectRules</td>
<td>89</td>
</tr>
<tr>
<td>createFinalNetlist</td>
<td>93</td>
</tr>
<tr>
<td>createNetlist</td>
<td>94</td>
</tr>
<tr>
<td>dc</td>
<td>96</td>
</tr>
<tr>
<td>definitionFile</td>
<td>98</td>
</tr>
<tr>
<td>delete</td>
<td>99</td>
</tr>
<tr>
<td>design</td>
<td>101</td>
</tr>
<tr>
<td>desVar</td>
<td>103</td>
</tr>
<tr>
<td>discipline</td>
<td>105</td>
</tr>
<tr>
<td>displayNetlist</td>
<td>107</td>
</tr>
<tr>
<td>envOption</td>
<td>108</td>
</tr>
<tr>
<td>evcdFile</td>
<td>110</td>
</tr>
<tr>
<td>evcdInfoFile</td>
<td>111</td>
</tr>
<tr>
<td>forcenode</td>
<td>112</td>
</tr>
<tr>
<td>globalSigAlias</td>
<td>113</td>
</tr>
<tr>
<td>globalSignal</td>
<td>114</td>
</tr>
<tr>
<td>ic</td>
<td>116</td>
</tr>
<tr>
<td>includeFile</td>
<td>117</td>
</tr>
<tr>
<td>modelFile</td>
<td>118</td>
</tr>
<tr>
<td>nodeset</td>
<td>119</td>
</tr>
<tr>
<td>noise</td>
<td>120</td>
</tr>
<tr>
<td>ocnCloseSession</td>
<td>121</td>
</tr>
<tr>
<td>ocnDisplay</td>
<td>122</td>
</tr>
<tr>
<td>ocnGetAdjustedPath</td>
<td>124</td>
</tr>
<tr>
<td>ocnWaveformTool</td>
<td>125</td>
</tr>
<tr>
<td>off</td>
<td>126</td>
</tr>
<tr>
<td>option</td>
<td>127</td>
</tr>
<tr>
<td>restore</td>
<td>129</td>
</tr>
<tr>
<td>resultsDir</td>
<td>130</td>
</tr>
<tr>
<td>run</td>
<td>131</td>
</tr>
</tbody>
</table>
# Data Access Commands

<table>
<thead>
<tr>
<th>Command</th>
<th>Page</th>
</tr>
</thead>
<tbody>
<tr>
<td>save</td>
<td>135</td>
</tr>
<tr>
<td>saveOption</td>
<td>137</td>
</tr>
<tr>
<td>simulator</td>
<td>139</td>
</tr>
<tr>
<td>solver</td>
<td>140</td>
</tr>
<tr>
<td>stimulusFile</td>
<td>141</td>
</tr>
<tr>
<td>store</td>
<td>143</td>
</tr>
<tr>
<td>temp</td>
<td>144</td>
</tr>
<tr>
<td>tran</td>
<td>145</td>
</tr>
<tr>
<td>vcdFile</td>
<td>146</td>
</tr>
<tr>
<td>vcdInfoFile</td>
<td>147</td>
</tr>
<tr>
<td>vecFile</td>
<td>148</td>
</tr>
<tr>
<td>hlcheck</td>
<td>149</td>
</tr>
<tr>
<td>dataTypes</td>
<td>151</td>
</tr>
<tr>
<td>getData</td>
<td>153</td>
</tr>
<tr>
<td>getResult</td>
<td>154</td>
</tr>
<tr>
<td>i</td>
<td>156</td>
</tr>
<tr>
<td>ocnHelp</td>
<td>157</td>
</tr>
<tr>
<td>ocnResetResults</td>
<td>159</td>
</tr>
<tr>
<td>openResults</td>
<td>161</td>
</tr>
<tr>
<td>outputParams</td>
<td>162</td>
</tr>
<tr>
<td>outputs</td>
<td>164</td>
</tr>
<tr>
<td>phaseNoise</td>
<td>166</td>
</tr>
<tr>
<td>pv</td>
<td>168</td>
</tr>
<tr>
<td>resultParam</td>
<td>170</td>
</tr>
<tr>
<td>results</td>
<td>172</td>
</tr>
<tr>
<td>selectResult</td>
<td>174</td>
</tr>
<tr>
<td>sp</td>
<td>174</td>
</tr>
<tr>
<td>sweepNames</td>
<td>177</td>
</tr>
<tr>
<td>sweepValues</td>
<td>179</td>
</tr>
<tr>
<td>sweepVarValues</td>
<td>181</td>
</tr>
<tr>
<td>v</td>
<td>182</td>
</tr>
<tr>
<td>vswr</td>
<td>184</td>
</tr>
</tbody>
</table>
8
Plotting and Printing Commands ............................................. 191
   addSubwindow .................................................................. 193
   addSubwindowTitle .......................................................... 194
   addTitle ............................................................................. 195
   addWaveLabel .................................................................... 196
   addWindowLabel ............................................................... 198
   clearAll .............................................................................. 199
   clearSubwindow .................................................................. 200
   currentSubwindow ............................................................ 201
   currentWindow .................................................................... 202
   dbCompressionPlot ............................................................. 203
   dcmatchSummary .................................................................. 204
   deleteSubwindow .................................................................. 208
   deleteWaveform .................................................................... 209
   displayMode ....................................................................... 210
   getAsciiWave ...................................................................... 211
   graphicsOff ........................................................................ 212
   graphicsOn ......................................................................... 213
   hardCopy ............................................................................ 214
   hardCopyOptions ............................................................... 215
   ip3Plot .............................................................................. 217
   newWindow ......................................................................... 218
   noiseSummary ..................................................................... 219
   ocnPrint ............................................................................ 223
   ocnSetAttrib ..................................................................... 226
   ocnYvsYplot ....................................................................... 228
   plot .................................................................................... 230
   plotStyle .......................................................................... 233
   pzPlot .............................................................................. 234
   pzSummary ........................................................................ 236
   removeLabel ...................................................................... 238
### OCEAN Reference

<table>
<thead>
<tr>
<th>Section</th>
<th>Page</th>
</tr>
</thead>
<tbody>
<tr>
<td>Report</td>
<td>239</td>
</tr>
<tr>
<td>xLimit</td>
<td>242</td>
</tr>
<tr>
<td>yLimit</td>
<td>243</td>
</tr>
<tr>
<td>Plotting and Printing SpectreRF Functions in OCEAN</td>
<td>245</td>
</tr>
<tr>
<td><strong>9</strong> OCEAN Aliases</td>
<td>247</td>
</tr>
<tr>
<td><strong>10</strong> Predefined and Waveform (Calculator) Functions</td>
<td>249</td>
</tr>
<tr>
<td>Predefined Arithmetic Functions</td>
<td>254</td>
</tr>
<tr>
<td>abs</td>
<td>256</td>
</tr>
<tr>
<td>acos</td>
<td>257</td>
</tr>
<tr>
<td>add1</td>
<td>258</td>
</tr>
<tr>
<td>asin</td>
<td>259</td>
</tr>
<tr>
<td>atan</td>
<td>260</td>
</tr>
<tr>
<td>cos</td>
<td>261</td>
</tr>
<tr>
<td>exp</td>
<td>262</td>
</tr>
<tr>
<td>int</td>
<td>263</td>
</tr>
<tr>
<td>linRg</td>
<td>264</td>
</tr>
<tr>
<td>log</td>
<td>265</td>
</tr>
<tr>
<td>logRg</td>
<td>266</td>
</tr>
<tr>
<td>max</td>
<td>267</td>
</tr>
<tr>
<td>min</td>
<td>268</td>
</tr>
<tr>
<td>mod</td>
<td>269</td>
</tr>
<tr>
<td>random</td>
<td>270</td>
</tr>
<tr>
<td>round</td>
<td>271</td>
</tr>
<tr>
<td>sin</td>
<td>272</td>
</tr>
<tr>
<td>sqrt</td>
<td>273</td>
</tr>
<tr>
<td>srandom</td>
<td>274</td>
</tr>
<tr>
<td>sub1</td>
<td>275</td>
</tr>
<tr>
<td>tan</td>
<td>276</td>
</tr>
<tr>
<td>xor</td>
<td>277</td>
</tr>
<tr>
<td>Waveform (Calculator) Functions</td>
<td>278</td>
</tr>
<tr>
<td>average</td>
<td>279</td>
</tr>
</tbody>
</table>
awvPlaceXMarker .................................................................................. 281
awvPlaceYMarker .................................................................................. 282
awvRefreshOutputPlotWindows ................................................................. 283
b1f .......................................................................................................... 284
bandwidth ............................................................................................... 285
click .......................................................................................................... 286
compare .................................................................................................... 288
compression .............................................................................................. 290
compressionVRI ......................................................................................... 292
compressionVRIcurves ........................................................................... 294
complex ...................................................................................................... 296
complexp .................................................................................................... 297
conjugate ................................................................................................. 298
convolve .................................................................................................... 299
cPwrContour ............................................................................................. 301
cReflContour ............................................................................................. 303
cross .......................................................................................................... 305
db10 .......................................................................................................... 307
db20 .......................................................................................................... 308
dbm ........................................................................................................... 309
delay .......................................................................................................... 310
deriv .......................................................................................................... 314
dft .............................................................................................................. 315
dftbb .......................................................................................................... 317
dnl .............................................................................................................. 319
dutyCycle ................................................................................................. 321
evmQAM .................................................................................................... 323
evmQpsk .................................................................................................... 325
eyeDiagram ............................................................................................... 327
flip ............................................................................................................. 329
fourEval ..................................................................................................... 330
deq ............................................................................................................. 332
deq_jitter ................................................................................................. 334
dfrequency ............................................................................................... 336
gaw ............................................................................................................ 337
gac ............................................................................................................ 338
<table>
<thead>
<tr>
<th>Function</th>
<th>Page</th>
</tr>
</thead>
<tbody>
<tr>
<td>gainBwProd</td>
<td>340</td>
</tr>
<tr>
<td>gainMargin</td>
<td>342</td>
</tr>
<tr>
<td>gmax</td>
<td>343</td>
</tr>
<tr>
<td>gmin</td>
<td>344</td>
</tr>
<tr>
<td>gmsg</td>
<td>345</td>
</tr>
<tr>
<td>gmux</td>
<td>346</td>
</tr>
<tr>
<td>gp</td>
<td>347</td>
</tr>
<tr>
<td>gpc</td>
<td>348</td>
</tr>
<tr>
<td>groupDelay</td>
<td>350</td>
</tr>
<tr>
<td>gt</td>
<td>351</td>
</tr>
<tr>
<td>harmonic</td>
<td>352</td>
</tr>
<tr>
<td>harmonicFreqList</td>
<td>354</td>
</tr>
<tr>
<td>harmonicList</td>
<td>356</td>
</tr>
<tr>
<td>histo</td>
<td>358</td>
</tr>
<tr>
<td>iinteg</td>
<td>359</td>
</tr>
<tr>
<td>imag</td>
<td>360</td>
</tr>
<tr>
<td>integ</td>
<td>361</td>
</tr>
<tr>
<td>intersect</td>
<td>363</td>
</tr>
<tr>
<td>ipn</td>
<td>364</td>
</tr>
<tr>
<td>ipnVRI</td>
<td>367</td>
</tr>
<tr>
<td>ipnVRICurves</td>
<td>370</td>
</tr>
<tr>
<td>kf</td>
<td>373</td>
</tr>
<tr>
<td>ln</td>
<td>374</td>
</tr>
<tr>
<td>log10</td>
<td>375</td>
</tr>
<tr>
<td>lsb</td>
<td>376</td>
</tr>
<tr>
<td>lshift</td>
<td>377</td>
</tr>
<tr>
<td>mag</td>
<td>378</td>
</tr>
<tr>
<td>nc</td>
<td>379</td>
</tr>
<tr>
<td>overshoot</td>
<td>381</td>
</tr>
<tr>
<td>peak</td>
<td>384</td>
</tr>
<tr>
<td>peakToPeak</td>
<td>386</td>
</tr>
<tr>
<td>period_jitter</td>
<td>387</td>
</tr>
<tr>
<td>phase</td>
<td>389</td>
</tr>
<tr>
<td>phaseDeg</td>
<td>390</td>
</tr>
<tr>
<td>phaseDegUnwrapped</td>
<td>391</td>
</tr>
<tr>
<td>phaseMargin</td>
<td>392</td>
</tr>
</tbody>
</table>
11

Parametric Analysis Commands

paramAnalysis ........................................ 448
paramRun .............................................. 452
## 12
### OCEAN Distributed Processing Commands

- deleteJob .............................................. 456
- digitalHostMode ...................................... 457
- digitalHostName ..................................... 458
- hostMode .............................................. 459
- hostName ............................................. 460
- killJob ............................................... 461
- monitor ............................................... 462
- remoteDir ............................................ 463
- resumeJob ............................................ 464
- suspendJob .......................................... 465
- wait .................................................. 466
- Sample Scripts ...................................... 467

## 13
### Language Constructs

- if ..................................................... 474
- unless ............................................... 476
- when .................................................. 477
- for ..................................................... 478
- foreach .............................................. 480
- while ............................................... 482
- case ................................................... 483
- cond .................................................. 485

## 14
### File Commands and Functions

- close .................................................. 488
- fscanf ................................................. 489
- gets ................................................... 491
- infile ............................................... 492
- load ................................................... 493
- newline ............................................... 495
15
OCEAN Commands in XL Mode

ocnSetXLMode
ocnxlBeginTest
ocnxlEndTest
ocnxlEndXLMode
ocnxlSelectTest
ocnxlSweepVar
ocnxlSweepParam
ocnxlCorner
ocnxlDisableTest
ocnxlDisableSweepVar
ocnxlDisableSweepParam
ocnxlDisableCornerForTest
ocnxlGlobalOptimizationOptions
ocnxlJobSetup
ocnxlLocalOptimizationOptions
ocnxlModelGroup
ocnxlOutputOceanScript
ocnxlOutputMatlabScript
ocnxlMonteCarloOptions
ocnxlPutToleranceSpec
ocnxlPutMinSpec
ocnxlPutMaxSpec
ocnxlPutGreaterthanSpec
ocnxlPutLessthanSpec
ocnxlPutRangeSpec
ocnxlPutTargetSpec
ocnxlResultsLocation
ocnxlRun
<table>
<thead>
<tr>
<th>Function</th>
<th>Page</th>
</tr>
</thead>
<tbody>
<tr>
<td>ocnxlRunSetupSummary</td>
<td>534</td>
</tr>
<tr>
<td>ocnxlSamplingOptions</td>
<td>535</td>
</tr>
<tr>
<td>ocnxlSetupLocation</td>
<td>536</td>
</tr>
<tr>
<td>ocnxlOutputExpr</td>
<td>537</td>
</tr>
<tr>
<td>ocnxlOutputSignal</td>
<td>538</td>
</tr>
<tr>
<td>ocnxlOutputTerminal</td>
<td>539</td>
</tr>
<tr>
<td>ocnxlOutputSummary</td>
<td>540</td>
</tr>
<tr>
<td>ocnxlTargetCellView</td>
<td>542</td>
</tr>
<tr>
<td>ocnxlYieldImprovementOptions</td>
<td>543</td>
</tr>
<tr>
<td>ocnxlEnableCornerForTest</td>
<td>546</td>
</tr>
<tr>
<td>ocnxlEnableSweepParam</td>
<td>547</td>
</tr>
<tr>
<td>ocnxlEnableSweepVar</td>
<td>548</td>
</tr>
<tr>
<td>ocnxlEnableTest</td>
<td>549</td>
</tr>
<tr>
<td>ocnxlGetBestPointParams</td>
<td>550</td>
</tr>
<tr>
<td>ocnxlGetCorners</td>
<td>551</td>
</tr>
<tr>
<td>ocnxlGetCurrentHistory</td>
<td>552</td>
</tr>
<tr>
<td>ocnxlGetCurrentHistoryId</td>
<td>553</td>
</tr>
<tr>
<td>ocnxlGetSession</td>
<td>554</td>
</tr>
<tr>
<td>ocnxlGetSpecs</td>
<td>555</td>
</tr>
<tr>
<td>ocnxlGetTests</td>
<td>556</td>
</tr>
<tr>
<td>ocnxlRemoveSpec</td>
<td>557</td>
</tr>
<tr>
<td>ocnxlRenameCurrentHistory</td>
<td>558</td>
</tr>
<tr>
<td>ocnxlRun</td>
<td>559</td>
</tr>
<tr>
<td>ocnxlHistoryPrefix</td>
<td>561</td>
</tr>
<tr>
<td>ocnxlLoadSetupState</td>
<td>562</td>
</tr>
<tr>
<td>ocnxlStartingPoint</td>
<td>565</td>
</tr>
<tr>
<td>ocnxlOutputAreaGoal</td>
<td>566</td>
</tr>
<tr>
<td>ocnxlConjugateGradientOptions</td>
<td>567</td>
</tr>
<tr>
<td>ocnxlMTSEnable</td>
<td>568</td>
</tr>
<tr>
<td>ocnxlMTSBlock</td>
<td>569</td>
</tr>
<tr>
<td>ocnxlProjectDir</td>
<td>571</td>
</tr>
<tr>
<td>ocnxlSimResultsLocation</td>
<td>572</td>
</tr>
<tr>
<td>ocnxlDisableCorner</td>
<td>572</td>
</tr>
<tr>
<td>ocnxlEnableCorner</td>
<td>574</td>
</tr>
<tr>
<td>ocnxlSaveSetupAs</td>
<td>575</td>
</tr>
<tr>
<td>ocnxlSetAllParametersDisabled</td>
<td>576</td>
</tr>
</tbody>
</table>
ocnxISetAllVarsDisabled ........................................... 577

16 OCEAN 4.4.6 Issues ..................................................... 579
  Mixed-Signal in OCEAN 4.4.6 ........................................ 579

Index ................................................................. 581
Preface

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

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

The preface discusses the following:

- Licensing in OCEAN on page 18
- Related Documents for OCEAN on page 18
- Typographic and Syntax Conventions on page 19
- Identifiers Used to Denote Data Types on page 22
Licensing in OCEAN

For information on licensing in OCEAN, see *Virtuoso Software Licensing and Configuration Guide*.

Related Documents for OCEAN

OCEAN is based on the Virtuoso® SKILL programming language. The following manuals give you more information about the SKILL language and other related products.

**Installation, Environment, and Infrastructure**

- For information on installing Cadence products, see the *Cadence Installation Guide*.
- For information on the Virtuoso design environment, see the *Virtuoso Design Environment User Guide*.
- The *Cadence 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 *Cadence SKILL Language Reference* provides descriptions, syntax, and examples for the SKILL and SKILL++ functions.
- The *Cadence SKILL++ Object System Reference* provides descriptions, syntax, and examples for the object system functions.
- The *Virtuoso Design Environment SKILL Reference* describes database SKILL functions, including data access functions.
- The *Virtuoso Design Environment SKILL Reference* describes database SKILL functions, including data access functions.
- The *Virtuoso Analog Design Environment L SKILL Language Reference* provides descriptions, syntax, and examples for the SKILL commands supported by Virtuoso Analog Design Environment L.
- The *Virtuoso Analog Design Environment XL SKILL Language Reference* provides descriptions, syntax, and examples for the SKILL commands supported by Virtuoso Analog Design Environment XL and Virtuoso Analog Design Environment XL.
Virtuoso Tools

- The *Virtuoso Analog Design Environment L User Guide* explains how to design and simulate analog circuits using Virtuoso Analog Design Environment L.

- The *Virtuoso Analog Design Environment XL User Guide* explains how to design and simulate analog circuits using Virtuoso Analog Design Environment XL.

- The *Virtuoso Analog Design Environment GXL User Guide* explains how to design and simulate analog circuits using Virtuoso Analog Design Environment GXL.


- 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.

**literal**
Nonitalic words indicate keywords that you must type literally. These keywords represent command (function, routine) or option names.

**argument (z_argument)**
Words in italics indicate user-defined arguments for which you must substitute a name or a value. (The characters before the underscore (_) in the word indicate the data types that this argument can take. Names are case sensitive. Do not type the underscore (z_) before your arguments.) For a listing of data types, see “Data Types Used in OCEAN” on page 30.

| Vertical bars (OR-bars) separate possible choices for a single argument. They take precedence over any other character.

[ ] Brackets denote optional arguments. When used with OR-bars, they enclose a list of choices. You can choose one argument from the list.

{ } Braces are used with OR-bars and enclose a list of choices. You must choose one argument from the list.
Three dots (…) 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.

\textit{argument}... Specify at least one, but more are possible.

\textit{[argument]}... Specify zero or more.

A comma and three dots together indicate that if you specify more than one argument, you must separate those arguments by commas.

\texttt{=>} 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.

\texttt{/} A slash separates the possible values that can be returned by a SKILL function.

\texttt{<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.

\textbf{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. For information on the SKILL language, see the \textit{Cadence SKILL Language User Guide}.

**Example 1**

\begin{verbatim}
list( g_arg1 [g_arg2] ... )
  => l_result
\end{verbatim}
Example 1 illustrates the following syntax characters.

- **list**
  Plain type indicates words that you must type literally.

- **g_arg1**
  Words in italics indicate arguments for which you must substitute a name or a value.

- **( )**
  Parentheses separate names of functions from their arguments.

- **_**
  An underscore separates an argument type (left) from an argument name (right).

- **[ ]**
  Brackets indicate that the enclosed argument is optional.

- **=>**
  A right arrow points to the return values of the function. Also used in code examples in SKILL manuals.

- **...**
  Three dots indicate that the preceding item can appear any number of times.

### Example 2

```plaintext
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.
Identifiers Used to Denote Data Types

OCEAN commands support several data types to identify the type of value you can assign to an argument.

Data types are identified by a single letter followed by an underscore; for example, t is the data type in t_viewNames. Data types and the underscore are used as identifiers only; they should not be typed.

<table>
<thead>
<tr>
<th>Prefix</th>
<th>Internal Name</th>
<th>Data Type</th>
</tr>
</thead>
<tbody>
<tr>
<td>a</td>
<td>array</td>
<td>array</td>
</tr>
<tr>
<td>b</td>
<td>ddUserType</td>
<td>Boolean</td>
</tr>
<tr>
<td>C</td>
<td>opfcontext</td>
<td>OPF context</td>
</tr>
<tr>
<td>d</td>
<td>dbobject</td>
<td>Cadence database object (CDBA)</td>
</tr>
<tr>
<td>e</td>
<td>envobj</td>
<td>environment</td>
</tr>
<tr>
<td>f</td>
<td>flonum</td>
<td>floating-point number</td>
</tr>
<tr>
<td>F</td>
<td>opffile</td>
<td>OPF file ID</td>
</tr>
<tr>
<td>g</td>
<td>general</td>
<td>any data type</td>
</tr>
<tr>
<td>G</td>
<td>gdmSpecIIUserType</td>
<td>gdm spec</td>
</tr>
<tr>
<td>h</td>
<td>hdbobject</td>
<td>hierarchical database configuration object</td>
</tr>
<tr>
<td>l</td>
<td>list</td>
<td>linked list</td>
</tr>
<tr>
<td>m</td>
<td>nmpIIUserType</td>
<td>nmpII user type</td>
</tr>
<tr>
<td>M</td>
<td>cdsEvalObject</td>
<td>—</td>
</tr>
<tr>
<td>n</td>
<td>number</td>
<td>integer or floating-point number</td>
</tr>
<tr>
<td>o</td>
<td>userType</td>
<td>user-defined type (other)</td>
</tr>
<tr>
<td>p</td>
<td>port</td>
<td>I/O port</td>
</tr>
<tr>
<td>q</td>
<td>gdmspecListIIUserType</td>
<td>gdm spec list</td>
</tr>
<tr>
<td>r</td>
<td>defstruct</td>
<td>defstruct</td>
</tr>
<tr>
<td>R</td>
<td>rodObj</td>
<td>relative object design (ROD) object</td>
</tr>
<tr>
<td>s</td>
<td>symbol</td>
<td>symbol</td>
</tr>
<tr>
<td>S</td>
<td>stringSymbol</td>
<td>symbol or character string</td>
</tr>
<tr>
<td>Prefix</td>
<td>Internal Name</td>
<td>Data Type</td>
</tr>
<tr>
<td>--------</td>
<td>---------------</td>
<td>-----------</td>
</tr>
<tr>
<td>t</td>
<td>string</td>
<td>character string (text)</td>
</tr>
<tr>
<td>u</td>
<td>function</td>
<td>function object, either the name of a function (symbol) or a lambda function body (list)</td>
</tr>
<tr>
<td>U</td>
<td>funobj</td>
<td>function object</td>
</tr>
<tr>
<td>v</td>
<td>hdbpath</td>
<td>—</td>
</tr>
<tr>
<td>w</td>
<td>wtype</td>
<td>window type</td>
</tr>
<tr>
<td>x</td>
<td>integer</td>
<td>integer number</td>
</tr>
<tr>
<td>y</td>
<td>binary</td>
<td>binary function</td>
</tr>
<tr>
<td>&amp;</td>
<td>pointer</td>
<td>pointer type</td>
</tr>
</tbody>
</table>
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 26
- OCEAN Online Help on page 26
- OCEAN Syntax Overview on page 27
- Parametric Analysis on page 33
- Distributed Processing on page 34

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 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.

- **Simulation Set-up Commands**
  - Specify the analyses to be run
  - Specify the nets and currents to save
  - Specify the simulator option values
  - Specify the circuit stimulus

- **Simulator Run Command**
  - Run the simulator

- **Data Access Commands**
  - Perform calculations on the results
  - Print information
  - Plot waveforms

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

You can use the `history` command to view the command history from the current session and the most recently terminated session.

**OCEAN Online Help**

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

```
ocnHelp('commandName')
```

This command returns an explanation of the command and examples of how the command can be used.
To get a listing of all the different types of commands in OCEAN, type the following:

\texttt{ocnHelp()}

For more information, see “\texttt{ocnHelp}” on page 159.

**OCEAN Syntax Overview**

OCEAN is based on the Virtuoso\textsuperscript{®} SKILL programming language and uses SKILL syntax. All the SKILL language commands can be used in OCEAN. This includes \texttt{if} statements, \texttt{case} statements, \texttt{for} loops, \texttt{while} loops, \texttt{read} commands, \texttt{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 \texttt{path} command.

\begin{verbatim}
path ( "~/simulation1/schematic/psf" "~/simulation2/schematic/psf" )
\end{verbatim}

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

\begin{verbatim}
path ( "~/simulation1/schematic/psf" "~/simulation2/schematic/psf" )
\end{verbatim}
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, \texttt{tran} is a symbol and must be preceded by a single quotation mark. The symbol \texttt{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 85, you see that the valid values for the first argument include \texttt{'tran}.

\begin{verbatim}
analysis( 'tran ... )
\end{verbatim}

The list of items you can save with the \texttt{save} command is also predefined. You must choose from this predefined list. See “\texttt{save}” on page 135 and refer to the valid values for the \texttt{s_saveType} argument. The \texttt{'v} symbol indicates that the item to be saved is the voltage on a net.

\begin{verbatim}
save( 'v "net1" )
\end{verbatim}

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 \texttt{s}, you must substitute a symbol (preceded by a single quotation mark) for the argument. Consider the following syntax statement:

\begin{verbatim}
selectResults( s_resultsName ) => t / nil
\end{verbatim}

In this case, there is one symbol that must be supplied: \texttt{s_resultsName}. For the \texttt{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:

\begin{verbatim}
results() => ( dc tran ac )
\end{verbatim}

\textbf{Note:} Depending on which results are selected, the values returned by the \texttt{results} command vary.

Question Mark

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

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.

```
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.

<table>
<thead>
<tr>
<th>Data Type</th>
<th>Internal Name</th>
<th>Prefix</th>
</tr>
</thead>
<tbody>
<tr>
<td>floating-point number</td>
<td>flonum</td>
<td>f</td>
</tr>
<tr>
<td>any data type</td>
<td>general</td>
<td>g</td>
</tr>
<tr>
<td>linked list</td>
<td>list</td>
<td>l</td>
</tr>
<tr>
<td>integer, floating-point number, or complex number</td>
<td>n</td>
<td></td>
</tr>
</tbody>
</table>
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.

<table>
<thead>
<tr>
<th>Assigning a Return Value to a Variable</th>
<th>Resulting Value for the Variable</th>
</tr>
</thead>
<tbody>
<tr>
<td>a=desVar(&quot;r1&quot; 1k)</td>
<td>a=1k</td>
</tr>
<tr>
<td>a=desVar(&quot;r1&quot; 1k &quot;r2&quot; 2k)</td>
<td>a=2k</td>
</tr>
<tr>
<td>a=desVar(&quot;r1&quot;)</td>
<td>a=1k, assuming r1 was set in a desVar command</td>
</tr>
<tr>
<td>a=desVar(&quot;r2&quot;)</td>
<td>a=2k, assuming r2 was set in a desVar command</td>
</tr>
</tbody>
</table>

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 )
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:

\[
\begin{align*}
\text{desVar( "b" 1 )} \\
\text{desVar( "a" "b+1" )}
\end{align*}
\]

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:

\[
\text{desVar("b" 2)}
\]

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:

\[
\text{R1 1 0 resistor r=b}
\]

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.

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

**outputs() in OCEAN**

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

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

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:

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

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

- Loop through `rl` 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 `rl=200`, `rs` equals 300, 500 and then 700. Then, for `rl=400`, `rs` equals 300, 500, and then 700. Finally, for `rl=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.

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

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

### Data Access Without Running a Simulation

You can retrieve and use data from previous simulations at any time by opening the data with the `openResults` command. After opening the data, you can use any data access commands on this data. For more information, see 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**

The simulation results can be plotted in WaveScan, which is supported in the OCEAN environment.
Using OCEAN

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

- OCEAN Use Models on page 37
- Using OCEAN Interactively on page 38
- License Requirements
- Creating OCEAN Scripts on page 42
- Running Multiple Simulators on page 48
- OCEAN Tips on page 48

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 42.

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 26.

Note: The current version of OCEAN has some specific issues that are addressed in Appendix 16, “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 command loads and reads your .cdsinit file first and then loads and reads the .oceanrc file. You can place OCEAN commands in your .oceanrc file, which is similar to the .cdsinit file. This file can contain any valid OCEAN command, function or SKILL initialization routine (excluding graphical dfII 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 40.

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

```
extit
```

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:

```
#!/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.

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

The -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 followed by a newline. Otherwise, OCEAN hangs. The reason for this is that when the workbench is started in the non-graphical mode, it does not redirect standard I/O as it normally does; instead, it lets the SKILL human interface (HI) handle the standard I/O. HI expects an 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 L 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 `virtuoso` command). Therefore, you might want to load your `.oceanrc` file manually in the CIW if you need information that it contains.

You can also use the `history` command from the CIW to list and reuse the most recently used commands.

**Interactive Session Demonstrating the OCEAN Use Model**

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

- **Start OCEAN and specify your simulator.**
  ```
  ocean
  simulator( ’spectre )
  ```

- **Specify an AC analysis.**
  ```
  analysis( ’ac ?start 1 ?stop 1000 ?lin 100 )
  or
  ac( 1 1000 ”linear” 100 )
  ```

- **Set a design variable.**
  ```
  desVar( ”rs” 1k )
  ```

- **Perform the first simulation run.**
  ```
  run( )
  ```

- **Change a design variable.**
  ```
  desVar( ”rs” 2k )
  ```

- **Perform the second simulation run.**
  ```
  run( )
  ```

- **Specify a transient analysis.**
  ```
  analysis( ’tran ?stop 1u ) (Spectre only)
  or
  tran( 0 1u 1n )
  ```

- **Perform the third simulation run.**
  ```
  run( )
  ```
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.

License Requirements

You need licenses to run the simulator() and ocnSetXLMode() OCEAN commands. For more information on these commands, see simulator on page 139 and ocnSetXLMode on page 504.

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

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

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

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

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

**Note:** If the alwaysTryHigherTieredLicenseInOcean environment variable is set to nil, errors are displayed if OCEAN is unable to checkout a license.
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 45.

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:

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

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 Virtuoso Analog Design Environment L or ,

1. Start the Virtuoso software,
   ```
   virtuoso
   ```
   The CIW appears.
2. From the CIW, choose *Tools – ADE L – Simulation*.
   The Virtuoso Analog Design Environment window appears.
3. Perform all of the design environment tasks that you want to capture in the script.
4. Choose *Session – Save Script*.
   The Save Ocean Script to File form appears.
5. Click *OK* to accept the default file name (`./oceanScript.ocn`), or change the name for the file and click *OK*.

   A script containing the OCEAN commands for the tasks you performed is created. For information about how to load this script, see “Loading OCEAN Scripts” on page 45.

**Controlling What Is Included in Scripts**

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

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

```
asimenv.misc saveDefaultsToOCEAN boolean t
```

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 L SKILL Language Reference*.

3. Load the functions in your .cdsinit file.

   For example, you might add the following lines to your .cdsenv file.
   ```
asimenv.misc preSaveOceanScript string "MYfirstProc"
asimenv.misc postSaveOceanScript string "MYlastProc"
```

   The functions `MYfirstProc` and `MYlastProc` might be defined like this.
procedure( MYfirstProc( session fp)
    fprintf(fp "; This will be the first line in the ocean script.\n")
)

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

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

load "myOceanProcs.il"

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

Loading OCEAN Scripts

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

From a UNIX Shell

To load an OCEAN script,
1. Type the following command to start OCEAN:
   ocean
   The OCEAN prompt appears.
2. Use the SKILL load command to load your script:
   load( "script_name.ocn" )
   Messages about the progress of your script appear.

From the CIW

To load an OCEAN script,
1. Start the Virtuoso software
   virtuoso &
   The CIW appears.
2. In the text entry field, use the SKILL load command to load your script:
   load( "script_name.ocn" )
Messages about the progress of your script appear in the CIW.

**Selecting Results**

You may use OCEAN to run several simulations on the same design and save the results in different result directories. You can then use 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 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 `virtuoso`
- `testSchemName` is your test schematic

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

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

1. Choose *Artist -> Results -> Select*

2. The *Select Results* form opens. Click *Browse*. A Unix Browser form appears.

3. Navigate to the directory that contains your Ocean generated directories c1, c2, and c3.

4. Click *OK* on the Unix Browser form. Now the *Select Results* Form should show c1, c2 and c3.

5. Double click 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 -
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 virtuoso

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".

   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:
  ```skill
  *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:
  ```skill
  analysis('tran ?from fromVal)
  ```
  However, the following pair of statements work correctly because `fromVal` has a value (is bound).
  ```skill
  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:
  ```skill
  ocean-Debug 2>
  ```
  You can continue working. However, if you would like to get out of the debugger, you can type
  ```skill
  debugQuit()
  ```
  Now you are back to the normal prompt:
  ```skill
  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 28.

- 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 51
- Naming Conventions on page 52
- Arithmetic Operators on page 52
- Scaling Factors on page 52
- Relational and Logical Operators on page 54
- SKILL Syntax on page 56
- Arithmetic and Logical Expressions on page 59

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**

<table>
<thead>
<tr>
<th>Operators in Descending Precedence</th>
<th>Underlying Function</th>
</tr>
</thead>
<tbody>
<tr>
<td>** exponentiation</td>
<td></td>
</tr>
<tr>
<td>* multiply</td>
<td></td>
</tr>
<tr>
<td>/ divide</td>
<td></td>
</tr>
<tr>
<td>+ plus</td>
<td></td>
</tr>
<tr>
<td>- minus</td>
<td></td>
</tr>
<tr>
<td>== equal</td>
<td></td>
</tr>
<tr>
<td>!= nequal</td>
<td></td>
</tr>
<tr>
<td>= assignment</td>
<td></td>
</tr>
</tbody>
</table>

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.
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**

<table>
<thead>
<tr>
<th>Character</th>
<th>Name</th>
<th>Multiplier</th>
<th>Examples</th>
</tr>
</thead>
<tbody>
<tr>
<td>Y</td>
<td>Yotta</td>
<td>1024</td>
<td>10Y [10e+25]</td>
</tr>
<tr>
<td>Z</td>
<td>Zetta</td>
<td>1021</td>
<td>10Z [10e+22]</td>
</tr>
<tr>
<td>E</td>
<td>Exa</td>
<td>1018</td>
<td>10E [10e+19]</td>
</tr>
<tr>
<td>P</td>
<td>Peta</td>
<td>1015</td>
<td>10P [10e+16]</td>
</tr>
<tr>
<td>T</td>
<td>Tera</td>
<td>10^{12}</td>
<td>10T [1.0e13]</td>
</tr>
<tr>
<td>G</td>
<td>Giga</td>
<td>10^{9}</td>
<td>10G [10,000,000,000]</td>
</tr>
<tr>
<td>M</td>
<td>Mega</td>
<td>10^{6}</td>
<td>10M [10,000,000]</td>
</tr>
<tr>
<td>K</td>
<td>Kilo</td>
<td>10^{3}</td>
<td>10K [10,000]</td>
</tr>
<tr>
<td>%</td>
<td>percent</td>
<td>10^{-2}</td>
<td>5% [0.05]</td>
</tr>
<tr>
<td>m</td>
<td>milli</td>
<td>10^{-3}</td>
<td>5m [5.0e-3]</td>
</tr>
<tr>
<td>u</td>
<td>micro</td>
<td>10^{-6}</td>
<td>1.2u [1.2e-6]</td>
</tr>
<tr>
<td>n</td>
<td>nano</td>
<td>10^{-9}</td>
<td>1.2n [1.2e-9]</td>
</tr>
<tr>
<td>p</td>
<td>pico</td>
<td>10^{-12}</td>
<td>1.2p [1.2e-12]</td>
</tr>
<tr>
<td>f</td>
<td>femto</td>
<td>10^{-15}</td>
<td>1.2f [1.2e-15]</td>
</tr>
<tr>
<td>a</td>
<td>atto</td>
<td>10^{-18}</td>
<td>1.2a [1.2e-18]</td>
</tr>
<tr>
<td>z</td>
<td>zepto</td>
<td>10^{-21}</td>
<td>1.2z [1.2e-21]</td>
</tr>
<tr>
<td>y</td>
<td>yocto</td>
<td>10^{-24}</td>
<td>1.2y [1.2e-24]</td>
</tr>
</tbody>
</table>

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

This section introduces you to

- Relational Operators: <, <=, >, >=, ==, !=
- Logical Operators: !, &&, ||

Relational Operators

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

Sample Relational Operators

<table>
<thead>
<tr>
<th>Operator</th>
<th>Arguments</th>
<th>Function</th>
<th>Example</th>
<th>Return Value</th>
</tr>
</thead>
<tbody>
<tr>
<td>&lt;</td>
<td>numeric</td>
<td>lessp</td>
<td>3 &lt; 5</td>
<td>t</td>
</tr>
<tr>
<td></td>
<td></td>
<td></td>
<td>3 &lt; 2</td>
<td>nil</td>
</tr>
<tr>
<td>&lt;=</td>
<td>numeric</td>
<td>leqp</td>
<td>3 &lt;= 4</td>
<td>t</td>
</tr>
<tr>
<td>&gt;</td>
<td>numeric</td>
<td>greaterp</td>
<td>5 &gt; 3</td>
<td>t</td>
</tr>
<tr>
<td>&gt;=</td>
<td>numeric</td>
<td>geqp</td>
<td>4 &gt;= 3</td>
<td>t</td>
</tr>
<tr>
<td>==</td>
<td>numeric</td>
<td>equal</td>
<td>3.0 == 3</td>
<td>t</td>
</tr>
<tr>
<td></td>
<td>string</td>
<td></td>
<td>&quot;abc&quot; == &quot;ABc&quot;</td>
<td>nil</td>
</tr>
<tr>
<td>!=</td>
<td>numeric</td>
<td>nequal</td>
<td>&quot;abc&quot; != &quot;ABc&quot;</td>
<td>t</td>
</tr>
<tr>
<td></td>
<td>string</td>
<td></td>
<td>&quot;abc&quot; != &quot;ABc&quot;</td>
<td>t</td>
</tr>
</tbody>
</table>

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

```
l > "abc"
Message: *Error* greaterp: can’t handle (l > "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

<table>
<thead>
<tr>
<th>Operator</th>
<th>Arguments</th>
<th>Function</th>
<th>Example</th>
<th>Return Value</th>
</tr>
</thead>
<tbody>
<tr>
<td>&amp;&amp;</td>
<td>general</td>
<td>and</td>
<td>3 &amp; 5</td>
<td>5</td>
</tr>
<tr>
<td></td>
<td></td>
<td></td>
<td>5 &amp; 3</td>
<td>3</td>
</tr>
<tr>
<td></td>
<td></td>
<td></td>
<td>t &amp; nil</td>
<td>nil</td>
</tr>
<tr>
<td></td>
<td></td>
<td></td>
<td>nil &amp; t</td>
<td>nil</td>
</tr>
<tr>
<td></td>
<td></td>
<td></td>
<td>general</td>
<td>or</td>
</tr>
<tr>
<td></td>
<td></td>
<td></td>
<td>5</td>
<td></td>
</tr>
<tr>
<td></td>
<td></td>
<td></td>
<td>t</td>
<td></td>
</tr>
<tr>
<td></td>
<td></td>
<td></td>
<td>nil</td>
<td></td>
</tr>
</tbody>
</table>

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 nonalphabetic 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**

<table>
<thead>
<tr>
<th>Character</th>
<th>Name</th>
<th>Meaning</th>
</tr>
</thead>
<tbody>
<tr>
<td>\</td>
<td>backslash</td>
<td>Escape for special characters</td>
</tr>
<tr>
<td>( )</td>
<td>parentheses</td>
<td>Grouping of list elements, function calls</td>
</tr>
<tr>
<td>[ ]</td>
<td>brackets</td>
<td>Array index, super right bracket</td>
</tr>
<tr>
<td>'</td>
<td>single quotation mark</td>
<td>Specifies a symbol (quoting the expression prevents its evaluation)</td>
</tr>
<tr>
<td>&quot;</td>
<td>quotation mark</td>
<td>String delimiter</td>
</tr>
<tr>
<td>,</td>
<td>comma</td>
<td>Optional delimiter between list elements</td>
</tr>
</tbody>
</table>
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:

```plaintext
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

```c
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

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

as well as in the more conventional algebraic form

```plaintext
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
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.
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 63
- Data Types on page 63
- Arrays on page 66
- Concatenating Strings (Lists) on page 66
- Declaring a SKILL Function on page 68
- Skill Function Return Values on page 70
- Syntax Functions for Defining Functions on page 70

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

**Sample SKILL Data Items**

<table>
<thead>
<tr>
<th>Data Type</th>
<th>Syntax Example</th>
</tr>
</thead>
<tbody>
<tr>
<td>integer</td>
<td>5</td>
</tr>
<tr>
<td>floating point number</td>
<td>5.3</td>
</tr>
<tr>
<td>text string</td>
<td>&quot;Mary had a little lamb&quot;</td>
</tr>
</tbody>
</table>

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**

<table>
<thead>
<tr>
<th>Data Type</th>
<th>Internal Name</th>
<th>Prefix</th>
</tr>
</thead>
<tbody>
<tr>
<td>array</td>
<td>array</td>
<td>a</td>
</tr>
<tr>
<td>boolean</td>
<td></td>
<td>b</td>
</tr>
<tr>
<td>floating-point number</td>
<td>flonum</td>
<td>f</td>
</tr>
<tr>
<td>any data type</td>
<td>general</td>
<td>g</td>
</tr>
<tr>
<td>linked list</td>
<td>list</td>
<td>l</td>
</tr>
<tr>
<td>floating-point number or integer</td>
<td>n</td>
<td></td>
</tr>
<tr>
<td>user-defined type</td>
<td>o</td>
<td></td>
</tr>
<tr>
<td>I/O port</td>
<td>port</td>
<td>p</td>
</tr>
<tr>
<td>symbol</td>
<td>symbol</td>
<td>s</td>
</tr>
<tr>
<td>symbol or character string</td>
<td>S</td>
<td></td>
</tr>
<tr>
<td>character string (text)</td>
<td>string</td>
<td>t</td>
</tr>
<tr>
<td>window type</td>
<td></td>
<td>w</td>
</tr>
<tr>
<td>integer number</td>
<td>fixnum</td>
<td>x</td>
</tr>
</tbody>
</table>

**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 52.

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.
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 true 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.

<table>
<thead>
<tr>
<th>Return Value</th>
<th>Meaning</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td><code>string1</code> is alphabetically greater than <code>string2</code>.</td>
</tr>
<tr>
<td>0</td>
<td><code>string1</code> is alphabetically equal to <code>string2</code>.</td>
</tr>
<tr>
<td>-1</td>
<td><code>string1</code> is alphabetically less than <code>string2</code>.</td>
</tr>
</tbody>
</table>

```
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 `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.

<table>
<thead>
<tr>
<th>Return Value</th>
<th>Meaning</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td><code>arg1</code> is alphanumerically greater than <code>arg2</code>.</td>
</tr>
<tr>
<td>0</td>
<td><code>arg1</code> is alphanumerically identical to <code>arg2</code>.</td>
</tr>
<tr>
<td>-1</td>
<td><code>arg2</code> is alphanumerically greater than <code>arg1</code>.</td>
</tr>
</tbody>
</table>

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.

```skill
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:

```skill
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:

```plaintext
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.
The actual arguments in a function call correspond to the formal arguments in the declaration of the function.

<table>
<thead>
<tr>
<th>term</th>
<th>definition</th>
</tr>
</thead>
<tbody>
<tr>
<td>expression</td>
<td>A use of a SKILL function, often by means of an operator supplying required parameters.</td>
</tr>
<tr>
<td>function body</td>
<td>The collection of SKILL expressions that define the function’s algorithm.</td>
</tr>
</tbody>
</table>
OCEAN Environment Commands

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

appendPath on page 74

path on page 75

prependPath on page 76

setup on page 77

history on page 79

ocnSetSilentMode on page 80
appendPath

appendPath( _dirName1 ... _dirNameN)
   => _dirNameN/nil

Description

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

Arguments

t_dirName1        Directory path.
t_dirNameN        Additional directory path.

Value Returned

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

Example

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

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

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

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

**OCEAN Environment Commands**

---

**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**

- `t_dirName1` Directory path.
- `t_dirNameN` Additional directory path.

**Value Returned**

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

**Example**

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

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

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

Returns the search path last set.
prependPath

prependPath( \[t\_dirName1 \ldots \ [t\_dirNameN]\])
  => undefined/nil

Description

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

Arguments

\[t\_dirName1\] Directory path.
\[t\_dirNameN\] Additional directory path.

Value Returned

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

Example

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

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

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

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

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

Returns the search path last set.
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:

<table>
<thead>
<tr>
<th>Name1</th>
<th>Param1</th>
<th>Param2</th>
<th>Param3</th>
</tr>
</thead>
<tbody>
<tr>
<td>value</td>
<td>value</td>
<td>value</td>
<td></td>
</tr>
</tbody>
</table>

<table>
<thead>
<tr>
<th>Name2</th>
<th>Param1</th>
<th>Param2</th>
<th>Param3</th>
</tr>
</thead>
<tbody>
<tr>
<td>value</td>
<td>value</td>
<td>value</td>
<td></td>
</tr>
</tbody>
</table>

<table>
<thead>
<tr>
<th>Name3</th>
<th>Param1</th>
<th>Param2</th>
<th>Param3</th>
</tr>
</thead>
<tbody>
<tr>
<td>value</td>
<td>value</td>
<td>value</td>
<td></td>
</tr>
</tbody>
</table>
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

\texttt{x\_charsPerLine} \quad \text{Specifies the number of characters per line output to the display.}
\text{Default value: 80}

\texttt{g\_messageOn} \quad \text{Specifies whether error messages are turned on.}
\text{Valid values: \texttt{t, nil}}
\text{Default value: \texttt{t}, which specifies that messages are turned on.}

\textbf{Value Returned}

\texttt{t} \quad \text{Returns \texttt{t} if the value is assigned to the name.}

\texttt{nil} \quad \text{Returns \texttt{nil} if there is a problem.}

\textbf{Example}

\texttt{setup( ?numberNotation 'engineering )}
\Rightarrow \texttt{t}

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

\texttt{setup( ?precision 5 )}
\Rightarrow \texttt{t}

Specifies that 5 significant digits are to be printed.

\texttt{setup(?numberNotation 'suffix ?charsPerLine 40 ?reportStyle 'spice ?messageOn t)}

Sets up number notation to \texttt{suffix} format, characters per line to 40, reporting style to \texttt{Spice}, and error message to \texttt{ON}. 
**history**

`history( [x_number] )`  =>  `t`

**Description**

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

**Arguments**

`x_number`  
The number of previously entered commands to be listed.  
Default value: 20

**Value Returned**

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

**Example**

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

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

- `ocean> !1`
  
  This executes the command numbered 1, which in this example is `simulator('spectre')`.
  
  **Note:** This command currently works from the ocean prompt but not from the CIW.
  
- `ocean> !des`
  
  This executes the last command whose prefix starts with `des` in the history. In this example, it is the second command listed, that is, `design( "tests" "simple" "schematic")`.  

**ocnSetSilentMode**

\[
\text{ocnSetSilentMode( b\_silentMode )} \\
=> t
\]

**Description**

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

**Arguments**

\[
b\_silentMode \\
\text{Accepts boolean values t or nil.}
\]

- Set to \(t\) to suppress the OCEAN warning and information messages.
- Set to \(\text{nil}\) to allow all OCEAN messages to be displayed.

**Value Returned**

\[
t \\
\text{Returns t to indicate the successful assignment of the passed argument.}
\]

**Example**

\[
\text{ocnSetSilentMode(t) => t}
\]

Suppresses the ocean warning messages

\[
\text{ocnSetSilentMode(nil) => t}
\]

Displays the ocean warning messages
Simulation Commands

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

- `ac` on page 83
- `analysis` on page 85
- `converge` on page 88
- `connectRules` on page 89
- `createFinalNetlist` on page 93
- `createNetlist` on page 94
- `dc` on page 96
- `definitionFile` on page 98
- `delete` on page 99
- `design` on page 101
- `desVar` on page 103
- `discipline` on page 105
- `displayNetlist` on page 107
- `envOption` on page 108
- `evcdFile` on page 110
- `evcdInfoFile` on page 111
- `forcenode` on page 112
- `globalSigAlias` on page 113
- `globalSignal` on page 114
- `ic` on page 116
- `includeFile` on page 117
modelFile on page 118
nodeset on page 119
noise on page 120
ocnCloseSession on page 121
ocnDisplay on page 122
ocnGetAdjustedPath on page 124
ocnWaveformTool on page 125
off on page 126
option on page 127
restore on page 129
resultsDir on page 130
run on page 131
save on page 135
saveOption on page 137
simulator on page 139
solver on page 140
stimulusFile on page 141
store on page 143
temp on page 144
tran on page 145
vcdFile on page 146
vcdInfoFile on page 147
vecFile on page 148
hlcheck on page 149
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

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

Value Returned

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

Example

ac(1 10000 2)

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

ac(1 10000 "Linear" 100)
Specifies an AC analysis from 1 to 10,000 by 100.

\texttt{ac(1 5000 "Logarithmic" 10)}

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

```plaintext
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 *Virtuoso Spectre Circuit Simulator Reference* 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.)
**Value Returned**

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

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

**Example**

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

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

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

Specifies that a transient analysis be performed.

```lisp
```

Sweeps temperature for the Spectre® circuit simulator.

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

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

```lisp
```

Sets the Spectre transfer function analysis.

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

Sets the Spectre sensitivity analysis.

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

Sets the Spectre noise analysis.

```lisp
analysis('dcmatch ?oprobe "/PR1" )
```

```lisp
analysis('dcmatch ?param "temp" ?start "24" ?stop "26 ?lin "5"
```

Sets the Spectre domatch analysis.

```lisp
```

Sets the Spectre pz analysis.

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" ?reffieldband "0" )

Sets the Spectre pnoise RF analysis.

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

Sets the Spectre pac RF analysis.


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.
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 L User Guide*.

Arguments

- **s_convName**  
  Name of the convergence type. Valid values are one of *nodeset* and *forcenode*. Note that *forcenode* is not supported for the spectre simulator.

- **t_netName1**  
  Name of the net to which you want to set convergence criteria.

- **f_value1**  
  Voltage value for the net

- **t_netNameN**  
  Name of the additional net

- **f_value**  
  Voltage value for the additional net

Value Returned

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

- **nil**  
  Returns nil and prints an error message if the function fails

Example

- `converge( 'ic "/I0/net1" 5 )`
  Sets the convergence name for the initial condition net1 to 5 volts.

- `converge( 'nodeset "/I0/net1" 5 )`
  Sets the convergence name for nodeset of net1 to 5 volts.
connectRules

connectRules( t.ruleName [?lib t_libName] [?view t_viewName] [baseRule t_baseRule] [?moduleInfo l_moduleInfo] [?resolutionInfo l_resolutionInfo] [?commonParam l_commonParam] [userDefined s_userDefined] )
=> t / nil

connectRules( t_ruleName )
=> t / nil

connectRules( ?none s_tag )
=> t / nil

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

l_moduleInfo: ((s.moduleName1 [?mode s_mode] [?paramInfo l_paramInfo] [?direction1 s_direction1][?discipline1 s_discipline1] [?direction2 s_direction2] [?discipline2 s_discipline2]) [(s.moduleName2 - ) -]

l_paramInfo: ((s.paramName1 s_value1) [(s.paramName2 s_value2) -])

l_resolutionInfo: ((s.resolvedDiscipline1 s_equivalentDisciplines1) [(s.resolvedDiscipline2 s_equivalentDisciplines2) -])

l_commonParam: ((s.paramName1 s_value1 [(s.moduleName1 s_moduleName2 - )]) [(s.paramName2 s_value2 ) -])

Description

Sets connect rules for a given AMS OCEAN session required by the elaborator. To specify multiple connect rules, use this command multiple times. To add a connect rule to an OCEAN session, you can either choose a built-in rule from the connectLib library (by specifying t_ruleName, t_libName and t_viewName) or one of your own compiled built-in connect rules (by specifying t_ruleName, t_libName and t_viewName). To add a user defined connect rule to an OCEAN session specify s_userDefined. To modify an existing built-in rule, you need to specify t_baseRule (the name of the built-in rule that needs be modified), specify a new name (by specifying t_ruleName, t_libName and t_viewName) and also specify one or more of the optional arguments.

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

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

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

\textit{t\_ruleName}\hspace{1cm} Name of the connect rule that you want to use in the current session.

\textit{t\_libName}\hspace{1cm} Name of the library that contains a list of user-compiled connect rules. If you do not specify this, the connect rules are assumed to be in the default location.

\textit{t\_viewName}\hspace{1cm} Name of the view of the selected cell.

\textit{t\_baseRule}\hspace{1cm} Name of the connect rule that you want to modify.

\textit{l\_moduleInfo}\hspace{1cm} Arguments that need to be updated for a specified connect rule. The arguments may include \texttt{s\_mode}, \texttt{s\_direction1}, \texttt{s\_direction2}, \texttt{s\_discipline1} and \texttt{s\_discipline2}.

Valid values for \texttt{s\_mode} are: \texttt{null}, \texttt{split}, \texttt{merged}.

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

\texttt{s\_discipline1} and \texttt{s\_discipline2} also work as a pair. Either they should both be \texttt{null} or they should both have values.

\textit{t\_resolutionInfo}\hspace{1cm} Names of disciplines that need to be resolved to another discipline. The value specified overwrites the \textit{l\_resolutionInfo} in the base rule or in the existing connect rule.

\textit{t\_commonParam}\hspace{1cm} One or more parameters that you want to modify for all modules or a set of modules. Although the same result can be achieved by using the \textit{l\_moduleInfo} argument, \textit{l\_commonParam} facilitates updating parameters for all modules in one go.

\textit{s\_userDefined}\hspace{1cm} Name of the user defined connect rule that you want to use in the current session. Specify \texttt{3step} as the value of \textit{s\_userDefined} and specify \textit{t\_ruleName}, \textit{t\_libName} and \textit{t\_viewName} to add a user defined connect rule for the Cellview-based netlister flow. Specify \texttt{irun} as the value of \textit{s\_userDefined} and specify \textit{t\_ruleName}, \texttt{s\_fileName} or both to add a user defined connect rule for the OSS-based
netlister with irun flow. Any other argument specified when adding a user defined connect rule will be ignored.

s_tag

The option used to indicate that no connect rules are to be used for the current session.

Value Returned

t

Returns t if the specified connect rules are set.

nil

Returns nil and prints an error message otherwise.

Example

connectRules("ConnRules_5V_full")

Sets ConnRules_5V_full as the current connect rule from the default connectLib located in your hierarchy.

connectRules("CustomRules_9V_high" ?lib "myConnectLib" ?view "myViewName")

Sets CustomRules_9V_high from myConnectLib, where myConnectLib contains a list of user-compiled connect rules and myViewName is the specified view name.


Checks if connRule3 exists in the session. If it does, it updates direction1 to input and direction2 to output for E2L and description for this rule. If this rule does not exist, then it takes the base values as values from ConnRules_18V_full and updates direction1, direction2, and description and names the new rule as connRule3.


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


If connRule3 does not exist and the base rule is not specified but description, moduleInfo and resolutionInfo are specified, the connect rule connRule3 is added with the values specified for moduleInfo, resolutionInfo and description. Note that in this case no checks are done (that is, module names and parameter names are not
checked against base information as no base rule information is available). This command is applicable while using the connectRules command as saved in ocean.

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

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

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

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

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

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

connectRules("connRule3" ?userDefined irun)

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

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

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

connectRules(?userDefined irun ?file "file1")

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

connectRules(?none t)
=> t

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

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

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

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

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

createFinalNetlist()
   => t / nil

Description

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

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

Arguments

None.

Value Returned

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

Example

createFinalNetlist()

Creates the final netlist for the current simulation run.
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 netlisted, 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 it set to `t (true).

Note: This command does not work with socket simulators.

Arguments

b_recreateAll

If set and the design is specified as lib/cell/view, the entire netlist is recreated.

Value Returned

t_fileName

Returns the name of the simulator input file on success.

nil

otherwise nil is returned

Example

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.
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/nl
dc

\[ dc( t\_compName [ t\_compParam ] g\_fromValue g\_toValue g\_byValue ) \]
\[ => \text{undefined/nil} \]

**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 and spectreVerilog simulators.

**Arguments**

- **`t\_compName`**
  - Name of the source (or component, for the Spectre® circuit simulator) to sweep.

- **`t\_compParam`**
  - For the Spectre® circuit simulator, the component parameter to be swept.

- **`g\_fromValue`**
  - Starting value for the DC analysis.

- **`g\_toValue`**
  - Ending value.

- **`g\_byValue`**
  - The increment at which to step through the analysis.

**Value Returned**

- **`undefined`**
  - The return value for this command/function is undefined.

- **`nil`**
  - Returns nil and prints an error message if the analysis is not specified.

**Example**

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

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

```
cdc("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 L User Guide*.

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

**Arguments**

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

**Value Returned**

- **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.
delete

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

Description
Deletes all the information specified.

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

Arguments

s_command
  Command that was initially used to add the items that are now being deleted.
  Valid values: analysis, connectRules, discipline, globalSignal, desVar, path, save, ic, forcenode, nodeset

g_commandArg1
  Argument corresponding to the specified command.

g_commandArg2
  Additional argument corresponding to the specified command.

Value Returned

  t
  Returns t if the information is deleted.

  nil
  Returns nil if there is an error.

Example

delete( 'save ')
  => t

Deletes all the saves.

delete( 'save 'v ')
  => t

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_lib t_cell t_view [t_mode])
    => t_cktFile/nil | (t_lib t_cell t_view)/nil

Description

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

You can use the lib, cell, view version of the design command only if you are running
OCEAN within virtuoso. You cannot use this version of the command within the OCEAN
process itself.

Arguments

t_cktFile  For the direct simulator, the name of the netlist. The name must
end in netlist. Note that the netlistHeader and
netlistFooter files are also needed in the same directory.

For socket simulators, this is the name of the raw circuit file. If
generated in the Virtuoso® Analog Design Environment, the file
is named design.c and is found in the netlist directory.

Otherwise, cktFile is a pre-existing netlist file from another
source. In this case, you might need to remove the .cards from
the netlist because the OCEAN commands are converted to
.cards and appended to the final netlist. The simulator might
give an error or warning if the .cards are read twice.

t_lib  Name of the Virtuoso® Analog Design Environment library that
contains the design.

t_cell  Name of the design.

t_view  View of the design (typically schematic).

t_mode  The mode in which the design should be opened. The value can
be r, w or a, representing read, write and append,
respectively. The default mode is append. Read-only designs
can be netlisted only by direct netlisters, and not socket. The w
mode should not be used as it overwrites the design.
Value Returned

\texttt{t\_cktFile} \quad \text{Returns the name of the design if successful.}

\texttt{l\_(lib\ cell\ view)} \quad \text{Returns the name of the view for an Virtuoso\textsuperscript{®} Analog Design Environment design if successful.}

\texttt{nil} \quad \text{Returns nil and prints an error message if there is a problem using the specified design.}

Example

For the Spectre\textsuperscript{®} circuit simulator,

\begin{verbatim}
design( "netlist" ) => netlist

\text{specifies that netlist, a netlist file, be used in the simulation.}

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

\text{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)

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

design()
=> (mylib ampTest schematic)

\text{Returns the lib-cell-view being used in the current session. If a design has not been specified, it returns nil.}
\end{verbatim}
desVar

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

**Description**

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

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

**Arguments**

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

**Value Returned**

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

**Example**

```lisp
(desVar() )
```

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

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

Sets the `rs` design variable to 1k.

```lisp
(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.

```lisp
a = evalstring( desVar( "rs") ) / 2
```
Sets a to $\frac{1k}{2}$ or 500.

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

\[
\text{discipline}( \ g\textunderscore\text{discipline1} \ [g\textunderscore\text{discipline2} \ ...] \ ) \\
=> t / nil
\]

**Description**

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

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

**Arguments**

- `g\textunderscore\text{discipline1}` Name of a discrete discipline to be added.
- `g\textunderscore\text{discipline2}` Names of additional discrete disciplines to be added.

**Value Returned**

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

**Example**

```
\text{discipline}( \ "logic1" \ "logic2" \ "("logic3") \ )
```

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

```
\text{discipline}("LL")
```

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

```
\text{delete('discipline \ "logic2" \ "LL\")}
```

Deletes disciplines `logic2` and `LL` from the session.
Deletes all the specified disciplines in the session.
**displayNetlist**

displayNetlist()
=> t / nil

**Description**

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

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

**Arguments**

None.

**Value Returned**

- **t**
  Returns `t` if the concatenated design information file.

- **nil**
  Returns `nil` or prints an error message otherwise.

**Example**

displayNetlist()
=> t

Displays the concatenated design information file.
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 L User Guide.

Arguments

s_envOption1 Name of the first environment option to set.
g_value1 Value for the option.

s_envOptionN Name of an additional environment option to set.
g_valueN Value for the option.

Value Returned

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

Example

envOption( 'paramRangeCheckFile "./myDir/range.check" )
Sets the paramRangeCheckFile environment option.

envOption( 'initFile "./myDotSFiles/init" )
Sets the initFile environment option.
envOption( 'updateFile "./myDotSFiles/update"' )

Sets the updateFile environment option.
evcdFile

`evcdFile( t_evcdFileName )
       => t_evcdFileName/nil`

Description

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

Note: This command is applicable for the UltraSim simulators.

Arguments

`t_evcdFileName`  The name of the EVCD file to be used for session.

Value Returned

`t_evcdFileName`  The EVCD file name is the output if the command is successful.

`nil`  Otherwise, nil is returned.

Example

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

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

evcdInfoFile(t_evcdInfoFileName)
   => t_evcdInfoFileName/nil

Description

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

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

Arguments

t_evcdInfoFileName
   The name of the EVCD info file to be included.

Value Returned

 t_evcdInfoFileName
   The EVCD info file name is the output if the command is successful.

nil
   Otherwise, nil is returned.

Example

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

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

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

Arguments

<table>
<thead>
<tr>
<th>t_netName1</th>
<th>Name of the net.</th>
</tr>
</thead>
<tbody>
<tr>
<td>f_value1</td>
<td>Voltage value for the net.</td>
</tr>
<tr>
<td>t_netNameN</td>
<td>Name of an additional net.</td>
</tr>
<tr>
<td>f_valueN</td>
<td>Voltage value for the net.</td>
</tr>
</tbody>
</table>

Value Returned

<table>
<thead>
<tr>
<th>undefined</th>
<th>The return value for this command/function is undefined.</th>
</tr>
</thead>
<tbody>
<tr>
<td>nil</td>
<td>Returns nil and prints an error message.</td>
</tr>
</tbody>
</table>

Example

forcenode( "net1" 5 "net34" 2 )

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

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

Description

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

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

Arguments

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

Value Returned

t Returns t when previous signal aliases have been removed successfully and new aliases are created according to the signal lists provided.

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

Example

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

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

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

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

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

**Description**

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

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

**Arguments**

<table>
<thead>
<tr>
<th>Argument</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>t_signalName</td>
<td>The name of the global signal.</td>
</tr>
<tr>
<td>t_langName</td>
<td>The namespace within which the signal is entered. It is used to map the signal name to Verilog-AMS.</td>
</tr>
<tr>
<td></td>
<td><strong>Valid Values:</strong> CDBA, Spectre, Spice, Verilog-AMS</td>
</tr>
<tr>
<td></td>
<td><strong>Default Value:</strong> CDBA</td>
</tr>
<tr>
<td>t_wireType</td>
<td>Indicates the Verilog type of the signal declaration.</td>
</tr>
<tr>
<td></td>
<td><strong>Valid Values:</strong> wire, supply0, supply1, tri, tri0, tri1, triand, trior, trireg, wand, wor, wreal</td>
</tr>
<tr>
<td></td>
<td><strong>Default Value:</strong> wire</td>
</tr>
<tr>
<td>t_discipline</td>
<td>A string value to indicate the discipline of the signal.</td>
</tr>
<tr>
<td>t_ground</td>
<td>Indicates if the signal is a ground signal or not.</td>
</tr>
<tr>
<td></td>
<td><strong>Valid Values:</strong> YES, NO</td>
</tr>
<tr>
<td></td>
<td><strong>Default Value:</strong> NO</td>
</tr>
</tbody>
</table>

**Value Returned**

<table>
<thead>
<tr>
<th>Value</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>t</td>
<td>Returns ( t ) when a global signal has been successfully added or modified.</td>
</tr>
<tr>
<td>nil</td>
<td>Returns ( \text{nil} ) and prints an error message if the function was unsuccessful.</td>
</tr>
</tbody>
</table>
Example

globalSignal("signal1" ?wireType "tri")

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

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

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

delete('globalSignal "sig1" "sig2")

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

delete('globalSignal)

Deletes all user-specified global signals.
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 L User Guide.

Arguments

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

Value Returned

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

Example

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

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

includeFile( t_fileName ) => t_fileName/nil

Description

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

Notes:

1. This command is not available for the direct simulator. Use the modelFile or stimulusFile command instead.

2. Using this command is comparable to using the Environment Options form of the Virtuoso® Analog Design Environment to name an include file and specify that the syntax for the file be that of the target simulator. If you want the include file to be in Cadence-SPICE circuit simulator syntax, you must edit the raw netlist file (which has a .c or .C suffix), and manually add the include file.

Arguments

 t_fileName Name of the file to include in the final netlist.

Value Returned

 t_fileName Returns the name of the file if successful.

 nil Returns nil and prints an error message otherwise.

Example

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

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

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

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

Arguments

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

Value Returned

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

Example

nodeset( "net1" 5 "net34" 2 )

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

noise( t_output t_source )
   => undefined/nil

Description

Specifies a noise analysis.

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

Arguments

t_output            Output node.
t_source            Input source.

Value Returned

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

Example

noise( "n1" "v1" )

Specifies a noise analysis.
ocnCloseSession

ocnCloseSession()

=> t / nil

Description

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

Value Returned

<table>
<thead>
<tr>
<th>Value</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>t</td>
<td>Returns t when the current session is successfully closed.</td>
</tr>
<tr>
<td>nil</td>
<td>Returns nil if there is a problem closing the active session.</td>
</tr>
</tbody>
</table>

Example

ocnCloseSession()

=> t

Closes the current OCEAN session.
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

*t_filename*  
File in which to write the information. The ocnDisplay command opens the file, writes to the file, then closes the file. If you specify the filename without a path, the ocnDisplay 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.

*s_command*  
Command that was initially used to add the items that are now being displayed. Valid values: Most simulation setup commands. The commands that are supported include design, analysis, tran, ac, dc, noise, resultsDir, temp, option, desVar, path, includeFile, modelFile, stimulusFile, definitionFile, saveOption, envOption, save, converge, ic, forcenode, nodeset, simulator, setup, restore

*g_commandArg1*  
Argument corresponding to the specified command.
**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.

**Example**

- `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 'analyis )
  => t`
  Displays and writes all the analyses to the port named myPort.
ocnGetAdjustedPath

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

Description

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

Arguments

<table>
<thead>
<tr>
<th>Argument</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>t_libName</td>
<td>Library name of the top cellview of the design.</td>
</tr>
<tr>
<td>t_cellName</td>
<td>Cell name of the top cellview of the design.</td>
</tr>
<tr>
<td>t_viewName</td>
<td>View name of the top cellview of the design.</td>
</tr>
<tr>
<td>t_netName</td>
<td>A single concatenated string for the instance hierarchy with &quot;/&quot; as the hierarchy separator in the string.</td>
</tr>
</tbody>
</table>

Value Returned

<table>
<thead>
<tr>
<th>Value</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>t_adjustedPath</td>
<td>The reduced net name. If the net is local to this cell view only, the reduced net name is the same as the provided net name.</td>
</tr>
<tr>
<td>nil</td>
<td>Returns nil if there is a problem returning the adjusted path.</td>
</tr>
</tbody>
</table>

Example

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

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

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

**Description**

Turns off the specified information.

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

**Arguments**

- `s_command` Command that was initially used to add the items that are now being turned off.  
  Valid value: `restore`

- `g_commandArg1` Argument corresponding to the specified command.

- `g_commandArg2` Additional argument corresponding to the specified command.

**Value Returned**

- `t` Returns `t` if the information is turned off.

- `nil` Returns `nil` and prints an error message if there are problems turning off the information.

**Example**

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

Turns off the `restore` command.

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

Turns off the transient `restore` command.
option

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

Description

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

Arguments

s_categ Type of simulator to be used.
Valid values: analog if the options are for an analog simulator,
digital for a digital simulator, or mixed for a mixed-signal simulator
Default value: analog

s_option1 Name of the simulator option.

g_value1 Value for the option.

s_option2 Name of an additional simulator option.

g_value2 Value for the option.

Value Returned

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

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

Example

option( ‘abstol 1e-10 )

Sets the abstol option to 1e-10.

option( ‘delmax 50n )

Sets the delmax option to 50n.

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

```plaintext
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

s_analysisType Type of the analysis.
Valid values: dc or tran

t_filename Name of the file containing the saved state.

Value Returned

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

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

Example

restore( 'dc "./storeFile" )
=> ./storeFile

Initializes the simulator to the state saved in the storeFile file.

restore( 'tran "./tranStoreFile" )
=> ./tranStoreFile

Initializes the simulator to the state of a transient analysis saved in the tranStoreFile file.
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/spectre/schematic/psf") =>
                "~/simulation/ckt/spectre/schematic/psf"

  Specifies the psf directory as the directory in which to store the PSF files.
  resultsDir() => "~/simulation/ckt/spectre/schematic/psf"

  Returns the results directory.
run

run([?jobName t_jobName] [?drmsCmd t_drmsCmd])
    => s_jobName/nil

run([analysisList][?jobName t_jobName][?host t_hostName][?queue t_queueName][?startTime t_startTime][?termTime t_termTime][?dependentOn t_dependentOn][?mail t_mailingList][?block s_block][?notify s_notifyFlag][?lsfResourceStr s_lsfResourceStr])
    => s_jobName/nil

run( )
    => t_dirName/nil

run(s_analysisType1 - s_analysisTypeN)
    => t_dirName/nil

Description

Starts the simulation or specifies a time after which an analysis should start. If distributed processing is not available on the system or is not enabled, the arguments specific to distributed processing (see Arguments section below for list of arguments specific to distributed processing) are ignored and the simulation runs locally. If distributed processing is available and is enabled, the environment default values are used if not specified in the run command arguments. The environmental default values are stored in the .cdsenv file.

Do not use the run command to start the parametric analysis. Instead, use the command that is specific to the analysis.

<table>
<thead>
<tr>
<th>To start</th>
<th>Use this command</th>
</tr>
</thead>
<tbody>
<tr>
<td>parametric analyses</td>
<td>paramRun</td>
</tr>
</tbody>
</table>

Arguments

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

s_analysisType1 Name of a prespecified analysis to be simulated.

s_analysisTypeN Name of another prespecified analysis to be simulated.

The following arguments apply only when the distributed processing mode is enabled:
### OCEAN Reference
Simulation Commands

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

- **t_hostName**: Name of the host on which to run the analysis. If no host is specified, the system assigns the job to an available host.

- **t_queueName**: Name of the queue. If no queue is defined, the analysis is placed in the default queue.

- **t_startTime**: Desired start time for the job. If dependencies are specified, the job does not start until all dependencies are satisfied.

- **t_termTime**: Termination time for job. If the job has not completed by the specified termination time, the job is aborted.

- **t_dependentOn**: List of jobs on which the specified job is dependent. The job is not started until dependent jobs are completed.

- **t_mailingList**: List of users to be notified when the analysis is complete.

- **s_block**: When `s_block` is not set to `nil`, the OCEAN script halts until the job is complete.
  Default value: `nil`

- **s_notifyFlag**: When not set to `nil`, the job completion message is echoed to the OCEAN interactive window.
  Default value: `t`

- **s_lsfResourceStr**: Specifies an LSF Resource Requirement string to submit a job. It is effective only in the LSF mode.

- **t_drmsCmd**: A DRMS (Distributed Resource Management System) command, such as a `bsub` command for LSF or a `qsub` command for SGE (Sun Grid Engine) used to submit a job. When this argument is used, all other arguments, except `?jobName` will be ignored. Moreover, it will not be possible to call the OCEAN function `wait` on the jobs submitted using this argument.

To know more about the command option, refer to the section Submitting a Job in the chapter Using the Distributed Processing Option in the Analog Design Environment of
### Value Returned

<table>
<thead>
<tr>
<th>Function</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td><code>s_jobName</code></td>
<td>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.</td>
</tr>
<tr>
<td><code>t_dirName</code></td>
<td>Returns the name of the directory in which the results are stored. This value is returned for local and blocking distributed modes.</td>
</tr>
<tr>
<td><code>nil</code></td>
<td>Returns <code>nil</code> and prints an error message if there is an error in the simulation. In this case, look at the <code>yourSimulator.out</code> file for more information. (This file is typically located in the <code>psf</code> directory.)</td>
</tr>
</tbody>
</table>

### Example

```plaintext
run( ?jobName "job1" ?drmsCmd "bsub -q lnx32")
=> s_jobName/nil
```

where `lnx32` is the name of the queue to which the job is submitted.

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

Starts the simulation.

```plaintext
run('tran, 'ac)
```

Runs only the `tran` and `ac` analyses.

```plaintext
run('dc)
```

Runs only the `dc` analysis.

```plaintext
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.

```plaintext
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.

Submits the current design and enabled analyses as a jobName job1 on the fast queue host menaka with the job start time as 22:59 and termination time as 23:25. A mail will be sent to preampGroup after the job ends.

run( ?jobName "job1" ?queue "fast" ?host "menaka" ?lsfResourceStr "mem>500")

Submits the current design and enabled analyses as a jobName job1 on the fast queue host menaka, if the host has at least 500 MB of RAM memory.
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:

<table>
<thead>
<tr>
<th>Valid Values</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>v</td>
<td>Specifies that a list of subsequent net names be saved.</td>
</tr>
<tr>
<td>i</td>
<td>Specifies that a list of subsequent currents be saved.</td>
</tr>
<tr>
<td>all</td>
<td>Specifies that all nets and all currents are to be saved.</td>
</tr>
<tr>
<td>allv</td>
<td>Specifies that all voltages are to be saved.</td>
</tr>
<tr>
<td>alli</td>
<td>Specifies that all currents are to be saved.</td>
</tr>
</tbody>
</table>

Default value: allv

t_saveName1 Name of the net, device, or other object.

t_saveNameN Name of another net, device, or object.
Value Returned

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

*nil*  
Returns nil and prints an error message if there is a problem saving the outputs.

Example

```lisp
save( 'v "net34" "net45" )
```

Saves the outputs for net34 and net45.

```lisp
save( 'i "R1" "/Q1/b" )
```

Saves the currents for R1 and Q1/b.

```lisp
save( 'all )
```

Saves all the nets and currents.

```lisp
save( 'i "q1:b" "r1:p" "mn1:d" )
```

For the spectre simulator, saves the current through the specified devices.

```lisp
save( ?categ 'analog 'v "/vin" "/vout" )
```

Saves the output for vin and vout.

```lisp
save( 'i "i(q1,b)" "i(r1)" "i(mn1,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 *Virtuoso Spectre Circuit Simulator User Guide* for more information on these options.

**Note:** The `saveOption` command does not work with socket simulators. If you are using a socket simulator, you must instead specify save options with the `save` command described in “save” on page 135.

**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.)

\texttt{g\_optionValueN}

Value for the save option.

**Value Returned**

\texttt{undefined}

The return value for this command/function is undefined.

\texttt{nil}

Returns \texttt{nil} if there are problems specifying options.

**Example**

```plaintext
saveOption( 'save "lvl" 'nestlvl 10 'currents "selected" 'useprobes "yes"
 'subcktprobelvl 2 ?saveahdlvars "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.

**Example**

```
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.
solver

solver( s_solver )
   => s_solver/nil

Description

Sets a solver for a given AMS OCEAN session. The valid values for solver are Spectre and UltraSim. You select Spectre if you want to use an accurate AMS-Spectre analog engine. You select UltraSim if you want to use the AMS-Ultrasim or FastSPICE(UltraSim) solver for a given AMS simulation.

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

Arguments

s_solver Name of the solver.

Value Returned

s_solver Returns the name of the solver.
nil Returns nil and prints an error message if the specified solver is not registered with the Virtuoso® Analog Design Environment through OASIS. If the solver is not registered, the solver from the preceding session is retained.

Example

solver( 'spectre )
=> spectre

Specifies AMS-Spectre as the solver to be used for the current AMS session.

solver( 'ultraSim )
=> ultraSim

Specifies AMS-UltraSim (UltraSim FastSPICE) as the solver to be used for the current AMS session.
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

- t_fileName
  - The name of the stimulus file to be included.

- t_fileName2...t_fileNameN
  - The names of the additional stimulus files to be included.

- b_xlate
  - If set to t, net and instance expressions are translated to simulator names. The default value of the b_xlate variable is t.

Value Returned

- l_fileNames
  - A list of the stimulus file names is the output if the command is successful.

- nil
  - Otherwise nil is returned

Example

stimulusFile( "tran.stimulus rf.stimulus" ?xlate nil)
=> ("tran.stimulus rf.stimulus")

Includes tran.stimulus and rf.stimulus in the simulator input file. No net and instance expressions are translated.

stimulusFile()
=> ("tran.stimulus" "rf.stimulus")
Returns the stimulusFile, if one was set earlier. Otherwise, it returns nil.
store

store( s_analysisType t_filename )
    => t_filename/nil

Description

Requests that the simulator store its node voltages to a file.

You can restore this file in a subsequent simulation to help with convergence or to specify a certain starting point. This command is not available for the Spectre® circuit simulator, with which you can use the store/restore options: readns, readforce, write, or writefinal.

Note: store is available for the cdsSpice and hspiceS simulators.

Arguments

s_analysisType        Type of the analysis.
    Valid values: dc or tran

 t_filename           Name of the file in which to store the simulator’s node voltages.

Value Returned

 t_filename           Returns the filename.

 nil                   Returns nil and prints an error message if there are problems storing the information to a file.

Example

store( 'dc "./storeFile" )
    => ./storefile

Stores the simulator’s node voltages in a file named storeFile in the current directory.

store( 'tran "./tranStoreFile" )
    => ./transtorefile

Stores the node voltages for a transient analysis in a file named tranStoreFile in the netlist (design) directory unless a full path is specified.
**temp**

temp( *f_tempValue*)

=> *s_tempValue*/nil

**Description**

Specifies the circuit temperature.

**Arguments**

*f_tempValue*  
Temperature for the circuit.

**Value Returned**

*s_tempValue*  
Returns the temperature specified.

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

**Example**

temp( 125 )

=> ?125?

atof(temp( 125 ))

=> 125.0

Sets the circuit temperature to 125.

temp()

=> 125

Gets the value you had set for the circuit temperature. If you have not set a value for the temperature, it returns the default value.
**tran**

```plaintext
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 and spectreVerilog simulators.

**Arguments**

- `g_fromValue` Starting time for the analysis.
- `g_toValue` Ending time.
- `g_byValue` Increment at which to step through the analysis.

**Value Returned**

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

**Example**

```plaintext
tran( 1u)
=> “1e-06”
```

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

```plaintext
tran( 0 1u 1n )
=> “1e-09”
```

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

vcdFile( t_vcdFileName )

=> t_vcdFileName/nil

**Description**

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

**Note:** This command is applicable for AMS and UltraSim simulators. For AMS, it works only when UltraSim is the solver.

**Arguments**

- **t_vcdFileName**
  
  The name of the VCD file to be used for session.

**Value Returned**

- **t_vcdFileName**
  
  The VCD file name is the output if the command is successful.

- **nil**
  
  Otherwise, nil is returned.

**Example**

vcdFile("/tmp/vcdFile.dat")

=> "/tmp/vcdFile.dat"

Specifies /tmp/vcdFile.dat as the VCD file to be used for current AMS-UltraSim OCEAN session.
### vcdInfoFile

vcdInfoFile( t_vcdInfoFileName )

=> t_vcdInfoFileName/nil

**Description**

Sets a VCD info file for a given AMS or UltraSim OCEAN session when you have set UltraSim as the solver. You also need to specify a VCD file while using this command. You can specify only one VCD info file for a session. You may use ocnDisplay('vcdInfoFile) to view the currently active VCD info file.

**Note:** This command is applicable for AMS and UltraSim simulators. For AMS, it works only when UltraSim is the solver.

**Arguments**

- **t_vcdInfoFileName**  
The name of the VCD info file to be included.

**Value Returned**

- **t_vcdInfoFileName**  
The VCD info file name is the output if the command is successful.
- **nil**  
Otherwise, nil is returned.

**Example**

vcdInfoFile("/tmp/vcdInfoFile.dat")

=> " /tmp/vcdInfoFile.dat"

Specifies /tmp/vcdInfoFile.dat as the VCD file to be used for current AMS-UltraSim OCEAN session.
vecFile

vecFile( t_vecFile [t_vecFile1 ... t_vecFileN] )
=> t_vecFile(s)/nil

Description

Sets the vector files to be used in an AMS or UltraSim OCEAN session. You use the vecFile command to specify a list of vector files which go to control file. You may use ocnDisplay('vecFile) to view the currently active vector files in an OCEAN session.

Note: This command is applicable for AMS and UltraSim simulators. For AMS, it works only when UltraSim is the solver.

Arguments

_\texttt{t\_vecFile} \quad \text{The name of the vector file to be included.}

_\texttt{t\_vecFile1...t\_vecFileN} \quad \text{The names of the additional vector files to be included.}

Value Returned

_\texttt{t\_vecFile} \quad \text{The names of the vector file(s) are listed if the command is successful.}

_\texttt{nil} \quad \text{Otherwise, nil is returned.}

Example

vecFile("/tmp/vec.dat" "/tmp/vec2.dat")
=> ("/tmp/vec1.dat" "/tmp/vec2.dat")

Specifies /tmp/vec.dat and /tmp/vec.dat2 as the vector files to be used for the current AMS-UltraSim OCEAN session.
hlcheck

hlcheck( t_value )
   => t / nil

Description

Sets or gets the value of the hlcheck option used in the vec_include statement in a netlist. You may use the ocnDisplay('hlcheck) command to view the current value of hlcheck in an OCEAN session associated with vector files.

Note: This command is applicable only when one or more vector files are specified in a given 'spectre' OCEAN session.

Arguments

   t_value       Value to be set for the hlcheck option. Possible values include "off", "0", and "1". The value "off" disables the hlcheck option in the vec_include statement.

Value Returned

   t          Returns t if the hlcheck option is set with the value supplied as argument

   nil       Otherwise, returns nil and an error message is displayed

Example

hlcheck( "1" )
   => t

Sets the value of the hlcheck option as 1 in the vec_include statement

hlcheck()
   => "1"

Returns the value of the hlcheck option
Data Access Commands

The data access commands let you open results and select different types of results to analyze. You can get the names and values of signals and components in the selected results, and you can print different types of reports.

In this chapter, you can find information on the following data access commands:

dataTypes on page 153
getData on page 154
getResult on page 156
i on page 157
ocnHelp on page 159
ocnResetResults on page 161
openResults on page 162
outputParams on page 164
outputs on page 166
phaseNoise on page 168
pv on page 170
resultParam on page 172
results on page 174
selectResult on page 175
sp on page 177
sweepNames on page 179
sweepValues on page 181
sweepVarValues on page 182
v on page 184
vswr on page 186
zm on page 188
zref on page 190
**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

- `l_dataTypes` Returns the list of data types.
- `nil` Returns `nil` and an error message if the list of datatypes cannot be returned.

### Example

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

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

ggetData( 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

t_name Name of the signal.

s_resultName Results from an analysis. When specified, this argument will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult command.

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

Value Returned

x_number Returns an integer simulation result.

o_waveform Returns a waveform object. A waveform object represents simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave:XXXXX.)

nil Returns nil and an error message if the value cannot be returned.

Example

ggetData( "/net6" ) => srrWave:25178234
OCEAN Reference
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 )
=> srrWave:96879364
```

Returns the number or waveform for V1. In this example, the return value is equivalent to:
\( i( "/V1" ?result 'ac ) \).

```
selectResult( 'tran ) =>
ocnPrint( getData( "net1" ) ) =>
```

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".
getResult

getResult ( [?result s_resultName [%resultsDir t_resultsDir]] )  
  => o_results/nil

Description

Gets the data object for a specified analysis without overriding the status of any previously executed selectResult() or openResults() commands.

Returns the data object for a particular analysis similar to the selectResult() function does. Unlike the selectResult() function, all subsequent data access commands will not internally use this information.

Arguments

s_resultName  Results from an analysis. When specified, this argument will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult command.

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

Value Returned

o_results  Returns the object representing the selected results.

nil  Returns nil and an error message if there are problems accessing the analysis.

Example

getResult( ?result 'tran )
i

i(  t_component [?result s_resultName [?resultsDir t_resultsDir]])
   => o_waveform/nil

Description

Returns the current through the specified component.

Arguments

t_component  Name of the component.

s_resultName  Results from an analysis. When specified, this argument will only
be used internally and will not alter the current result which was
set by the selectResult command. The default is the current
result selected with the selectResult command.

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

Value Returned

o_waveform  Returns a waveform object. A waveform object represents
simulation results that can be displayed as a series of points on
a grid. (A waveform object identifier looks like this:
srrWave:XXXXX.).

nil  Returns an error message and nil if there is a problem.

Example

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

Returns the current through the R1 component.

ocnPrint( i( "/R1" ) )
Prints the current through the R1 component.
\[ \text{ocnPrint( i( "/R1" ?result 'dc ) )} \]

Prints the current through the R1 component with respect to the dc swept component.
\[ \text{ocnPrint( i( "/R1" ?resultsDir "./test2/psf" ?result 'dc ) )} \]

Prints the current through the R1 component with respect to dc for the results from a different run (stored in test2/psf).
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

- **t_filename**
  File in which to write the information. The ocnHelp command opens the file, writes to the file, and closes the file. If you specify the filename without a path, the ocnHelp command creates the file in the directory pointed to by your Skill Path. To find out what your Skill path is, type `getSkillPath()` at the OCEAN prompt.

- **p_port**
  Port (previously opened with outfile) through which to append the information to a file. You are responsible for closing the port. See the outfile command for more information.

- **s_command**
  Command for which you want help.

Value Returned

- **t**
  Displays the online help and returns t.

- **nil**
  Returns nil and an error message if help cannot be displayed.

Example

ocnHelp()
=> t

Displays information about using online help.
ocnHelp('analysis')
=> t

Displays help for the analysis command.

ocnHelp( ?output "helpInfo" )
=> t

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

ocnResetResults()
=> t

Description
Unsets the results opened by the openResults command. Use this command to return to the state that existed prior to using the openResults command.

Arguments
None.

Value Returned
t
   Resets the results and returns t.

Example

getResult( ?result 'tran )  Returns nil when no results have been opened.
openResults( "./psf" )      Makes getResult return valid object.
ocnResetResults()          Resets the results opened by openResults and makes getResult return nil.
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 a psf directory and must contain `runObjFile`. The default value for this argument is `t`. 
Value Returned

\textit{t\_dirName} \hspace{1cm} The directory containing the PSF files.

\texttt{nil} \hspace{1cm} Returns \texttt{nil} and an error message if there are problems opening the results.

Example

\begin{verbatim}
openResults( "./simulation/opamp/spectre/schematic/psf" )
=> "./simulation/opamp/spectre/schematic/psf"
\end{verbatim}

Opens the results in the \texttt{psf} directory within the specified path.

\begin{verbatim}
openResults( "./psf" )
=> psf
\end{verbatim}

Opens the results in the \texttt{psf} directory in the current working directory.

\begin{verbatim}
openResults( "./psf" t )
=> psf
\end{verbatim}

Opens the results in the \texttt{psf} directory in the current working directory. It also allows the evaluation of the Calculator expression.
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

- **t_compType**
  - Name of a component.

- **s_resultName**
  - Results from an analysis. When specified, this argument will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult command.

- **t_resultsDir**
  - Directory containing the PSF files (results). If you supply this argument, you must also supply the resultName argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the openResults command.

Value Returned

- **l_outputParams**
  - Returns the list of parameters.

- **nil**
  - Returns nil and an error message if there are no associated parameters or if the specified component (compType) does not exist.
Example

```
selectResult( '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" )
Returns all the signal names that are node voltages. The dataType (signal) returns the data type of the signal.

```lisp
outputs(?result "tran" ?resultsDir "./psf")
=> ("net11" "net15" "net17")
```

Returns the names of the outputs for the tran results stored at the location "./psf."
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.
Example

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

- **t_name**
  - Name of the node or component.

- **t_param**
  - Name of the parameter.

- **s_resultName**
  - Results from an analysis. When specified, this argument will only be used internally and will not alter the current result which was set by the `selectResult` command. The default is the current result selected with the `selectResult` command.

- **t_resultsDir**
  - Directory containing the PSF files (results). If you supply this argument, you must also supply the `resultName` argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the `openResults` command. The default is the current results directory set by the `openResults` command.

**Value Returned**

- **g_value**
  - Returns the requested parameter value.

- **nil**
  - Returns `nil` and prints an error message.

**Example**

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

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

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

\texttt{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.
resultParam

resultParam( S_propertyName [?result s_resultName [?resultsDir t_resultsDir]] ) => L_value/nil

Description

Returns the value of a header property from the selected result data.

Arguments

s_propertyName
Name of the parameter

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
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

L_value
Value of the parameter. The data type depends on the data type of the parameter.

nil
Returns nil and an error message if there are problems returning the results.

Example

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

Returns the name of the positive and negative output signals from PSS-noise analysis result. In this case, the data type of the returned value is a string.

resultParam("port1.r.value" ?result "sp")
=> 40.0
resultParam("port2.r.value" ?result "sp")
=> 40.0

Returns the reference impedance of the ports in a two-port network from the S-parameter analysis result. In this case, the data type of the returned value is a floating point number.

resultParam("positive output signal" ?result "pnoise.pss" ?resultsDir "./psf")
=> "0"

Returns the names of the positive output signals from the PSS-noise analysis results stored at the location ./psf.
results
results( [ ?resultsDir t_resultsDir ] )
   => l_results/nil

Description

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

Arguments

t_resultsDir

Directory containing the PSF files (results). When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the openResults command.

Value Returned

l_results
Returns the list of result types.

nil
Returns nil and an error message if there are problems returning the results.

Example

results()
=> ( dc tran ac )

Returns the list of results available.

results( ?resultsDir "./psf" )

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

s_resultsName  Results from an analysis.
n_sweepValue   The sweep value you wish to select for an analysis.

Value Returned

o_results      Returns the object representing the selected results.
nil            Returns nil and an error message if there are problems selecting the analysis.

Example

selectResult( 'tran')

Selects the results for a transient analysis.
sweepValues(3.0 3.333333 3.666667 4.0 4.333333 4.666667 5.0)
selectResult("tran" "3.333333")

The sweepValues command prints a list of sweep values.

The selectResult command selects a specific value for a transient analysis.

selectResult( 'tran')

Selects the results for a transient analysis.

paramAnalysis("supply" ?start 3 ?stop 5 ?step 1.0/3)
paramRun("supply")
selectResult(( ’tran car sweepValues() )

Selects the data corresponding to the first parametric run.

**Note:** `selectResult(’tran)` would select the entire family of parametric data.
sp

sp( x_iIndex x_jIndex [?result s_resultName [?resultsDir t_resultsDir]] ) => o_waveform/nil

Description

Returns S-parameters for N port networks.

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

Arguments

x_iIndex
The \textit{i}th index of the coefficient in the scattering matrix.

x_jIndex
The \textit{j}th index of the coefficient in the scattering matrix.

s_resultName
Results from an analysis. When specified, this argument will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult command.

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

Value Returned

o_waveform
Waveform object representing the S-parameter.

nil
Returns \textit{nil} if there is an error.

Example

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

s11 = sp(1 1 ?result "sp" ?resultsDir "./simResult/psf")
Returns the S-parameter s11 for results of S-parameter(sp) analysis stored at the location .simResult/psf.
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

o_waveForm

Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave:XXXXX). When this argument is used, the t_resultsDir and s_resultName arguments are ignored.

s_resultName

Results from an analysis. When specified, this argument will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult command.

t_resultsDir

Directory containing the PSF files (results). If you supply this argument, you must also supply the resultName argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the openResults command.

Value Returned

l_sweepName

Returns a list of the sweep names.

nil

Returns nil and prints an error message if the sweep names cannot be returned.

Example

selectResult('tran)
sweepNames()
=> ( "TEMPDC" "time" )
Returns a list of sweep variables for the selected results. In this case, the return values indicate that the data was swept over temperature and time.

```lisp
sweepNames(?result 'ac)
=> ("TEMPDC" "freq")
sweepNames()
=> ("TEMPDC" "time")
w = VT("/vout")
sweepNames( w )
=> ( "r" "time")
```

Returns the sweep variables for the waveform w.

```lisp
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

o_waveForm Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave:XXXXX.)

Value Returned

l_sweepValues Returns the list of sweep values.

nil Returns nil and an error message if the list of sweep values cannot be returned.

Example

sweepValues()
=> ( -50 -15 20 55 90.0 )

Returns a list of sweep values for the selected results. In this case, the return values indicate the temperature over which the data was swept.

w = VT("/vout")

sweepNames( w )
=> ( "r" "time" )

sweepValues( w )
=> ( 2000 4000 6000 )

Returns a list of sweep values for the wave 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

t_varName        Name of the specific variable from which the values are retrieved.

s_resultName     Results from an analysis. When specified, this argument will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult command.

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

Value Returned

l_sweepValues    Returns the list of sweep values.

nil              Returns nil and an error message if the list of sweep values cannot be returned.

Example

selectResult('tran)
sweepNames()
=> ("TEMPDC" "Vs supply" "time")
sweepVarValues("TEMPDC")
=> (0 32)
sweepNames(?result 'ac)
=> ("TEMPDC" "Vsupply" "freq")
sweepVarValues("Vsupply" ?result 'ac)
=> (5 12 15)
sweepNames(?result 'ac ?resultsDir "./simResult/psf")
=> ("TEMPDC" "freq")
sweepVarValues("TEMPDC" ?result 'ac ?resultsDir "./simResult/psf")
=> (-15 20 55)
v

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

Description

Returns the voltage of the specified net.

Arguments

t_net Name of the net.
s_resultName Results from an analysis. When specified, this argument will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult command.
t_resultsDir Directory containing the PSF files (results). If you supply this argument, you must also supply the resultName argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the openResults command.

Value Returned

o_waveform Returns a waveform object. A waveform object represents simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave:XXXXX.).

nil Returns an error message and nil if there is a problem.

Example

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

Returns the voltage for net56.

ocnPrint( v( "/net56" ) )
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`).
**vswr**

\[
vswr( x\_index \ [?result \ s\_resultName \ [?resultsDir \ t\_resultsDir]] ) \implies \text{o\_waveform/nil}
\]

**Description**

Computes the voltage standing wave ratio.

This function is a higher level wrapper for the OCEAN expression

\[
(1 + \text{mag}(s(x\_index \ x\_index ))) / (1 - \text{mag}(s(x\_index \ x\_index )))
\]

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

**Arguments**

- **x\_index**: Index of the port.
- **s\_resultName**: Results from an analysis. When specified, this argument will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult command.
- **t\_resultsDir**: Directory containing the PSF files (results). If you supply this argument, you must also supply the resultName argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the openResults command.

**Value Returned**

- **o\_waveform**: Waveform object representing the voltage standing wave ratio.
- **nil**: Returns an error message or nil if there is a problem.

**Example**

```
plot(vswr(2))
vswr1 = vswr(1 ?result "sp" ?resultsDir "./simResult/psf")
```
Returns the voltage standing wave ratio value at port 1 for the results of S-parameter(sp) analysis stored at the location ./simResult/psf.
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

\[
\frac{1 + s(x_{\text{index}} x_{\text{index}})}{1 - s(x_{\text{index}} x_{\text{index}})} \times \text{or}(\text{zref}(x_{\text{index}}) 50)
\]

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

Arguments

x_index

Index of the port.

s_resultName

Results from an analysis. When specified, this argument will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult command.

t_resultsDir

Directory containing the PSF files (results). If you supply this argument, you must also supply the resultName argument. When specified, this argument will only be used internally and will not alter the current results directory which was set by the openResults command. The default is the current results directory set by the openResults command.

Value Returned

o_waveform

Waveform object representing the port input impedance.

nil

Returns an error message and nil if there is a problem.

Example

plot(zm(2))
zm1 = zm(1 ?result "sp" ?resultsDir "./simResult/psf")
Returns input impedance at port 1 for results of S-parameter (sp) analysis stored at the location ./simResult/psf.
zref

zref( x_portIndex [?result s_resultName [?resultsDir t_resultsDir]] ) => f_impedance/nil

Description

Returns the reference impedance for an N-port network.

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

Arguments

x_portIndex  Index of the port.

s_resultName  Results from an analysis. When specified, this argument will only be used internally and will not alter the current result which was set by the selectResult command. The default is the current result selected with the selectResult command.

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

Value Returned

f_impedance  Reference impedance.
nil  Returns an error message and nil if there is a problem.

Example

Zref = zref(2)
zref1 = zref(1 ?result "sp" ?resultsDir "./simResult/psf")

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

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

- addSubwindow
- addSubwindowTitle
- addTitle
- addWaveLabel
- addWindowLabel
- clearAll
- clearSubwindow
- currentSubwindow
- currentWindow
- dbCompressionPlot
- dcmatchSummary
- deleteSubwindow
- deleteWaveform
- displayMode
- getAsciiWave
- graphicsOff
- graphicsOn
- hardCopy
- hardCopyOptions
- ip3Plot
This chapter also includes a topic, Plotting and Printing SpectreRF Functions in OCEAN.
**addSubwindow**

`addSubwindow()`

`=> x_subwindowID/nil`

**Description**

Adds a subwindow to the current Waveform window and returns the number for the new subwindow, which is found in the upper right corner.

**Arguments**

None.

**Value Returned**

- `x_subwindowID` Returns the window ID of the new subwindow.
- `nil` Returns `nil` and an error message if there is no current Waveform window.

**Example**

`addSubwindow() => 3`

Adds a new subwindow to the Waveform window.
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.
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.
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**

- **x_waveIndex** Integer identifying the waveform curve.
- **l_location** List of two waveform coordinates that describe the location for the label.
- **t_label** Label for the waveform.
- **l_textOffset** An offset of the label from `l_location`, in screen units of the current subwindow. If `l_textOffset` is not specified, it defaults to `0:0` and the label is displayed at the location. If `l_textOffset` 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.
- **x_color** Label color specified as an index in the technology file. Default value: `10`
- **t_justify** Justification, which is specified as "upperLeft", "centerLeft", "lowerLeft", "upperCenter", "centerCenter", "lowerCenter", "upperRight", "centerRight", or "lowerRight". Default value: "lowerLeft"
- **t_fontStyle** Font style, which is specified as "euroStyle", "gothic", "math", "roman", "script", "stick", "fixed"
"swedish", "raster", or "milSpec".
Default value: the font style of the current subwindow

\textit{\texttt{x\_height}}
Height of the font.
Default value: the font height of the current subwindow

\textit{\texttt{t\_orient}}
Orientation of the text, specified as either "R0" or "R90".
Default value: "R0"

\textit{\texttt{g\_drafting}}
Boolean that specifies whether the label stays backwards or upside-down. If set to \texttt{t}, a backwards or upside-down label is displayed in a readable form. If set to \texttt{nil}, a backwards or upside-down label stays the way it is.
Default value: \texttt{t}

\textit{\texttt{g\_overbar}}
Boolean that specifies whether underscores in labels are displayed as overbars. If set to \texttt{t}, underscores in labels are displayed as overbars. If set to \texttt{nil}, underbars are displayed as underbars.
Default value: \texttt{nil}

\textbf{Value Returned}

\textit{\texttt{s\_labelId}}
Returns an identification number for the waveform label.

\textit{\texttt{nil}}
Returns \texttt{nil} if there is an error.

\textbf{Example}

\texttt{addWaveLabel( 1 list( 0 0.5 ) "R5 = " )}

Attaches the "R5 = " label to the specified coordinates on waveform curve 1.

\texttt{addWaveLabel( 2 list( 0 0.5 ) "R_6 = " ?textOffset 0:20 ?justify "lowerCenter" ?fontStyle "roman" ?height 10 ?orient "R20" ?drafting \texttt{t} ?overbar \texttt{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**

```plaintext
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**

- `l_location`: List of two waveform coordinates that describe the location for the label.
  - Valid values: 0 through 1
- `t_label`: Label for the waveform.

**Value Returned**

- `s_labelId`: Returns an identification number for the subwindow label.
- `nil`: Returns `nil` if there is an error.

**Example**

```plaintext
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

\begin{itemize}
\item \texttt{t} \quad \text{Returns \texttt{t} if the contents of the subwindow are erased.}
\item \texttt{nil} \quad \text{Returns \texttt{nil} and an error message otherwise.}
\end{itemize}

Example

clearSubwindow()
clearSubwindow()
   => t

Erases the contents of the current subwindow.
currentSubwindow

currentSubwindow( x_subwindow )
    => t / nil

Description

Specifies \textit{x\_subwindow} as the current subwindow.

Arguments

\textit{x\_subwindow} \hspace{1cm} Number of the subwindow, found in the upper right corner, that is to become the current subwindow.

Value Returned

\textit{t} \hspace{1cm} Returns \textit{t} when the subwindow is set to \textit{x\_subwindow}.

\textit{nil} \hspace{1cm} Returns \textit{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.
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

  o_wave
    The waveform for which to plot the compression.

  x_harmonic
    Harmonic frequency index.

  x_extrapolationPoint
    The extrapolation point.

  x_compression
    The amount of dB compression.
    Default value: 1

Value Returned

  t
    Returns t if the point is plotted

  nil
    returns nil 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").
**dcmatchSummary**

```scad
dcmatchSummary([?resultsDir t_resultsDir] [?resultName 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 ls_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**

- **t_resultsDir**
  The directory containing the dcmatch-analysis results.

- **S_resultName**
  Results from an analysis for which you want to print the dcmatchSummary report.

- **t_fileName**
  File in which to write the information. The `dcmatchSummary` command opens the file, writes to the file and closes the file. If you specify the filename without a path, the `dcmatchSummary` command creates the file in the directory pointed to by your Skill Path. To find out what your Skill path is, use `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.

- **ln_paramValues**
  List of values for swept parameters at which the `dcmatchSummary` is to be printed. In case there is just one swept parameter the value can be specified as is.
ls_deviceType

List of device type strings to be included. Valid values are a list of strings or ‘all’ or a single device name. Default value is ‘all’.

ls_variations

An association list containing the device name and the associated variations to print. You can also specify the value ‘all’ to print all available variations for a device. Default value is ‘all’. For Example: ‘( (“bsim3v3” (“sigmaOut” “sigmaVth”)) (“resistor” (“sigmaOut”))

lt_includeInst

List of instance name strings to definitely include in the dcmatchSummary.

lt_excludeInst

List of instance name strings to exclude in the dcmatchSummary.

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

<table>
<thead>
<tr>
<th>Valid Values</th>
<th>Description</th>
<th>Sample Values for truncateData</th>
</tr>
</thead>
<tbody>
<tr>
<td>'top'</td>
<td>Saves information for the number of components specified with truncateData. The components with the highest contributions are saved.</td>
<td>10</td>
</tr>
<tr>
<td>'relative'</td>
<td>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</td>
<td>1.9n</td>
</tr>
<tr>
<td>'absolute'</td>
<td>Saves information for all the components in the selected set whose contribution are more than truncateData.</td>
<td>0.1</td>
</tr>
</tbody>
</table>
Ocean Reference
Plotting and Printing Commands

<table>
<thead>
<tr>
<th>ls_sortType</th>
<th>Specifies how the printed results are to be sorted. The valid values are nil, 'name, 'output.</th>
</tr>
</thead>
<tbody>
<tr>
<td>Value Returned</td>
<td>Returns the name of the port.</td>
</tr>
<tr>
<td>t_fileName</td>
<td>Returns nil and an error message if the summary cannot be printed.</td>
</tr>
</tbody>
</table>

Example

dcmatchSummary( ?result 'dcmatch-mine )

Prints a report for non-swept DC-Mismatch analysis.
dcmatchSummary( ?resultsDir "/usr/simulation/lowpass/spectre/schematic" ?result 'dcmatch )

Prints a report for non-swept DC-Mismatch analysis for the results from a different run (stored in the schematic directory).

Prints a report for swept DC-Mismatch analysis at swept parameter value of 25.
dcmatchSummary( ?result dcmatch-mine ?output "../summary.out")

Prints a report for non-swept DC-Mismatch analysis in the output file summary.out.
dcmatchSummary( ?paramValues 25 ?deviceType "bsim3v3" ?variations "(("bsim3v3" ("sigmaOut "sigmaVth" )))")

Prints a report for swept DC-Mismatch analysis at swept parameter value of 25 for bsim3v3 deviceType and sigmaOut and sigmaVth variations.
dcmatchSummary( ?paramValues 25 ?truncateType 'top ?truncateData 1)

Prints a report for swept DC-Mismatch analysis at swept parameter value of 25 printing only the component having the highest contribution.
dcmatchSummary( ?paramValues 25 ?sortType 'name )
Prints a report for swept DC-Mismatch analysis at swept parameter value of 25 sorted on name.
deleteSubwindow

deleSubwindow()
=> t / nil

Description

Deletes the current subwindow from the current Waveform window.

Arguments

None.

Value Returned

| t | Returns t if the current subwindow is deleted. |
| nil | Returns nil and an error message if there is no current subwindow. |

Example

deleSubwindow()
=> t

Deletes the current subwindow from the Waveform window.
**deleteWaveform**

```c
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**

- `x_index` Integer identifying a particular waveform curve.
- `all_string` The string "all" specifying that all waveform curves are to be deleted.

**Value Returned**

- `t` Returns `t` if the curves are deleted.
- `nil` Returns `nil` and an error message if the curves are not deleted.

**Example**

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

Deletes waveform 1 from the current subwindow.

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

Deletes all the curves from the current subwindow.
**displayMode**

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

**Description**

Sets the display mode of the current subwindow.

**Arguments**

- `t_mode`  
  String representing the display mode for the subwindow.  
  Valid values: `strip`, `smith`, or `composite`

**Value Returned**

- `t`  
  Returns `t` when the display mode of the subwindow is set.

- `nil`  
  Returns `nil` and an error message if the display mode cannot be set.

**Example**

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

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

  t_filename
    The name of the Ascii file to be read in.

  x_xColumn
    The column in the data file that contains the X-axis data.

  x_yColumn
    The column in the data file that contains the Y-axis data.

  x_xskip
    The number of lines to skip in the X column.

  x_yskip
    The number of lines to skip in the Y column.

Value Returned

  o_wave
    The DRL waveform object

  nil
    Returns nil if the function fails.

Example

getAsciiWave("~/mydatafile.txt " 1 2 )
 => srrWave:32538648

Reads in an ascii file ~/mydatafile.txt, which has x-axis data in the first column and y-axis data in the second column, and returns a DRL waveform object.

getAsciiWave("~/mydatafile.txt " 1 2 ?xskip 1 ?yskip 2)
 => srrWave:32538656

Reads in an ascii file ~/mydatafile.txt, which has x-axis data in the first column and y-axis data in the second column and skips 1 line in the xColumn and 2 lines in the yColumn, and returns a DRL waveform object.
**graphicsOff**

```lisp
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**

- **t**
  - Returns `t` if redrawing is disabled.

- **nil**
  - Returns `nil` if there is an error, such as there is no current Waveform window.

**Example**

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

Disables the redrawing of the Waveform window.
graphicsOn

```lisp
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**

```lisp
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 or a file. To plot to a printer specify a printer name using the \( \texttt{hcPrinterName} \) argument of the \texttt{hardCopyOptions} command. To plot to a file, specify a file name using the \( \texttt{hcOutputFile} \) argument of the \texttt{hardCopyOptions} command.

Note: You must first set any plotting options with the \texttt{hardCopyOptions} command.

Arguments

\( w\_\text{windowId} \)
The window ID of the waveform window whose plot is to be sent to a printer or a file. The default value is the window ID of the current window.

Value Returned

\( t \)
Returns \( t \) if successful.

\( \text{nil} \)
Returns \( \text{nil} \) if there is an error.

Example

\begin{verbatim}
hardCopy()
=> t
\end{verbatim}

Sends a waveform plot to the printer or to a file.

\begin{verbatim}
w = newWindow()
plot(v("/vout"))
hardCopy(w)
\end{verbatim}

Sends the waveform plot of \( w \) to the printer or to a file.
hardCopyOptions

hardCopyOptions( [?hcCopyNum x_hcCopyNum] [?hcOffsetHeight x_hcOffsetHeight] [?hcOffsetWidth x_hcOffsetWidth] [?hcOrientation s_hcOrientation] [?hcOutputFile g_hcOutputFile] [?hcPrinterName s_hcPrinterName] [?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

x_hcCopyNum The number of copies to plot.
Valid values: any positive integer
Default value: 1

x_hcOffsetHeight The vertical margin.
Valid values: any positive integer
Default value: 0

x_hcOffsetWidth The horizontal margin.
Valid values: any positive integer
Default value: 0

s_hcOrientation The plot orientation.
Valid values: ‘portrait,’ ‘landscape,’ ‘automatic
Default value: ‘automatic
**Ocean Reference**  
Plotting and Printing Commands

---

### g_hcOutputFile
Name of the output file. The output file can be created in one of the following file formats:
- BMP – Windows Device Independent Bitmap (.bmp)
- PNG – Portable Network Graphics (.png)
- PS – PostScript (.ps)
- TIFF – Tagged Image File Format (.tif)

Valid values: a string or nil
Default value: nil

### s_hcPrinterName
The name of the printer.
Valid values: a string or nil
Default value: nil

### 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.

---

**Example**

hardCopyOptions( ?hcNumCopy 1 )

Plots one copy of the window or subwindow.

hardCopyOptions(?hcNumCopy 3 ?hcOutputFile "myOutFile.bmp")

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

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

Plots 2 copies of the window in portrait orientation and sends them to the file myOutFile.png.
ip3Plot

```lisp
ip3Plot( o_wave x_sigHarmonic x_refHarmonic x_extrapolationPoint )
=> t / nil
```

**Description**

Plots the IP3 curves.

This command should be run on the results of the Spectre swept pss and pac analysis.

Refer to the “Simulating Mixers” chapter of the *Virtuoso Spectre Circuit Simulator RF Analysis User Guide* for more information on ip3Plot.

**Arguments**

- **o_wave**: Waveform for which to plot the ip3.
- **x_sigHarmonic**: Index of the third order harmonic.
- **x_refHarmonic**: Index of the first order (fundamental) harmonic.
- **x_extrapolationPoint**: Extrapolation point.

**Value Returned**

- **t**: Returns `t` if the curves are plotted.
- **nil**: Returns `nil` if there is an error.

**Example**

```lisp
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

<table>
<thead>
<tr>
<th>w_windowId</th>
<th>Returns the window ID of the new Waveform window.</th>
</tr>
</thead>
<tbody>
<tr>
<td>nil</td>
<td>Returns nil and an error message if the new Waveform window cannot be created.</td>
</tr>
</tbody>
</table>

Example

newWindow()

=> window:3

Creates a new Waveform window that is numbered 3 in the upper right corner.
noiseSummary

noiseSummary(s_type [?result s_resultName [?resultsDir t_resultsDir]]
       [?frequency f_frequency] [?weight f_weight] [?output t_fileName | p_port]
       [?noiseUnit t_noiseUnit] [?truncateData x_truncateData]
       [?truncateType s_truncateType] [?digits x_digits]
       [?percentDecimals x_percentDecimals] [?from f_from] [?to f_to]
       [?deviceType ls_deviceType] [?weightFile t_weightFile]
       [?paramValues ls_paramValues])
=> 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

s_type
Type of noise-analysis results for which to print the report.
Valid values: spot, to specify noise at a particular frequency, or
integrated, to specify noise integrated over a frequency range.

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.

f_frequency
Frequency value of interest.

f_weight
Waveform representing the function with which the integral is
weighted.
Default value: 1.0

File in which to write the information. The noiseSummary
command opens the file, writes to the file, and closes the file. If
you specify the filename without a path, the `noiseSummary` command creates the file in the directory pointed to by your Skill Path. To find out what your Skill path is, type `getSkillPath()` at the OCEAN prompt.

**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 V^2/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.

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

**x_digits**  
Number of significant digits with which the contributors are printed.

**x_percentDecimals**  
Number of decimals printed for any relative contribution.
Ocean Reference
Plotting and Printing Commands

\textit{f_{from}} \quad \text{For integrated noise, the start value for frequency.}

\textit{f_{to}} \quad \text{For integrated noise, the end value for frequency.}

\textit{ls\_deviceType} \quad \text{List of device type strings to be included.}
\text{Valid values: a list of strings or ‘all}

\textit{t\_weightFile} \quad \text{Absolute or relative path of the file that contains information about weights. This data is used to compute weighted noise. If the values are provided for both parameters, weight and weightFile, the value for weight gets precedence.}

\textit{ls\_paramValues} \quad \text{List of values where each value co-relates to a specific sweep variable name. This field must be used when the data is parametric. The order of this list must coincide with the list returned by the \texttt{sweepNames} function excluding the frequency variable.}

\textbf{Value Returned}

\textit{t\_fileName} \quad \text{Returns the name of the port.}

\textit{p\_port} \quad \text{Returns the name of the file.}

\textit{nil} \quad \text{Returns \texttt{nil} and an error message if the summary cannot be printed.}

\textbf{Example}

\begin{verbatim}
noiseSummary( ‘integrated ?result ‘noiseSweep-noise )
\end{verbatim}

Prints a report for an integrated noise analysis.

\begin{verbatim}
noiseSummary( ‘integrated ?resultsDir 
"/usr/simulation/lowpass/spectre/schematic
 ?result ‘noise)
\end{verbatim}

Prints a report for an integrated noise analysis for the results from a different run (stored in the \texttt{schematic} directory).

\begin{verbatim}
noiseSummary( ‘spot ?resultsDir 
"/usr/simulation/lowpass/spectre/schematic
 ?result ‘noise ?frequency 100M )
\end{verbatim}

Prints a report for a spot noise analysis at a frequency of 100M.
noiseSummary('integrated ?truncateType 'none ?digits 10
?weightFile "./weights.dat")

Prints the weighted noise for an integrated noise analysis using information in the weight
file weights.dat.

noiseSummary('integrated ?output "./NoiseSum1" ?noiseUnit "V" ?truncateData 20
?truncateType 'top ?from 10 ?to 10M ?deviceType list("bjt" "mos" "resistor")

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

noiseSummary( 'integrated ?from 1 ?to 100M ?truncateType 'top ?truncateData 20
?deviceType 'all ?noiseUnit "V^2" ?output "./filename.ns" ?paramValues list(2.47e-9))

Prints a report for an integrated noise analysis at a specific swept value.
ocnPrint

ocnPrint( [?output t_filename | p_port] [?precision x_precision]
[?numberNotation s_numberNotation] [?numSpaces x_numSpaces] [?width
x_width] [?from x_from] [?to x_to] [?step x_step] [?linLog t_linLog]
o_waveform1 [o_waveform2 ...] )
=> t / nil

Description

Prints the text data of the waveforms specified in the list of waveforms.

If you provide a filename as the ?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

t_filename
File in which to write the information. The ocnPrint command opens the file, writes to the file, and closes the file. If you specify the filename without a path, the OCEAN environment creates the file in the directory pointed to by your Skill Path. To find out what your Skill path is, type getSkillPath() at the OCEAN prompt.

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.

x_precision
The number of significant digits to print. This value overrides any global precision value set with the setup command.
Valid values: 1 through 16
Default value: 6
Note: To print the specified significant number of digits, ensure that the value of the x_width argument is the same or greater than the value of the x_precision argument.

s_numberNotation
The notation for printed information. This value overrides any global format value set with the setup command.
Valid values: ‘suffix,’ ‘engineering,’ ‘scientific,’ ‘none
Default value: ‘suffix
The format for each value is ‘suffix: 1m, 1u, 1n, etc.;’ ‘engineering: 1e-3, 1e-6, 1e-9, etc.;’ ‘scientific: 1.0e-2, 1.768e-5, etc.;’ ‘none.’

The value ‘none’ is provided so that you can turn off formatting and therefore greatly speed up printing for large data files. For the fastest printing, use the ‘none’ value and set the ?output argument to a filename or a port, so that output does not go to the CIW.

`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.

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

`t_linLog` The scale to be used for printing.
Valid values: Linear, Log
Default value: Linear

`o_waveform1` Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave:XXXXX.)

`o_waveform2` Additional waveform object.

**Value Returned**

`t` Returns t if the text for the waveforms is printed.

`nil` Returns nil and an error message if the text for the waveforms cannot be printed.
Example

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

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

```
ocnPrint( vm( "/net56" ) vp( "/net56" ) )
=> t

Prints the text for the waveforms for the magnitude of the voltage of net56 and the phase of the voltage of net56.
```

```
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.
```
ocnSetAttrib


Description

Sets the waveform window plotting attributes.

Arguments

**XAxisLabel**

Label (symbol or string) for the X axis in the waveform window.

**YAxisLabel**

Label (symbol or string) for the Y axis associated with the `stripNumber` and `yNumber` in the waveform window.

**XScale**

Scale of the X axis in the waveform window.

Valid values (symbols): `'auto`, `'log`, and `'linear`

**YScale**

Scale of the Y axis associated with the `stripNumber` and `yNumber` in the waveform window.

Valid values (symbols): `'log` and `'linear`

**XLimit**

Displays limits of the X axis in the waveform window.

Valid values: List of two numbers or `'auto` (symbol).

The first number in the list indicates the minimum limit and the second indicates the maximum limit.

 `'auto` sets the limit to autoscale.

**YLimit**

Displays limits of the Y axis associated with the `stripNumber` and `yNumber` in the waveform window.

Valid values: List of two numbers or `'auto` (symbol).

The first number in the list indicates the minimum limit and the second indicates the maximum limit.

 `'auto` sets the limit to autoscale.

**YRange**

Y range (integer) of the waveforms associated with the `stripNumber` and `yNumber` in the waveform window.

**Origin**

Axes origin of the waveform window.

Valid values: List of two numbers.
Note: The valid range for stripNumber is 1-20 and that for yNumber is 1-4.

Value Returned

\[ t \] Returns \( t \) if the values of all arguments are set successfully.

\[ \text{nil} \] Returns \( \text{nil} \) if one or more arguments fail to set as specified.

Example

\[ \text{ocnSetAttrib(?XAxisLabel } 'XMylabel \ ?YAxisLabel 'YMyLabelt \ ?stripNumber 2 \ ?yNumber 1 ) } \]
\[ \Rightarrow t \]

Sets the X and Y axis labels to \( X\text{Mylabel} \) and \( Y\text{MyLabel} \), respectively.

\[ \text{ocnSetAttrib(?XScale } '\text{log} \ ?YScale 'linear \ ?stripNumber 2 \ ?yNumber 1 ) } \]
\[ \Rightarrow t \]

Sets the scale of X and Y axis to \( \text{log} \) and \( \text{linear} \), respectively.

\[ \text{ocnSetAttrib(?XScale } '\text{auto} \ ?XLimit '(3 7) \ ?YLimit 'auto \ ?stripNumber 2 \ ?yNumber 1) } \]
\[ \Rightarrow t \]

Sets the scale of X axis to autoscale. Sets the Y display limits to autoscale.
ocnYvsYplot

=> 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

- **o_wavex**: Reference wave against which the wave provided needs to be plotted.
- **o_wavey**: Wave to be plotted against the reference wave.
- **o_exprx**: Reference expression against which the expression provided needs to be plotted.
- **o_expry**: Expression to be plotted against the reference expression.
- **l_titleList**: List of waveform titles. If the waveform is simple, only one label will be required. If the waveform is param, a list of labels needs to be provided.
- **l_colorList**: 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

- **wave**: Returns the waveform specified.
- **nil**: Returns nil if the plot could not be generated.
Example
wy = VT("/vout")
wx = VT("/vin")
ex = "VT('/vin')"
ey = "VT('/vout')"

ocnYvsYplot(?wavex wx ?wavey wy ?titleList '("simpleWave") ?colorList '(3))

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

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

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

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

Description

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

**Note:** `plot` is implemented as a macro and not as a SKILL function. Therefore, the functions that expect a function name as an argument will not accept `plot` as a valid argument. For example, the following call to the function `apply` is not valid:

```
apply('plot)
```

Arguments

<table>
<thead>
<tr>
<th>Argument</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>o_waveform1</td>
<td>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>srrWave:XXXXX</code>.)</td>
</tr>
<tr>
<td>o_waveform2</td>
<td>Additional waveform object.</td>
</tr>
<tr>
<td>l_yNumberList</td>
<td>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</td>
</tr>
<tr>
<td>l_exprList</td>
<td>List of strings used to give names to the waveform objects.</td>
</tr>
<tr>
<td>x_stripNumber</td>
<td>An integer using which you can plot waveforms selectively on different strips and subwindows. If you specify an integer, it is used as the strip for all waveforms. To use the strip option for multiple waveforms, you can specify a list of strip numbers.</td>
</tr>
</tbody>
</table>

**Important**

*WaveScan* does not support `stripNumber`. 
### Value Returned

<table>
<thead>
<tr>
<th>Value</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td><code>t</code></td>
<td>Returns <code>t</code> if the waveforms are plotted.</td>
</tr>
<tr>
<td><code>nil</code></td>
<td>Returns <code>nil</code> and an error message if the waveforms cannot be plotted.</td>
</tr>
</tbody>
</table>
Example

plot(v("/net56") )

Plots the waveform for the voltage of net56.

plot( vm("/net56") vp("/net56") )

Plots the waveforms for the magnitude of the voltage of net56 and the phase of the voltage of net56.

plot( v("OUT") i("VFB") ?expr list("voltage" "current") )

Plots the waveforms, but changes one legend label from v("OUT") to voltage and changes the other legend label from i("VFB") to current.

plot( v("OUT") i("VFB") ?yNumber list(1 2) )

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

plot(wave1 wave2 wave3 ?strip list(1 2 2))

Plots wave1 to strip 1, and wave2 and wave3 to strip 2.
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

<table>
<thead>
<tr>
<th>Argument</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>auto</td>
<td>The appropriate plotting style is automatically chosen.</td>
</tr>
<tr>
<td>scatterplot</td>
<td>Data points are not joined.</td>
</tr>
<tr>
<td>bar</td>
<td>Vertical bars are drawn at each data point that extend from the point to the bottom of the graph.</td>
</tr>
<tr>
<td>joined</td>
<td>Each data point is joined to adjacent data points by straight-line segments.</td>
</tr>
</tbody>
</table>

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.
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

<table>
<thead>
<tr>
<th>Argument</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>t_resultsDir</td>
<td>Directory containing the results. If you specify a directory with resultsDir, the pzPlot command plots the results for that directory.</td>
</tr>
<tr>
<td>S_resultName</td>
<td>Pointer to results from the analysis for which you want to plot the report.</td>
</tr>
<tr>
<td>S_toPlot</td>
<td>Use this option to plot only poles, only zeros or both poles and zeros information. Valid values: ‘poles,’ ‘zeros,’ ‘polesZeros.’</td>
</tr>
<tr>
<td>f_fval</td>
<td>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.</td>
</tr>
<tr>
<td>f_rval</td>
<td>Real value which is used to filter out poles and zeros whose real value are less than or equal to the value specified.</td>
</tr>
</tbody>
</table>

Value Returned

<table>
<thead>
<tr>
<th>Value</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>t</td>
<td>Returns t if it plots a report.</td>
</tr>
<tr>
<td>nil</td>
<td>Returns nil otherwise.</td>
</tr>
</tbody>
</table>
Example

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).
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

t_resultsDir Directory containing the results. If you specify a directory with
resultsDir, the pzSummary command plots the results for that
directory.

S_resultName Pointer to results from the analysis for which you want to print the
report.

S_toPlot Use this option to plot only poles, only zeros or both poles and
zeros information.
Valid values: 'poles, 'zeros, 'polesZeros.

f_fval Maximum pole and zero frequency value to filter out poles and
zeros that are outside the frequency band of interest (FBOI) and
that do not influence the transfer function in the FBOI.

f_rval Real value which is used to filter out poles and zeros whose real
value are less than or equal to the value specified.

Value Returned

t Returns t if it prints a report.

nil Returns nil otherwise.
**Example**

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

Prints a report for all the poles and zeros for the specified results.

pzSummary(?resultsDir "/usr/simulation/lowpass/spectre/schematic" ?print 'poles)

Prints a report containing only poles for the specified results.

pzSummary( ?print 'zeros ?realfilter -1.69e-01)

Prints a report for all those zeros whose real values are 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).
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.

Example
label = addWindowLabel( list( 0.75 0.75 ) "test" )
Adds the "test" label to the current subwindow at the specified coordinates and stores the
label identification number in label.
removeLabel( label )
Removes the label whose identification number is stored in label. In this case, the "test"
label is removed.
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

<table>
<thead>
<tr>
<th>Argument</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>t_filename</td>
<td>File in which to write the information. The report command opens the file, writes to the file, and closes the file. If you specify the filename without a path, the OCEAN environment creates the file in the directory pointed to by your Skill Path. To find out what your Skill path is, type getSkillPath() at the OCEAN prompt.</td>
</tr>
<tr>
<td>p_port</td>
<td>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.</td>
</tr>
<tr>
<td>t_type</td>
<td>Type of information to print, such as all bjts.</td>
</tr>
<tr>
<td>t_name</td>
<td>Name of the node or component.</td>
</tr>
<tr>
<td>t_param</td>
<td>Name of the parameter to print. It is also a list.</td>
</tr>
<tr>
<td>s_reportStyle</td>
<td>Specifies the format of the output. Valid values: spice and paramValPair Default value: paramValPair</td>
</tr>
</tbody>
</table>
The `spice` format looks like this:

<table>
<thead>
<tr>
<th>Name</th>
<th>Param1</th>
<th>Param2</th>
<th>Param3</th>
</tr>
</thead>
<tbody>
<tr>
<td>Name1</td>
<td>value</td>
<td>value</td>
<td>value</td>
</tr>
<tr>
<td>Name2</td>
<td>value</td>
<td>value</td>
<td>value</td>
</tr>
<tr>
<td>Name3</td>
<td>value</td>
<td>value</td>
<td>value</td>
</tr>
</tbody>
</table>

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.

**Example**

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

Prints all the operating-point parameters.

```c
report( ?type "bjt" )
=> t
```

Prints all the `bjt` operating-point parameters.

```c
report( ?type "bjt" ?param "ib" )
=> t
```

Prints the `ib` parameter for all bjts.
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.
**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.
**yLimit**

```plaintext
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**

- `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`.
- `x_yNumber`: Specifies the Y axis that will have limited display with the range specified by `l_minMax`. Valid values: 1 through 4
- `x_stripNumber`: Specifies the strip in which the y display is to be limited as specified by `x_yNumber`. Valid values: 1 through 20

**Value Returned**

- `t`: Returns `t` if the Y axis display limits are set.
- `nil`: Returns `nil` and an error message if the Y axis display limits cannot be set.

**Example**

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

Sets Y axis 1 to display from 4.5 to 7.5.

```plaintext
yLimit( list( 4.5 7.5 ) ?yNumber 1 ?stripNumber 3)
```
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), NF (noise figure), and RN (noise resistance) results follows. Plotting NNR (normalized noise resistance) is very similar to plotting RN.

```ocean
; start ocean with Spectre as the as the simulator.
simulator( 'spectre )
; specify design and model path
design( "/usr1/mnt4/myhome/simulation/myckt/schematic/netlist/myckt.c")
path( "/usr1/mnt4/myhome/models" )
; specify analysis used: sp with noise
analysis(’sp ?start "100M" ?stop "10G" ?donoise "yes"
?oprobe "/PORT1" ?iprobe "/PORT0" )
;set design variables
desVar( "r2" 37)
desVar( "r1" 150)
;set temperature
temp( 25 )
;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 Virtuoso Spectre Circuit Simulator RF Analysis User Guide.

For more information on these functions, click these links: getData, sp, ocnPrint, and plot.
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.

\[ \text{vm( } t\_\text{net } [?\text{result } s\_\text{resultName}] \text{ )} \]

where \( s\_\text{resultName} \) 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:

\[ \text{vm( } t\_\text{net } [?\text{resultsDir } t\_\text{resultsDir} ][?\text{result } s\_\text{resultName}] \text{ )} \]

where \( t\_\text{resultsDir} \) is the name of a different directory containing PSF results, and \( s\_\text{resultName} \) is the name of a datatype contained in that directory. (If you specify another directory with \( t\_\text{resultsDir} \), you must also specify the particular results with \( s\_\text{resultName} \).)

List of Aliases

<table>
<thead>
<tr>
<th>Alias</th>
<th>Syntax</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>( \text{vm} )</td>
<td>( \text{vm( } t_\text{net } [?\text{resultsDir } t_\text{resultsDir}][?\text{result } s_\text{resultName}] \text{ )} \Rightarrow o_\text{waveform/nil} )</td>
<td>Aliased to ( \text{mag(}v()\text{)} ). Gets the magnitude of the voltage of a net.</td>
</tr>
<tr>
<td>( \text{vdb} )</td>
<td>( \text{vdb( } t_\text{net } [?\text{resultsDir } t_\text{resultsDir}][?\text{result } s_\text{resultName}] \text{ )} \Rightarrow o_\text{waveform/nil} )</td>
<td>Aliased to ( \text{db20(}v()\text{)} ). Gets the power gain in decibels from net in to net out.</td>
</tr>
<tr>
<td>( \text{vp} )</td>
<td>( \text{vp( } t_\text{net } [?\text{resultsDir } t_\text{resultsDir}][?\text{result } s_\text{resultName}] \text{ )} \Rightarrow o_\text{waveform/nil} )</td>
<td>Aliased to ( \text{phase(}v()\text{)} ). Gets the phase of the voltage of a net.</td>
</tr>
</tbody>
</table>
## List of Aliases, continued

<table>
<thead>
<tr>
<th>Alias</th>
<th>Syntax</th>
<th>Aliased to</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>vr</td>
<td>vr(t_net [[?resultsDir t_resultsDir][?result s_resultName]) =&gt; o_waveform/nil</td>
<td><code>real(v())</code></td>
<td>Gets the real part of a complex number representing the voltage of a net.</td>
</tr>
<tr>
<td>vim</td>
<td>vim(t_net [[?resultsDir t_resultsDir][?result s_resultName]) =&gt; o_waveform/nil</td>
<td><code>imag(v())</code></td>
<td>Gets the imaginary part of a complex number representing the voltage of a net.</td>
</tr>
<tr>
<td>im</td>
<td>im(t_component [[?resultsDir t_resultsDir][?result s_resultName]) =&gt; o_waveform/nil</td>
<td><code>mag(i())</code></td>
<td>Gets the magnitude of the AC current through a component.</td>
</tr>
<tr>
<td>ip</td>
<td>ip(t_component [[?resultsDir t_resultsDir][?result s_resultName]) =&gt; o_waveform/nil</td>
<td><code>phase(i())</code></td>
<td>Gets the phase of the AC current through a component.</td>
</tr>
<tr>
<td>ir</td>
<td>ir(t_component [[?resultsDir t_resultsDir][?result s_resultName]) =&gt; o_waveform/nil</td>
<td><code>real(i())</code></td>
<td>Gets the real part of a complex number representing the AC current through a component.</td>
</tr>
<tr>
<td>iim</td>
<td>iim(t_component [[?resultsDir t_resultsDir][?result s_resultName]) =&gt; o_waveform/nil</td>
<td><code>imag(i())</code></td>
<td>Gets the imaginary part of a complex number representing the AC current through a component.</td>
</tr>
</tbody>
</table>
Predefined 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 L SKILL Language Reference*.

- **Predefined Arithmetic Functions**
  - abs
  - acos
  - add1
  - asin
  - atan
  - cos
  - exp
  - int
  - linRg
  - log
  - logRg
  - logRg
  - max
  - min
  - mod
  - random
  - round
Predefined and Waveform (Calculator) Functions

- sin
- sqrt
- srandom
- sub1
- tan
- xor

- Waveform (Calculator) Functions
  - average
  - awvPlaceXMarker
  - awvPlaceYMarker
  - b1f
  - bandwidth
  - clip
  - compare
  - compression
  - compressionVRI
  - compressionVRICurves
  - conjugate
  - convolve
  - cPwrContour
  - cReflContour
  - cross
  - db10
  - db20
  - dbm
  - delay
  - deriv
dft

dftbb

dnl
dutyCycle
evrmQAM
evrmQpsk
eyeDiagram
flip
fourEval
freq
freq_jitter
frequency
ga
gac
gainBwProd
gainMargin
gmax
gmin
gmsg
gmux
gp
gpc
groupDelay
gt
harmonic
harmonicFreqList
harmonicList
histo
iinteg
imag
integ
intersect
ipn
ipnVRI
ipnVRICurves
kf
ln
log10
lsb
lshift
mag
nc
overshoot
peak
peakToPeak
period_jitter
phase
phaseDeg
phaseDegUnwrapped
phaseMargin
phaseRad
phaseRadUnwrapped
pow
psd
psdbb
pzbode
pzfilter
real
riseTime
rms
rmsNoise
root
rshift
sample
settlingTime
slewRate
spectralPower
spectrum
ssb
stddev
tangent
thd
unityGainFreq
value
xmax
xmin
xval
ymax
ymin
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**

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

**Trigonometric Functions**

<table>
<thead>
<tr>
<th>Synopsis</th>
<th>Result</th>
</tr>
</thead>
<tbody>
<tr>
<td>sin(n)</td>
<td>sine, argument n is in radians</td>
</tr>
<tr>
<td>cos(n)</td>
<td>cosine</td>
</tr>
<tr>
<td>tan(n)</td>
<td>tangent</td>
</tr>
<tr>
<td>asin(n)</td>
<td>arc sine, result is in radians</td>
</tr>
<tr>
<td>acos(n)</td>
<td>arc cosine</td>
</tr>
<tr>
<td>atan(n)</td>
<td>arc tangent</td>
</tr>
</tbody>
</table>
### Predefined Arithmetic Functions

<table>
<thead>
<tr>
<th>Synopsis</th>
<th>Result</th>
</tr>
</thead>
<tbody>
<tr>
<td><strong>Random Number Generator</strong></td>
<td></td>
</tr>
<tr>
<td>random(x)</td>
<td>Returns a random integer between 0 and x-1. If random is called with no arguments, it returns an integer that has all of its bits randomly set.</td>
</tr>
<tr>
<td>srandom(x)</td>
<td>Sets the initial state of the random number generator to x.</td>
</tr>
</tbody>
</table>
abs

\[
\text{abs( } n\_\text{number } \text{)}
\]

\[
=> n\_\text{result}
\]

Description

Returns the absolute value of a floating-point number or integer.

Arguments

\[
n\_\text{number}
\]

Floating-point number or integer.

Value Returned

\[
n\_\text{result}
\]

The absolute value of \( n\_\text{number} \).

Example

\[
\text{abs( } -209.625 \text{)}
\]

\[
=> 209.625
\]

\[
\text{abs( } -23 \text{)}
\]

\[
=> 23
\]
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
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.
asinh

\( \text{asinh} \left( n\_\text{number} \right) \)

\( \Rightarrow f\_\text{result} \)

**Description**

Returns the arc sine of a floating-point number or integer.

**Arguments**

\( n\_\text{number} \)

Floating-point number or integer.

**Value Returned**

\( f\_\text{result} \)

The arc sine of \( n\_\text{number} \).

**Example**

\( \text{asinh}(0.3) \)

\( \Rightarrow 0.3046927 \)
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 and Waveform (Calculator) Functions

---

**COS**

\[
\cos( \text{n\_number} ) \\
=> \text{f\_result}
\]

**Description**

Returns the cosine of a floating-point number or integer.

**Arguments**

\text{n\_number}  
Floating-point number or integer.

**Value Returned**

\text{f\_result}  
The cosine of \text{n\_number}.

**Example**

\[
\cos(0.3) \\
=> 0.9553365 \\
\cos(3.14/2) \\
=> 0.0007963
\]
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.

Example

exp( 1 )
 => 2.718282
exp( 3.0 )
 => 20.08554
**int**

```latex
\texttt{int( n\_arg )}
\rightarrow x\_result
```

**Description**

Returns the largest integer not larger than the given argument.

**Note:** This function works on vector as well as waveform data. The function is applied to individual elements of the vector and waveform data.

**Arguments**

- **n_arg**
  - A numeric value (which can be integer or floating point number).

**Value Returned**

- **x_result**
  - The value of the largest integer not larger than the value specified by `n_arg`.

**Example**

```plaintext
\texttt{int( 3.01 )}
\rightarrow 3
\texttt{int( 3.99 )}
\rightarrow 3
```
**linRg**

linRg( n_from n_to n_by )  
=> l_range/nil

**Description**

Returns a list of numbers in the linear range from \( n_{\text{from}} \) to \( n_{\text{to}} \) incremented by \( n_{\text{by}} \).

**Arguments**

- \( n_{\text{from}} \): Smaller number in the linear range.
- \( n_{\text{to}} \): Larger number in the linear range.
- \( n_{\text{by}} \): Increment value when stepping through the range.

**Value Returned**

- \( l_{\text{range}} \): 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)
```
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
**logRg**

\[
\text{logRg}( n_{\text{from}} \text{ n_to } n_{\text{by}} ) \\
\Rightarrow \ l\_\text{range}/\text{nil}
\]

**Description**

Returns a list of numbers in the log10 range from \( n_{\text{from}} \) to \( n_{\text{to}} \) advanced by \( n_{\text{by}} \).

The list is a geometric progression where the multiplier is 10 raised to the \( 1/n_{\text{by}} \) power. For example if \( n_{\text{by}} \) is 0.5, the multiplier is 10 raised to the 2nd power or 100.

**Arguments**

- \( n_{\text{from}} \) Smaller number in the linear range.
- \( n_{\text{to}} \) Larger number in the linear range.
- \( n_{\text{by}} \) Increment value when stepping through the range.

**Value Returned**

- \( l\_\text{range} \) List of numbers in the linear range.
- \( \text{nil} \) Returned if error.

**Example**

\[
\text{logRg}(1 \ 1M \ 0.5) \\
(1.0 \ 100.0 \ 10000.0 \ 1000000.0)
\]
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

n_num1        First value to check.
n_num2        Next value to check.
[n_num3...]    Additional values to check.

Value Returned

n_result        The maximum of the values passed in.

Example

max(3 2 1)
=> 3
max(-3 -2 -1)
=> -1
**min**

\[ \text{min}( n\text{\_}num1 \ n\text{\_}num2 \ [n\text{\_}num3 \ldots] ) \Rightarrow n\text{\_}result \]

**Description**

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

**Arguments**

- \( n\text{\_}num1 \)
  - First value to check.
- \( n\text{\_}num2 \)
  - Next value to check.
- \([n\text{\_}num3\ldots]\)
  - Additional values to check.

**Value Returned**

- \( n\text{\_}result \)
  - The minimum of the values passed in.

**Example**

\[
\begin{align*}
\text{min}(1 \ 2 \ 3) & \Rightarrow 1 \\
\text{min}(-1 \ -2.0 \ -3) & \Rightarrow -3.0
\end{align*}
\]
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
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.

ValueReturned

x_result
  Returns a random integer between 0 and x_number minus 1.

Example

random( 93 )
=> 26
**round**

\[
\text{round}\left(n_{\text{arg}}\right) \\
\Rightarrow x_{\text{result}}
\]

**Description**

Rounds a floating-point number to its closest integer value.

**Arguments**

\[
n_{\text{arg}} \quad \text{Floating-point number.}
\]

**Value Returned**

\[
x_{\text{result}} \quad \text{The integer whose value is closest to } n_{\text{arg}}.
\]

**Example**

\[
\text{round}(1.5) \\
\Rightarrow 2
\]

\[
\text{round}(-1.49) \\
\Rightarrow -1
\]

\[
\text{round}(1.49) \\
\Rightarrow 1
\]
sin

\[ \sin(n_{\text{number}}) \]
\[ \Rightarrow f_{\text{result}} \]

Description

Returns the sine of a floating-point number or integer.

Arguments

- \( n_{\text{number}} \):
  - Floating-point number or integer.

Value Returned

- \( f_{\text{result}} \):
  - The sine of \( n_{\text{number}} \).

Example

- \( \sin(3.14/2) \)
  \[ \Rightarrow 0.9999997 \]
- \( \sin(3.14159/2) \)
  \[ \Rightarrow 1.0 \]

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

Example

sqrt( 49 )
=> 7.0

sqrt( 43942 )
=> 209.6235
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
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.
**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**

\[
\tan( 3.0 ) \\
=> -0.1425465
\]
xox

xor( b_in1 b_in2 ) => b_res

Description

Returns the XOR value of the boolean inputs.

Arguments

b_in1 The first boolean input.

b_in2 The second boolean input.

Value Returned

b_res The resultant XOR value.

Example

xor(nil nil) => nil
xor(t nil) => t
xor(nil t) => t
xor(t t) => nil
Waveform (Calculator) Functions

The calculator commands are described in this section.
**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 $y=f(x)$, $\text{average}(y) =

\[
\frac{\int_{\text{from}}^{\text{to}} f(x) \, dx}{\text{to} - \text{from}}
\]

where $\text{from}$ is the initial value for $x$ and $\text{to}$ is the final value.

**Arguments**

- **o_waveform**
  
  Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: `srrWave:XXXXX`.)

**Value Returned**

- **n_average**
  
  Returns a number representing the average value of the input waveform.

- **o_waveformAverage**
  
  Returns a waveform (or family of waveforms) representing the average value if the input is a family of waveforms.

- **nil**
  
  Returns `nil` and an error message otherwise.
Example

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

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

- **w_windowId** Waveform window ID.
- **n_xLoc** The x-coordinate at which to place the marker.
- **x_subwindowId** Waveform subwindow ID.

**Value Returned**

- **t_xLoc** Returns a string of x-coordinates if the command is successful and the vertical marker info form is opened.
- **t** Returns this when the command is successful but the vertical marker info form is not opened.
- **nil** Returns nil or an error message.

**Example**

awvPlaceXMarker( window 5) => "5"

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

awvPlaceXMarker( window 6 ?subwindow 2) => t

Vertical marker info form is not opened.
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

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

Value Returned

  t_yLoc      Returns a string of y-coordinates if the command is successful
               and the horizontal marker info form is opened.
  t           Returns this when the command is successful but the horizontal
               marker info form is not opened.
  nil         Returns nil or an error message.

Example

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.
awvRefreshOutputPlotWindows

awvRefreshOutputPlotWindows(s_session)

Description

Refreshes all existing plot windows (with new simulation data, if any) attached with the session s_session.

Arguments

s_session Currently active environment variable.

Value Returned

None.
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

- o_s11: Waveform object representing s11.
- o_s12: Waveform object representing s12.
- o_s21: Waveform object representing s21.
- o_s22: Waveform object representing s22.

Value Returned

- o_waveform: Waveform object representing the alternative stability factor.
- nil: Returns nil and an error message otherwise.

Example

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

bandwidth( o_waveform n_db t_type )
   => n_value/o_waveform/nil

Description
Calculates the bandwidth of a waveform.

Arguments

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

  n_db        Positive number that defines the bandwidth.

  t_type      Type of input filter.
              Valid values: "low", "high" or "band".

Value Returned

  n_value     Returns a number representing the value of the bandwidth if the input argument is a single waveform.

  o_waveform  Returns a waveform (or family of waveforms) representing the bandwidth if the input argument is a family of waveforms.

  nil         Returns nil and an error message otherwise.

Example

bandwidth( v( "/OUT" ) 3 "low")

Gets the 3 dB bandwidth of a low-pass filter.

bandwidth( v( "/OUT" ) 4 "band" )

Gets the 4 dB bandwidth of a band-pass filter.
clip

clip( o_waveform n_from n_to ) => o_waveform/nil

Description
Restricts the waveform to the range defined by \textit{n_from} and \textit{n_to}.

You can use the \texttt{clip} function to restrict the range of action of other commands. If \textit{n_from} is \texttt{nil}, \textit{n_from} is taken to be the first X value of the waveform, and if \textit{n_to} is \texttt{nil}, \textit{n_to} is taken to be the last X value of the waveform. If both \textit{n_to} and \textit{n_from} are \texttt{nil}, the original waveform is returned.

Arguments

\textit{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: \texttt{srrWave:XXXXX}.)

\textit{n\_from} \\
Starting value for the range on the X axis.

\textit{n\_to} \\
Ending value for the range on the X axis.

Value Returned

\textit{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.

\texttt{nil} \\
Returns \texttt{nil} and an error message otherwise.

Example

\texttt{x = clip( v( "/net9" ) 2m 4m )}
\texttt{plot( x )}

Plots the portion of a waveform that ranges from 2 ms to 4 ms.

\texttt{plot( clip( v( "/net9" ) nil nil ) )}

Plots the original waveform.

\texttt{plot( clip( v( "/net9" ) nil 3m ) )}
Plots the portion of a waveform that ranges from 0 to 3 ms.
**compare**

```plaintext
compare( o_waveform1 o_waveform1 [f_abstol [f_reltol]] )
   => o_comparisonWaveform/nil
```

**Description**

Compares the two given waveforms based on the specified values for absolute and relative tolerances. This function compares only the sections of the two waveforms where the X or independent axes overlap.

The following situations are possible:

- If neither relative nor absolute tolerance is specified, the function returns the difference of the two waveforms (o_waveform1 - o_waveform2).

- If only the absolute tolerance is specified, the function returns the difference of the two waveforms only when the absolute value of the difference is greater than the absolute tolerance (|o_waveform1 - o_waveform2| > f_abstol); otherwise it returns a zero waveform.

- If only the relative tolerance is specified, the function returns the difference of the two waveforms only when the absolute value of the difference is greater than the product of the relative tolerance and the larger of the absolute values of the two waveforms (|o_waveform1 - o_waveform2| > f_reltol * max(|o_waveform1|, |o_waveform2|)); otherwise it returns a zero waveform.

- If both relative and absolute tolerances are specified, the function returns the difference of the two waveforms only when the absolute value of the difference is greater than the sum of the separately calculated tolerance components (|o_waveform1 - o_waveform2| > f_abstol + f_reltol * max(|o_waveform1|, |o_waveform2|)); otherwise it returns a zero waveform.

**Note:** The function also compares parametric waveforms. However, for a successful comparison of parametric waveforms, the family tree structures of the two input waveforms should be the same. For both the input waveforms, the number of child waveforms at each level should also be the same, except at the leaf level where the elements are simple scalars. This is an obvious condition to obtain a meaningful comparison.

**Arguments**

- `o_waveform1` Waveform 1.
- `o_waveform2` Waveform 2.
f_abstol
Absolute tolerance.
Default value: 0.0

f_reltol
Relative tolerance.
Default value: 0.0

Value Returned

o_comparisonWaveform
Returns the difference of the two given waveforms based on the specified values of the relative and absolute tolerances.

nil
Returns nil and an error message otherwise.

Example

compare( wave1 wave2 2.2 0.4 ) => srrWave:175051528

Returns the difference of the waveforms wave1 and wave2 based on the specified absolute and relative tolerances of 2.2 and 0.4, respectively.
**compression**

```lisp
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**

- **o_waveform**
  Waveform object representing output power (in dBm) versus input power (in dBm).

- **f_x**
  The X coordinate value (in dBm) used to indicate the point on the output power waveform where the constant-slope power line begins. This point should be in the linear region of operation. Default value: Unless \( f_y \) is specified, defaults to the X coordinate of the first point of the `o_waveform` wave.

- **f_y**
  The Y coordinate value (in dBm) used to indicate the point on the output power waveform where the constant-slope power line begins. This point should be in the linear region of operation. Default value: Unless \( f_x \) is specified, defaults to the Y coordinate of the first point of the `o_waveform` wave.

- **f_compression**
  The delta (in dB) between the power waveform and the ideal gain line that marks the compression point. Default value: 1

- **s_measure**
  Symbol indicating whether the measurement is to be input referred (‘input) or output referred (‘output). Default value: ‘input
Value Returned

\[ f\_compPoint \] Depending on the setting of \( s\_measure \), returns either input referred or output referred compression point.

nil Returns nil and an error message otherwise.

Example

\[
xloc = \text{compression( wave ?x -25 ?compress 1)}
\]
\[
yloc = \text{compression( wave ?x -25 ?measure "Output")}
\]
; Each of following returns a compression measurement:
\[
\text{compression(dB20(harmonic(v("/Pif" result "pss_fd") 2)))}
\]
\[
\text{compression(dbm(harmonic(spectralPower(v("/Pif"
  \ result "pss_fd")/ 50.0
  v("/Pif" result "pss_fd")(2))))}
\]
\[
\text{compression(dbm(harmonic(spectralPower(v("/Pif"
  \ result "pss_fd")/resultParam("rif:r"
  \ result "pss_td") v("/Pif"
  \ result "pss_fd")(2))))}
\]
\[
\text{compression(dbm(harmonic(spectralPower(i("/rif/PLUS"
  \ result "pss_fd") v("/Pif" result "pss_fd")(2))}
  \ x -25 ?compress 0.1 ?measure "Output")}
\]
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 \( \text{dBm(spectralPower((i or v/r),v))} \) 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

$o_vport$  
Voltage across the output port. This argument must be a family of spectrum waveforms (1 point per harmonic) created by parametrically sweeping an input power (in dBm) of the circuit.

$x_harm$  
Harmonic index of the voltage wave contained in $o_vport$. When $o_iport$ is specified, also applies to a current waveform contained in $o_iport$.

$o_iport$  
Current into the output port. This argument must be a family of spectrum waveforms (1 point per harmonic) created by parametrically sweeping an input power (in dBm) of the circuit. When specified, the output power is calculated using voltage and current.  
Default value: nil

$f_rport$  
Resistance into the output port. When specified and $o_iport$ is nil, the output power is calculated using voltage and resistance.  
Default value: 50
Ocean Reference
Predefined and Waveform (Calculator) Functions

- **f_epoint**: The X coordinate value (in dBm) used to indicate the point on the output power waveform where the constant-slope power line begins. This point should be in the linear region of operation. Default value: the X coordinate of the first point of the `o_waveform` wave

- **f_gcomp**: The delta (in dB) between the power waveform and the ideal gain line that marks the compression point. Default value: 1

- **s_measure**: Symbol indicating if measurement is to be input referred (`input`) or output referred (`output`). Default value: `input`

**Value Returned**

- **o_waveform**: Returns a waveform when `o_waveform1` is a family of waveforms.

- **f_number**: Returns a number when `o_waveform1` is a waveform.

- **nil**: Returns nil and an error message otherwise.

**Example**

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")
```
**CompressionVRICurves**

```ruby
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 nth 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**

- **o_vport**
  Voltage across the output port. This argument must be a family of spectrum waveforms (1 point per harmonic) created by parametrically sweeping an input power (in dBm) of the circuit.

- **x_harm**
  Harmonic index of the wave contained in `o_vport`. When `o_iport` is specified, also applies to a current waveform contained in `o_iport`.

- **o_iport**
  Current into the output port. This argument must be a family of spectrum waveforms (1 point per harmonic) created by parametrically sweeping an input power (in dBm) of the circuit. When specified, the output power is calculated using voltage and current.
  Default value: nil

- **f_rport**
  Resistance into the output port. When specified and `o_iport` is nil, the output power is calculated using voltage and
resistance.
Default value: 50

\( f_{\text{epoint}} \)
The X coordinate value (in dBm) used to indicate the point on the output power waveform where the constant-slope power line begins. This point should be in the linear region of operation. Default value: the X coordinate of the first point of the \( o_{\text{waveform}} \) wave

\( f_{\text{gcomp}} \)
The delta (in dB) between the power waveform and the ideal gain line that marks the compression point.
Default value: 1

Value Returned

\( o_{\text{waveform}} \)
Returns a family of waveforms containing the output power and tangent line.

\( \text{nil} \)
Returns \( \text{nil} \) and an error message otherwise.

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

\[
\text{compressionVRICurves}(v("/Pif" ?result "pss_fd") 2)
\]

\[
\text{compressionVRICurves}(v("/Pif" ?result "pss_fd") 2
  ?rport resultParam("rif:r" ?result "pss_td"))
\]

\[
\text{compressionVRICurves}(v("/Pif" ?result "pss_fd") 2
  ?iport i("/rif/PLUS" ?result "pss_fd") ?epoint -25
  ?gcomp 0.1)
\]
**complex**

`complex( f_real f_imaginary )`  
=> `o_complex`

**Description**

Creates a complex number of which the real part is equal to the real argument, and the imaginary part is equal to the imaginary argument.

**Arguments**

- `f_real`  
The real part of the complex number.
- `f_imaginary`  
The imaginary part of the complex number.

**Value Returned**

- `o_complex`  
Returns the complex number.

**Example**

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

Arguments

$g\_value$  
A skill object.

Value 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
conjugate

conjugate( {o_waveform | n_x} )
=> o_waveform/n_y/nil

Description

Returns the conjugate of a waveform or number.

Arguments

- **o_waveform**: Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: `srrWave:XXXXX`.)
- **n_x**: Complex or imaginary number.

Value Returned

- **o_waveform**: Returns the conjugate of a waveform if the input argument is a waveform.
- **n_y**: Returns the result of `n_x` being mirrored against the real axis (X axis) if the input argument is a number.
- **nil**: Returns `nil` and an error message otherwise.

Example

For this example, assume that the first three statements are true for the `conjugate` function that follows them.

```plaintext
x=complex(-1 -2)
real(x) = -1.0
imag(x) = -1.0
conjugate(x) = complex(-1, 2)
```

Returns the conjugate of the input complex number.
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

\[
\int_{\text{from}}^{\text{to}} f_1(s)f_2(t-s)ds
\]

\(f_1\) and \(f_2\) are the functions defined by the first and second waveforms.

**Note:** The `convolve` function is numerically intensive and might take longer than the other functions to compute.

Arguments

- **o_waveform1**
  Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: `srrWave:XXXXX`.)

- **o_waveform2**
  Additional waveform object.

- **n_from**
  Starting point (X-axis value) of the integration range.

- **n_to**
  Ending point (X-axis value) of the integration range.

- **t_type**
  Type of interpolation.
  Valid values: "linear" or "log".

- **n_by**
  Increment.
Value Returned

\textit{o\_waveform} \hspace{1cm} \textit{n\_number} \hspace{1cm} \textit{nil}

\textit{o\_waveform} \hspace{1cm} 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.

\textit{n\_number} \hspace{1cm} Returns a value representing the convolution if both of the input arguments are numbers.

\textit{nil} \hspace{1cm} Returns \texttt{nil} and an error message otherwise.

Example

\texttt{sinWave = expr( n sin( n ) linRg( 0 20 0.01 ) )}

\texttt{triWave = artListToWaveform( `( ( -4, 0 ) ( -3, 1 ) ( -2, 0 ) ( -1, -1 ) ( 0, 0 ) ( 1, 1 ) ( 2, 0 ) ( 3, -1 ) ( 4, 0 ) )}

\texttt{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.
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 $x_{\text{harm}}$ 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

- **o_iwave**
  Current used to calculate power, expected to be a two-dimensional family of harmonic waveforms.

- **o_vwave**
  Voltage used to calculate power, expected to be a two-dimensional family of harmonic waveforms.

- **x_harm**
  Harmonic index of the waves contained in $o_iwave$ and $o_vwave$.

- **o_iwaveLoad**
  Current used to calculate reflection coefficient, expected to be a two-dimensional family of harmonic waveforms.
  Default value: $o_iwave$

- **o_vwaveLoad**
  Voltage used to calculate reflection coefficient, expected to be a two-dimensional family of harmonic waveforms.
  Default value: $o_vwave$

- **f_maxPower**
  Maximum power magnitude value for contours.
  Default value: automatic

- **f_minPower**
  Minimum power magnitude value for contours.
  Default value: automatic
**x_numCont**
Total number of contours returned.
Default value: 8

**f_refImp**
Reference resistance used to calculate reflection coefficients.
Default value: 50

**b_closeCont**
Boolean indicating when to close the contours. When `nil`, largest segment of each contour is left open.
Default value: `nil`

**s_modifier**
Symbol indicating the modifier function to apply to the calculated power. The modifier function can be any single argument OCEAN function such as `'db10` or `'dBm`.
Default value: `'mag`

**Value Returned**

**o_waveform**
Returns a family of waveforms (contours) for Z-Smith plotting.

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

**Example**

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

```ocean
   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:

```ocean
   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:

```ocean
   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")
```
cRefContour

\[
cRefContour( \text{o_iwave} \text{ o_vwave} \text{x_harm} [\text{o_iwaveLoad} \text{ o_iwaveLoad}] [\text{o_vwaveLoad} \text{ o_vwaveLoad}] \text{[f_maxRefl f_minRefl]} [\text{f_maxRefl} f_{\text{minRefl}}] [\text{f_maxRefl} f_{\text{minRefl}}] \text{[x_numCont x_numCont]} [\text{f_maxRefl} f_{\text{minRefl}}] \text{[x_numCont b_closeCont]} ) 
\Rightarrow \text{o_waveform/nil}
\]

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 \(x_{\text{harm}}\) harmonic is extracted from all the input waveforms. Reflection coefficient magnitude is calculated using voltage, current, reference resistance, and the \(\text{mag}\) function. The reference reflection coefficients are calculated separately by using voltage, current, and a reference resistance.

Arguments

- **o_iwave**
  
  Current used to calculate reflection coefficient magnitude, expected to be a two-dimensional family of spectrum waveforms.

- **o_vwave**
  
  Voltage used to calculate reflection coefficient magnitude, expected to be a two-dimensional family of spectrum waveforms.

- **x_harm**
  
  Harmonic index of the waves contained in \(\text{o_iwave}\) and \(\text{o_vwave}\).

- **o_iwaveLoad**
  
  Current used to calculate reference reflection coefficient, expected to be a two-dimensional family of harmonic waveforms. Default value: \(\text{o_iwave}\)

- **o_vwaveLoad**
  
  Voltage used to calculate reference reflection coefficient, expected to be a two-dimensional family of spectrum waveforms. Default value: \(\text{o_vwave}\)

- **f_maxRefl**
  
  Maximum reflection coefficient magnitude value for contours. Default value: automatic

- **f_minRefl**
  
  Minimum reflection coefficient magnitude value for contours. Default value: automatic
**x_numCont**  
Total number of contours returned.  
Default value: 8

**f_refImp**  
Reference resistance used to calculate reflection coefficients.  
Default value: 50

**b_closeCont**  
Boolean indicating when to close the contours. When nil, the largest segment of each contour is left open.  
Default value: nil

**Value Returned**

**o_waveform**  
Returns a family of waveforms (contours) for Z-Smith plotting.

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

**Example**

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

```plaintext
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:

```plaintext
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:

```plaintext
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)
```
cross

cross( o_waveform n_crossVal x_n s_crossType [b_multiple [s_Xname]] )
=> o_waveform/g_value/nil

Description

Computes the X-axis value at which a particular crossing of the specified edge type of the threshold value occurs.

Arguments

- **o_waveform**: Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: `srrWave:XXXXX`.)

- **n_crossVal**: Y-axis value at which the corresponding values of X are calculated.

- **x_n**: Number that specifies which X value to return. If `x_n` equals 1, the first X value with a crossing is returned. If `x_n` equals 2, the second X value with a crossing is returned, and so on. If you specify a negative integer for `x_n`, the X values with crossings are counted from right to left (from maximum to minimum).

- **s_crossType**: Type of the crossing.
  Valid values: ‘rising,’ ‘falling,’ ‘either.’

- **b_multiple**: An optional boolean argument that takes the value `nil` by default. If set to `t`, the value specified for the `x_n` argument is ignored and the function returns all occurrences of the crossing event.

- **s_XName**: An optional argument that is used only when `b_multiple` is set to `t`. It takes the value `time` by default. It controls the contents of the x vector of the waveform object returned by the function.
  Valid values: ‘time,’ ‘cycle’
Value Returned

\textit{o\_waveform} \hspace{1cm} \text{Returns a waveform if the input argument is a family of waveforms.}

\textit{g\_value} \hspace{1cm} \text{Returns the X-axis value of the crossing point if the input argument is a single waveform.}

\textit{nil} \hspace{1cm} \text{Returns nil and an error message otherwise.}

Example

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

Gets the time value (X axis) corresponding to specified voltage "/net9"=2.5V (Y axis) for the second rising edge.

cross( v( "/net9" ) 1.2 1 'either )

Gets the time value (X axis) corresponding to specified voltage "/net9"=1.2V (Y axis) for the first edge, which can be a rising or falling edge.

cross(VT("/out") 2.5 0 0 t "time") (s)

Returns multiple occurrences of crossing events specified against time-points at which each crossing event occurs.

cross(VT("/out") 2.5 0 0 t "cycle") (s)

Returns multiple occurrences of crossing events specified against cycle numbers, where a cycle number refers to the n'th occurrence of the crossing event in the input waveform.
**db10**

\[
\text{db10(} \{\text{o\_waveform} \mid \text{n\_number}\} \) \\
\Rightarrow \text{o\_waveform}/\text{n\_number}/\text{nil} 
\]

**Description**

Returns 10 times the log10 of the specified waveform object or number. This function can also be written as dB10.

**Arguments**

- \text{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: \text{srrWave:XXXXX}.)
- \text{n\_number} Number.

**Value Returned**

- \text{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.
- \text{n\_number} Returns a number if the input argument is a number.
- \text{nil} Returns \text{nil} and an error message otherwise.

**Example**

\[
\text{db10( ymax(v("/net9")) )} \\
\text{Returns a waveform representing log10(ymax(v("/net9"))) multiplied by 10.} \\
\text{db10(1000) } \\
\Rightarrow 30.0 \\
\text{Gets the value log10(1000) multiplied by 10, or 30.} 
\]
db20

\[ \text{db20}( \{ \text{o\_waveform} \mid \text{n\_number} \} ) \]
\[ \Rightarrow \text{o\_waveform/n\_number/nil} \]

**Description**

Returns 20 times the \( \log_{10} \) of the specified waveform object or number. This function can also be written as dB20.

**Arguments**

- **o\_waveform**
  
  Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave:XXXXX.)

- **n\_number**
  
  Number.

**Value Returned**

- **o\_waveform**
  
  Returns a waveform object if the input argument is a waveform object or returns a family of waveforms if the input argument is a family of waveforms.

- **n\_number**
  
  Returns a number if the input argument is a number.

- **nil**
  
  Returns nil and an error message otherwise.

**Example**

\[ \text{db20}( \text{ymax( v("/net9") )} ) \]

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

\[ \text{db20}( 1000 ) \]
\[ \Rightarrow 60.0 \]

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

```plaintext
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**

- **o_waveform**
  
  Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: `srrWave:XXXXX`.)

- **n_number**
  
  Number.

**Value Returned**

- **o_waveform**
  
  Returns a waveform object if the input argument is a waveform object or returns a family of waveforms if the input argument is a family of waveforms.

- **n_number**
  
  Returns a number if the input argument is a number.

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

**Example**

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

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

delay( ?waveform1 o_waveform1 ?value1 n_value1 ?edge1 s_edge1 ?nth1 x_nth1 ?td1
n_td1 ?waveform2 o_waveform2 ?value2 n_value2 ?edge2 s_edge2 ?nth2 x_nth2
{[?td2 n_td2] | ![td2r0 n_td2r0]} ?stop n_stop @rest args
[b_histoDisplay][x_noOfHistoBins])
=> 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

- **o_waveform1**: First waveform object.
- **n_value1**: Value at which the crossing is significant for the first waveform object.
- **s_edge1**: Type of the edge that must cross n_value1. Valid values: ‘rising,’ ‘falling,’ ‘either’
- **x_nth1**: Number that specifies which crossing is to be the trigger event. For example, if x_nth1 is 2, the trigger event is the second edge of the first waveform with the specified type that crosses n_value1.
- **n_td1**: Time at which to start the delay measurement. The simulator begins looking for the trigger event, as defined by o_waveform1, n_value1, t_edge1, and x_nth1, only after the n_td1 time is reached.
- **o_waveform2**: Second waveform object.
- **n_value2**: Value at which the crossing is significant for the second waveform.
- **s_edge2**: Type of the edge for the second waveform. Valid values: ‘rising,’ ‘falling,’ ‘either’
**Ocean Reference**

**Predefined and Waveform (Calculator) Functions**

---

$x_{nth2}$  
Number that specifies which crossing is to be the target event. For example, if $x_{nth2}$ is 2, the target event is the second edge of the second waveform with the specified type that crosses $n_{value2}$.

$n_{td2}$  
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 $o_{waveform2}, n_{value2}, t_{edge2},$ 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 $o_{waveform1}$ and $o_{waveform2}$ are specified. This parameter cannot be specified at the same time with $n_{td2}$.

The simulator begins looking for the target event, as defined by $o_{waveform2}, n_{value2}, t_{edge2},$ and $x_{nth2}$, only after the $n_{td2r0}$ 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.

$args$  
Variable list of arguments passed to the delay function (as created from the Calculator UI). These variables also include support for multiple occurrences of the delay event.

$b_{histoDisplay}$  
When set to t, returns a waveform that represents the statistical distribution of the riseTime data in the form of a histogram. The height of the bars (bins) in the histogram represents the frequency of the occurrence of values within the range of riseTime data.
**Value Returned**

- **o_waveform**
  Returns a waveform representing the delay if the input argument is a family of waveforms.

- **n_value**
  Returns the delay value if the input argument is a single waveform.

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

**Example**

```plaintext
delay( ?waveform1 wf1 ?value1 2.5 ?nth1 2 ?edge1 'either ?waveform2 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.

```plaintext
delay( ?td1 5 ?waveform2 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.

```plaintext
delay( ?waveform1 wf1 ?value1 2.5 ?nth1 1 ?edge1 'rising ?td1 5 ?waveform2 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.

```plaintext
delay(VT("/out"), 2.5, 1, 'rising, VT("/in"), 2.5, 1, 'rising', 1, 1, t)
```

Computes the delay between the rising edges of `VT("/out")` and `VT("/in")` when the waveforms cross their respective threshold values (that is, 2.5).

```plaintext
delay(VT("/out") 1.5 1 "rising" VT("/out") 1.5 2 "rising" 1 1 t "trigger") (s)
```
Returns multiple occurrences of delay specified against trigger time-points at which each delay event occurs.

delay(VT("/out") 1.5 1 "rising" VT("/out") 1.5 2 "rising" 1 1 t "target") (s)

Returns multiple occurrences of delay specified against target time-points at which each delay event occurs.

delay(VT("/out") 1.5 1 "rising" VT("/out") 1.5 2 "rising" 1 1 t "cycle") (s)

Returns multiple occurrences of delay specified against cycle numbers, where a cycle number refers to the n'th occurrence of the delay event in the input waveform.
deriv

deriv( o_waveform )

=> o_waveform/nil

Description

Computes the derivative of a waveform with respect to the X axis.

Note the following:

- After the second derivative, the results become inaccurate because the derivative is obtained numerically.
- Use the magnitude value instead of dB in frequency domain.

Arguments

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

Value Returned

o_waveform Returns a waveform object representing the derivative with respect to the X axis of the input waveform. Returns a family of waveforms if the input argument is a family of waveforms.

nil Returns nil and an error message otherwise.

Example

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

Plots the waveform representing the derivative of the voltage of "/net8".
**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:

\[
\text{from, from + deltaT, from + 2 * deltaT, ...}, \\
\text{from + (N - 1) * deltaT}
\]

The output of \( \text{dft} \) 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, \( (\text{from} + (N - 1) \times \text{deltaT}) \), is \( (\text{to} - \text{deltaT}) \) rather than \( \text{to} \). The \( \text{dft} \) command assumes that \( w(\text{from}) \) equals \( w(\text{to}) \).

**Arguments**

- **o_waveform**
  Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: \text{srrWave:XXXXX}.)

- **n_from**
  Starting value for the \( \text{dft} \) computation.

- **n_to**
  Ending value for the \( \text{dft} \) computation.

- **x_num**
  Number of timepoints. If \( x_{num} \) is not a power of 2, it is forced to be the next higher power of 2.

- **t_windowName**
  Variable representing different methods for taking a \( \text{dft} \) computation.
  Valid values: \text{Rectangular, ExtCosBell, HalfCycleSine, Hanning or Cosine2, Triangle or Triangular, Half3CycleSine or HalfCycleSine3, Hamming, Cosine4, Parzen, Half6CycleSine or HalfCycleSine6, Blackman, or Kaiser.}
For more information about `windowName`, see the information about Discrete Fourier Transform (dft) in the *Virtuoso Analog Design Environment L 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.
### 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**

- **o_waveform1**: Time domain waveform object with units of volts or amps.
- **o_waveform2**: Time domain waveform object with units of volts or amps.
- **f_timeStart**: Start time for the spectral analysis interval. Use this parameter and **f_timeEnd** to exclude part of the interval. For example, you might set these values to discard initial transient data.
- **f_timeEnd**: End time for the spectral analysis interval.
- **x_num**: The number of time domain points to use. The maximum frequency in the Fourier analysis is directly proportionate to **x_num** and inversely proportional to the difference between **f_timeStart** and **f_timeEnd**.
- **t_windowName**: 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.
- **x_smooth**: The Kaiser window smoothing parameter. If there are no requests, there is no smoothing. **Valid values**: 0 <= x_smooth <= 15. **Default value**: 1.
- **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

t_spectrumbType
A string that can be either singleSided or doubleSided. When this option is single-sided, the resultant waveform is only on one side of the y axis starting from 0 to N-1. When it is double-sided, it is symmetric to the Y axis from -N/2 to (N/2) -1.

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")
dnl

dnl( o_dacSignal o_sample|o_pointList|n_interval [?mode t_mode] [?threshold n_threshold] [?crossType t_crossType] [?delay f_delay] [?method t_method] [?units x_units] [?nbsamples n_nbsamples] )
=> n_dnl/nil

Description

Computes the differential non-linearity of a transient simple or parametric waveform.

Arguments

o_dacSignal  Waveform for which the differential non-linearity is to be calculated.

o_sample    Waveform used to obtain the points for sampling the dacSignal. These are the points at which the waveform crosses the threshold while either rising or falling (defined by the crossType argument) with the delay added to them.

n_pointList  List of domain values at which the sample points are obtained from the dacSignal.

n_interval   The sampling interval.

t_mode       The mode for calculating the threshold.
Valid values: auto and user.
Default value: auto.
If set to user, an n_threshold value needs to be provided.
If set to auto, n_threshold is calculated internally.

n_threshold  The threshold value against which the differential non-linearity is to be calculated. It needs to be specified only when the mode is selected as user.

t_crossType  The points at which the curves of the waveform intersect with the threshold. While intersecting, the curve may be either rising or falling.
Valid values: rising and falling, respectively.
Default crossType is rising.
**Ocean Reference**  
Predefined and Waveform (Calculator) Functions

\[ f_{\text{delay}} \]  
The delay time after which the sampling begins.  
Valid values: Any valid time value.  
Default value: 0.

\[ t_{\text{method}} \]  
The method to be used for calculation.  
Valid values: end (end-to-end) and fit (straight line).  
Default value: end.

\[ x_{\text{units}} \]  
The unit for expressing the output waveform.  
Valid values: abs (absolute) and lsb (multiples of least significant bit).  
Default value: abs.

\[ n_{\text{nbsamples}} \]  
The number of samples used for calculating the non-linearity. If not specified, the samples are taken against the entire data window.

**Note:** For each of the three ways in which the sample points can be specified, only a few of the other optional arguments are meaningful, as indicated below:

- For \( o_{\text{sample}} \), the arguments \( t_{\text{mode}}, n_{\text{threshold}}, t_{\text{crossType}}, f_{\text{delay}}, t_{\text{method}}, \) and \( x_{\text{units}} \) are meaningful.
- For \( n_{\text{pointList}} \), the arguments \( t_{\text{method}} \) and \( x_{\text{units}} \) are meaningful.
- For \( n_{\text{interval}} \), the arguments \( t_{\text{method}}, x_{\text{units}}, \) and \( n_{\text{nbsamples}} \) are meaningful.

**Value Returned**

\[ n_{\text{dnl}} \]  
Returns the differential waveform.

\[ \text{nil} \]  
Returns \text{nil} and an error message otherwise.

**Example**

```plaintext
dnl( wave1 wave2 \( ?\text{crossType} \) "rising" \( ?\text{delay} \) 0.4 )  
=> srrWave:175051544
```

Returns the differential non-linearity for \( \text{wave1} \) by taking the points at which \( \text{wave2} \) crosses the internally calculated threshold while \( \text{rising} \) as the sample points and adding a delay of 0.4 to them.
dutyCycle

dutyCycle( o_waveform [?mode t_mode] [?threshold n_threshold] [?xName t_xName] [?outputType t_outputType] )
   => o_waveform/f_average/nil

Description

Computes the duty cycle for a given waveform as a function of time or cycle.

**Note:** Duty cycle is the ratio of the time for which the signal remains ‘high’ and the time period of the signal.

Arguments

- **o_waveform**: Waveform, expression, or a family of waveforms.
- **t_mode**: The mode for calculating the threshold.
  - Valid values: auto and user.
  - Default value: auto.
  - If set to user, an n_threshold value needs to be provided.
  - If set to auto, n_threshold is calculated internally.
- **n_threshold**: The threshold value. It needs to be specified only when the mode is selected as user.
- **t_xName**: The X-axis of the output waveform.
  - Valid values: time and cycle.
  - Default value: time.
- **outputType**: Type of output.
  - Valid values: average and plot.
  - If set to average, the output is an average value.
  - If set to plot, the output is a waveform.
  - In both the cases, the output is expressed in terms of a percentage.
  - Default value: plot.
Value Returned

- \textit{o\_waveform} \hspace{1cm} Returns a waveform that represents duty cycle as a function of time.

- \textit{f\_average} \hspace{1cm} Returns the average duty cycle value as a percentage.

- \textit{nil} \hspace{1cm} Returns \textit{nil} if the duty cycle cannot be calculated.

Example

dutyCycle( \textit{wave1} )
\hspace{1cm} \Rightarrow \textit{srrWave:175051552}

Returns the duty cycle as a function of time for the wave \textit{wave1}.

dutyCycle( \textit{wave1} ?\textit{outputType} "average" )
\hspace{1cm} \Rightarrow 52.1066

Returns the average (in percentage) of the duty cycle values for the wave \textit{wave1}. 
evmQAM

**Description**

Processes the I and Q waveform outputs from the transient simulation run to calculate the Error Vector Magnitude (EVM) for multi-mode modulations. The function plots the I versus Q scatterplot. EVM is a useful measurement to describe the overall signal amplitude and phase modulated signal quality. It is based on a statistical error distribution normalized from an ideal digital modulation. Quadrature Amplitude Modulation (QAM) is a typical modulation scheme where EVM is useful. The EVM is calculated by detecting the I and Q signal levels corresponding to the four possible I and Q symbol combinations and calculating the difference between the actual signal level and the ideal signal level.

**Note:** This function is not supported for families of waveforms.

**Arguments**

- **o_waveform**
  - The waveform for the I signal.
- **o_waveformQ**
  - The waveform for the Q signal.
- **n_tDelay**
  - The start time (a numerical value) for the first valid symbol. This can be obtained from the Waveform Viewer window by recording the time of the first minimum or first maximum (whichever is earlier) on the selected signal stream.
- **n_samplingT**
  - A sampling time (a numerical value) for the symbol. Each period is represented by a data rate. The data rate at the output is determined by the particular modulation scheme being used.
- **x_levels**
  - The modulation levels.
  - Valid values: 4, 16, 64, 256
  - Default value: 4
- **b_normalize**
  - An option to see the scatter plot normalized to the ideal values +1 and -1 (for example, when superimposing scatter plots from different stages in the signal flow, where the levels may be quite different but the you want to see relative degradation or improvement in the scatter). This option does not affect the calculation of the EVM number.
Valid values: nil, t  
Default value: t

Value Returned

{o_waveform} Returns a waveform object representing the EVM value computed from the input waveforms.

nil Returns nil and an error message if the function is unsuccessful.

Example

evmQAM( v("samp_out_Q"), v("samp_out_I") 1.5u, 181.81n, 4, t )

Calculates the EVM value for the modulation level 4 in normalized form.
evmQpsk

```c
evmQpsk( o_waveform1 o_waveform2 n_tDelay n_sampling b_autoLevelDetect
         n_voltage n_offset b_normalize )
=> o_waveform/nil
```

**Description**

Processes the I and Q waveform outputs from the transient simulation run to calculate the Error Vector Magnitude (EVM) and plot the I versus Q scatterplot. EVM is a useful measurement to describe the overall signal amplitude and phase modulated signal quality. It is based on a statistical error distribution normalized from an ideal digital modulation. Quadrature Phase Shift Keying (QPSK) is a typical modulation scheme where EVM is useful. The EVM is calculated by detecting the I and Q signal levels corresponding to the four possible I and Q symbol combinations and calculating the difference between the actual signal level and the ideal signal level.

**Note:** This function is not supported for families of waveforms.

**Arguments**

- `o_waveform1` The waveform for the I signal.
- `o_waveform2` The waveform for the Q signal.
- `n_tDelay` The start time for the first valid symbol. This can be obtained from the Waveform Viewer window by recording the time of the first minimum or first maximum (whichever is earlier) on the selected signal stream.
- `n_sampling` A period for the symbol. Each period is represented by a data rate. The data rate at the output is determined by the particular modulation scheme being used.
- `b_autoLevelDetect` An option to indicate that you want the amplitude (n_voltage) and DC offset (n_offset) to be automatically calculated. Amplitude is calculated by averaging the rectified voltage level of the signal streams and DC offset by averaging the sum of an equal number of positive and negative symbols in each signal stream. These values are used to determine the EVM value. If this value is set to nil, you must specify values for n_voltage and n_offset.
Valid values: `nil`, `true`
Default value: `true`

**n_voltage**

The amplitude of the signal.

**n_offset**

The DC offset value.

**b_normalize**

An option to see the scatter plot normalized to the ideal values +1 and -1 (for example, when superimposing scatter plots from different stages in the signal flow, where the levels may be quite different but you want to see relative degradation or improvement in the scatter). This option does not affect the calculation of the EVM number.

Valid values: `nil`, `true`
Default value: `nil`

### Value Returned

**o_waveform**

Returns a waveform object representing the EVM value computed from input waveforms.

nil

Returns nil and an error message if the function is unsuccessful.

### Example

evmQpsk( v("samp_out_Q"), v("samp_out_I") 1.5u, 181.81n, true, nil, nil, nil)

Calculates the EVM value when **b_autoLevelDetect** is set to `true`. In this case, no values are specified for **n_voltage** and **n_offset**.

evmQpsk( v("samp_out_Q"), v("samp_out_I") 1.5u, 181.81n, nil, 1.3, 0, nil)

Calculates the EVM value when **b_autoLevelDetect** is set to `nil`. In this case, values are specified for **n_voltage** and **n_offset**.
**eyeDiagram**

```plaintext
eyeDiagram ( o_waveform n_start n_stop n_period ?advOptions t_advOptions )
=> o_waveform/nil
```

**Description**

Returns an eye-diagram plot of the input waveform signal. It returns the waveform object of the eye-diagram plot. Using an advanced option, the function also calculates the maximum vertical and horizontal opening of the eye formed when the input waveform is folded by the specified period to form the eye.

**Arguments**

- **o_waveform**: Input waveform signal.
- **n_start**: The X-axis start value from where the eye-diagram plot is to begin.
- **n_stop**: The X-axis stop value where the eye-diagram plot is to terminate.
- **n_period**: The period after which the waveform is to be folded to form the eye.
- **t_advOptions**: The option to specify whether the vertical or horizontal opening of the eye is to be calculated.
  - **Valid values:** `vertical`, `horizontal`
  - **Default value:** `nil`

**Note:** If `t_advOptions` is specified, the function approximates vertical eye height and horizontal eye width to assume the symmetry of the eye. The function returns the most optimum results for single eye scenarios.

**Value Returned**

- **o_waveform**: Returns a waveform object representing the eye-diagram plot of the input waveform.
- **nil**: Returns `nil` and an error message otherwise.
Example
eyeDiagram( v("/out" ) 0n 500n 12.5n )

Returns a waveform that represents an eye-diagram plot.

eyeDiagram( v("/out" ) 0n 500n 12.5n ?advOptions "vertical" )

Calculates the maximum vertical opening of the eye that is formed when the input waveform is folded after 12.5n

eyeDiagram( v("/out" ) 0n 500n 12.5n ?advOptions "horizontal" )

Calculates the maximum horizontal opening of the eye that is formed when the input waveform is folded after 12.5n
**flip**

\[
\text{flip( o\_waveform )} \\
\quad \Rightarrow o\_waveform/\text{nil}
\]

**Description**

Returns a waveform with the X vector values negated.

**Arguments**

- **o\_waveform**
  
  Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave:XXXXX.)

**Value Returned**

- **o\_waveform**
  
  Returns a waveform object representing the input waveform mirrored about its Y axis. Returns a family of waveforms if the input argument is a family of waveforms.

- **nil**
  
  Returns nil and an error message otherwise.

**Example**

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

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

`fourEval( o_waveform n_from n_to n_by [?b_baseBand] )
=> o_waveform/nil`

**Description**

Evaluates the Fourier series represented by an expression.

This function is an inverse Fourier transformation and thus the inverse of the `dft` command. The `fourEval` function transforms the expression from the frequency domain to the time domain.

**Arguments**

- `o_waveform`: Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: `srrWave:XXXXX`.)
- `n_from`: Starting point on the X axis at which to start the evaluation.
- `n_to`: Increment.
- `n_by`: Ending point on the X axis.
- `b_baseBand`: Accepts boolean values `t` or `nil`. The default value is `nil`. When set to `t`, the function evaluates the baseband version of the inverse of the `dft` function by converting the unsymmetrical spectrum to a symmetrical one.

**Value Returned**

- `o_waveform`: Returns a waveform object representing the inverse Fourier transformation of the input waveform. Returns a family of waveforms if the input argument is a family of waveforms. Returns the baseband version of the inverse of the `dft` function if `baseBand` is set to `t`.
- `nil`: Returns `nil` 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.
freq

freq( o_waveform t_crossType [?threshold n_threshold] [?mode t_mode] 
    [?xName xName][b_histoDisplay][x_noOfHistoBins] 
) => o_outputWave/nil

Description

Computes the frequency of the input waveform(s) as a function of time or cycle.

Arguments

\( o\_\text{waveform} \)  
Waveform, expression, or a family of waveforms.

\( t\_\text{crossType} \)  
The points at which the curves of the waveform intersect with the threshold. While intersecting, the curve may be either rising or falling. For the \text{freq} function, you may specify the frequency to be calculated against either the rising points or the falling points by setting \text{crossType} to \text{rising} or \text{falling}, respectively. The default \text{crossType} is \text{rising}.

\( n\_\text{threshold} \)  
The threshold value against which the frequency is to be calculated. This needs to be specified only when the \text{mode} selected is \text{user}.

\( t\_\text{mode} \)  
The mode for calculating the threshold. This is \text{auto}, by default, in which case \( n\_\text{threshold} \) is calculated internally. It can alternatively be set to \text{user}, in which case, an \( n\_\text{threshold} \) value needs to be provided.

\( t\_\text{xName} \)  
The X-axis of the output waveform. The default value is \text{time} but \text{cycle} is also a valid value.

\( b\_\text{histoDisplay} \)  
When set to \text{t}, returns a waveform that represents the statistical distribution of the \text{riseTime} data in the form of a histogram. The height of the bars (bins) in the histogram represents the frequency of the occurrence of values within the range of \text{riseTime} data.

Valid values: \text{t} \, \text{nil}
Default value: \text{nil}

\( x\_\text{noOfHistoBins} \)  
Denotes the number of bins represented in the histogram representation.
Valid values: Any positive integer
Default value: nil

Note: b_histoDisplay and x_noOfHistObins are added for backward compatibility only. It will be deprecated in future releases. Use the histo function for plotting the histogram of the resulting function.

Value Returned

\textit{o\_outputWave} \quad \text{Returns the frequency as a function of time or cycle.}

\textit{nil} \quad \text{Returns nil if the frequency cannot be calculated.}

Example

\texttt{freq( wave1 "rising" ?mode "user" ?threshold 18.5 ?xName "cycle" )}
\texttt{=> srrWave: 170938688}

Returns the frequency for \texttt{wave1} with the \texttt{threshold} at 18.5 against \texttt{cycle} on the x-axis.
freq_jitter

freq_jitter( o_waveform t_crossType [?mode t_mode] [?threshold n_threshold] [binSize n_binSize] [?xName t_xName] [?outputType t_outputType] )
=> o_waveform/f_val/nil

Description

Calculates the frequency jitter.

Arguments

- **o_waveform**: Waveform, expression, or a family of waveforms.
- **t_crossType**: The points at which the curves of the waveform intersect with the threshold. While intersecting, the curve may be either rising or falling.
  - Valid values: rising and falling.
  - Default value: rising.
- **t_mode**: The mode for calculating the threshold.
  - Valid values: auto and user.
  - If set to user, an n_threshold value needs to be provided.
  - If set to auto, n_threshold is calculated internally.
  - Default value: auto.
- **n_threshold**: The threshold value against which the frequency is to be calculated. It needs to be specified only when the mode selected is user.
- **n_binSize**: The width of the moving average window. The deviation of value at the particular point from the average of this window is the jitter.
- **t_xName**: The X-axis of the output waveform.
  - Valid values: time and cycle.
  - Default value: time.
- **t_outputType**: Type of output.
  - Valid values: sd and plot.
  - If set to sd, the output is a standard deviation jitter.
  - If set to plot, the output is a waveform.
  - Default value: plot.
Value Returned

\texttt{o\_waveform} \quad Returns the frequency jitter values as a function of time or cycle when the \texttt{outputType} is set to \texttt{plot}.

\texttt{f\_val} \quad Returns the standard deviation value when the \texttt{outputType} is set to \texttt{sd}.

\texttt{nil} \quad Returns \texttt{nil} otherwise.

Example

\begin{verbatim}
freq\_jitter( \texttt{wave1 “rising” ?mode “user” ?threshold 1 ?binSize 2 ?xName “cycle” \texttt{?outputType “sd”} } 
=> 0.1338585
\end{verbatim}

Returns the standard deviation for the frequency jitter of \texttt{wave1} with the threshold of 1 against the cycle on the x-axis.
**frequency**

\[
\text{frequency}( \text{o\_waveform} ) \\
\Rightarrow \text{o\_waveform}/n_{\text{value}}/\text{nil}
\]

**Description**

Computes the reciprocal of the average time between two successive midpoint crossings of the rising waveform.

**Arguments**

- **o\_waveform**
  - Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: \text{srrWave:XXXXX}.)

**Value Returned**

- **o\_waveform**
  - Returns a waveform representing the frequency of a family of waveforms if the input argument is a family of waveforms.

- **n\_value**
  - Returns a number representing the frequency of the specified waveform.

- **nil**
  - Returns \text{nil} and an error message otherwise.

**Example**

\[
\text{frequency}( \text{v( "/net12" ) } )
\]

Returns the frequency of "/net12".
ga

ga( o_s11 o_s12 o_s21 o_s22 [ ?gs n_gs] )
   => o_waveform/nil

Description

Returns the available gain in terms of the supplied parameters and the optional source reflection coefficient (Gs).

Arguments

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

Value Returned

o_waveform  Waveform object representing the available gain.
nil          Returns nil and an error message otherwise.

Example

s11 = sp(1 1)
s12 = sp(1 2)
s21 = sp(2 1)
s22 = sp(2 2)
plot(ga(s11 s12 s21 s22))
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.
Value Returned

\texttt{o\_waveform} \hspace{1cm} \text{Waveform object representing the available gain circles.}

\texttt{nil} \hspace{1cm} \text{Returns \texttt{nil} and an error message otherwise.}

Example

\begin{verbatim}
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))
\end{verbatim}
**gainBwProd**

gainBwProd( o_waveform )
=> o_waveform/n_value/nil

**Description**

Calculates the gain-bandwidth product of a waveform representing the frequency response of interest over a sufficiently large frequency range.

Returns the product of the zero-frequency-gain and 3dB-gain-frequency.

\[
gainBwProd(\text{gain}) = A_o \times f_2
\]

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**

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

**Value Returned**

- **o_waveform**: Returns a waveform representing the gain-bandwidth product for a family of waveforms if the input argument is a family of waveforms.
- **n_value**: Returns a value for the gain-bandwidth product for the specified waveform.
- **nil**: Returns nil and an error message otherwise.
Example

gainBwProd( v( "/OUT" ) )

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

\[
gainMargin(\ o\_\text{waveform}[\ b\_\text{stable}])
\]

\[
=>\ o\_\text{waveform}/n\_\text{value}/\text{nil}
\]

**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.

\[
gainMargin(\ \text{gain}) = 20 \times \log_{10}(\ \text{value}(\ \text{gain} \ f0))
\]

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**

- **o_waveform**
  Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: `srrWave:XXXXX`.)

- **b_stable**
  Boolean optional value that takes the value `nil` by default.

**Value Returned**

- **o_waveform**
  Returns a waveform representing the gain margin for a family of waveforms if the input argument is a family of waveforms.

- **n_value**
  Returns the value for the gain margin of the specified waveform.

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

**Example**

\[
gainMargin(\ v(\ "/OUT\ )\ ) = -9.234
\]
\[
gainMargin(\ v(\ "/OUT\ )\ \text{nil}\ ) = -9.234
\]
\[
gainMargin(\ v(\ "/OUT\ )\ \text{t}\ ) = 9.234
\]
gmax

\[
gmax( o\_s11 \ o\_s12 \ o\_s21 \ o\_s22 )
\Rightarrow o\_\text{waveform}/\text{nil}
\]

**Description**

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

**Arguments**

- \( o\_s11 \)
  - Waveform object representing s11.
- \( o\_s12 \)
  - Waveform object representing s12.
- \( o\_s21 \)
  - Waveform object representing s21.
- \( o\_s22 \)
  - Waveform object representing s22.

**Value Returned**

- \( o\_\text{waveform} \)
  - Load reflection coefficient.
- \( \text{nil} \)
  - Returns \( \text{nil} \) and an error message otherwise.

**Example**

\[
s11 = \text{sp}(1\ 1)
s12 = \text{sp}(1\ 2)
s21 = \text{sp}(2\ 1)
s22 = \text{sp}(2\ 2)
\]

\[
\text{plot}(\text{gmax}(s11\ s12\ s21\ s22))
\]
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:

\[
y_{Opt} = o_{Gopt} + (\text{complex } 0 1) \ast o_{Bopt} \\
\text{return } ( \frac{1}{f_{zref}(1) - y_{Opt}} ) / ( \frac{1}{f_{zref}(1) + y_{Opt}} )
\]

Arguments

- $o_{Gopt}$: Waveform object representing the optimum source conductance.
- $o_{Bopt}$: Waveform object representing the optimum source susceptance.
- $f_{zref}$: Reference impedance.

Value Returned

- $o_{gminWave}$: Waveform object representing the optimum noise reflection coefficient.
- nil: Returns nil and an error message otherwise.

Example

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

gmsg( o_s11 o_s12 o_s21 o_s22 )
=> o_waveform/nil

Description

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

Arguments

- **o_s11**: Waveform object representing s11.
- **o_s12**: Waveform object representing s12.
- **o_s21**: Waveform object representing s21.
- **o_s22**: Waveform object representing s22.

Value Returned

- **o_waveform**: Waveform object representing the maximum stable power gain.
- **nil**: Returns nil and an error message otherwise.

Example

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

\[
\text{gmux( } o_{s11} \ o_{s12} \ o_{s21} \ o_{s22} \ ) \\
\Rightarrow o_{\text{waveform}}/\text{nil}
\]

Description

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

Arguments

\begin{itemize}
\item \textit{o\_s11} \hspace{1cm} \text{Waveform object representing s11.}
\item \textit{o\_s12} \hspace{1cm} \text{Waveform object representing s12.}
\item \textit{o\_s21} \hspace{1cm} \text{Waveform object representing s21.}
\item \textit{o\_s22} \hspace{1cm} \text{Waveform object representing s22.}
\end{itemize}

Value Returned

\begin{itemize}
\item \textit{o\_waveform} \hspace{1cm} \text{Waveform object representing the maximum unilateral power gain.}
\item \textit{nil} \hspace{1cm} \text{Returns nil and an error message otherwise.}
\end{itemize}

Example

\begin{verbatim}
s11 = sp(1 1) 
s12 = sp(1 2) 
s21 = sp(2 1) 
s22 = sp(2 2) 
plot(gmux(s11 s12 s21 s22))
\end{verbatim}
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

<table>
<thead>
<tr>
<th>Argument</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>o_s11</td>
<td>Waveform object representing s11.</td>
</tr>
<tr>
<td>o_s12</td>
<td>Waveform object representing s12.</td>
</tr>
<tr>
<td>o_s21</td>
<td>Waveform object representing s21.</td>
</tr>
<tr>
<td>o_s22</td>
<td>Waveform object representing s22.</td>
</tr>
<tr>
<td>n_gl</td>
<td>Load reflection coefficient. Default value: 0</td>
</tr>
</tbody>
</table>

Value Returned

<table>
<thead>
<tr>
<th>Value Returned</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>o_waveform</td>
<td>Waveform object representing the power gain.</td>
</tr>
<tr>
<td>nil</td>
<td>Returns nil and an error message otherwise.</td>
</tr>
</tbody>
</table>

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>))
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 $g\text{\_level}$ and $g\text{\_frequency}$ allows either the level or the frequency to be swept while the other remains fixed.

Arguments

- **$o\text{\_s11}$**: Waveform object representing $s_{11}$.
- **$o\text{\_s12}$**: Waveform object representing $s_{12}$.
- **$o\text{\_s21}$**: Waveform object representing $s_{21}$.
- **$o\text{\_s22}$**: Waveform object representing $s_{22}$.
- **$g\text{\_level}$**: 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 $g\text{\_level}$ argument can be specified as either. In that case, a power gain circle is calculated at each one of the 13 levels.
- **$g\text{\_frequency}$**: The frequency. It can be specified as a scalar or a linear range. If it is specified as a linear range, the frequency is swept. The linear range is specified as a list with three values: the start of the range, the end of the range, and the increment. For example, `list(100M 1G 100M)` specifies a linear range with the following values:

  $$\{ 100\text{M}, 200\text{M}, 300\text{M}, 400\text{M}, 500\text{M}, 600\text{M}, 700\text{M}, 800\text{M}, 900\text{M}, 1\text{G} \}$$

In that case, a power gain circle is calculated at each one of the 10 frequencies.
### Value Returned

<table>
<thead>
<tr>
<th>Value</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td><code>o_waveform</code></td>
<td>Waveform object representing the power gain circles.</td>
</tr>
<tr>
<td><code>nil</code></td>
<td>Returns <code>nil</code> and an error message otherwise.</td>
</tr>
</tbody>
</table>
groupDelay

\[
groupDelay( o\_waveform ) \\
\Rightarrow o\_waveform/nil
\]

Description

Computes the group delay of a waveform.

This command returns the derivative of the phase of \( o\_waveform / 2\pi \). 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 \( v_p \) 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: \texttt{srrWave:XXXXX}.)

Value Returned

- \( o\_waveform \)
  - Returns a waveform representing the group delay of the specified waveform.

- nil
  - Returns nil and an error message otherwise.

Example

\[
\text{plot( groupDelay( v( "/net3" ) ) ) }
\]

Plots the waveform representing the group delay of the voltage of "/net3".
gt

gt( o_s11 o_s12 o_s21 o_s22 [ ?gs n_gs] [ ?gl n_gl] )
  => o_waveform/nil

Description

Returns the transducer gain in terms of the supplied parameters and the optional source reflection coefficient (Gs) and the input reflection coefficient (Gl).

Arguments

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

Value Returned

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

Example

  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>))
**harmonic**

harmonic( o_waveform h_index ) => o_waveform/g_value/nil

**Description**

Returns the waveform for a given harmonic index.

**Arguments**

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

- **h_index** The index number that designates the harmonic information to be returned. For the ‘pss’, ‘pac’, and ‘pxf’ analyses, the index is an integer number. For the ‘pdisto’ analysis, the index is a list of integers that correspond with the frequency names listed in the funds analysis parameter in the netlist. If more than one h_index is desired at one time, a list can be specified.

**Value Returned**

- **o_waveform** Returns a waveform (when a single h_index is specified) or family of waveforms (when more than one h_index is specified) if the input argument is a family of waveforms.

- **g_value** Returns the harmonic value if the input is a single waveform with the X values being harmonics

- **nil** Returns nil and displays a message if there is an error.

**Example**

For each of the following commands:

```
harmonic(v("/net49" ?result "pss-fd.pss") 1)
harmonic(v("/Pif" ?result "pdisto-fi.pdisto") list(1 -1))
```

Each result is a complex number.
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 "sweepdisto_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 "sweepdisto_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 2)))
```

Each resulting waveform is not in a useful format.
**harmonicFreqList**

```
harmonicFreqList( [?resultsDir t_resultsDir] [?result S_resultName])
    => n_list/nil
```

**Description**

Returns a list of lists, with each sublist containing a harmonic index and the minimum and maximum frequency values that the particular harmonic ranges between.

If both of these frequency values are the same, just one frequency value is returned.

**Arguments**

- **t_resultsDir**
  Directory containing the PSF files (results). If you supply this argument, you must also supply the `resultName` argument.

- **S_resultName**
  Results from an analysis.

**Value Returned**

- **n_list**
  Returns a list of lists. For the ‘pss, ‘pac, and ‘pxf analyses, the first element of each sublist is an integer number. For the ‘pdisto analysis, the first element of each sublist is a list of integers that correspond with the frequency names listed in the `funds analysis` parameter in the netlist. For all sublists, the remaining entries are the minimum and maximum frequency values that the particular harmonic ranges between. If both of these frequency values are the same, just one frequency value is returned.

- **nil**
  Returns nil if no harmonics are found in the data.

**Example**

For each of the following commands:

```
harmonicFreqList( ?result "pss-fd.pss" )
harmonicFreqList( ?result "pac-pac" )
harmonicFreqList( ?result "sweeppss_pss_fd-sweep" )
harmonicFreqList( ?result "sweeppss_pac-sweep" )
```
Each result is a list of integers.

For each of the following commands:

```plaintext
harmonicFreqList( ?result "pdisto-fi.pdisto" )
harmonicFreqList( ?result "sweeppdisto_pdisto_fi-sweep" )
```

Each result is a list of lists, with each sublist containing a combination of integer numbers that correspond with the frequency names listed in the `funds` analysis parameter in the netlist. These names can also be extracted from the PSF data by using the `resultParam` function to find the `largefundname` and `moderatefundnames` values. For example:

```plaintext
strcat(resultParam('largefundname ?result "pdisto-fi.pdisto" ) " "
resultParam('moderatefundnames ?result "pdisto-fi.pdisto" ))
```

Returns a string representing the order of the frequency names.
harmonicList

harmonicList( [?resultsDir t_resultsDir] [?result S_resultName] ) => n_list

Description

Returns the list of harmonic indices available in the resultName or current result data.

Arguments

- **t_resultsDir**: Directory containing the PSF files (results). If you supply this argument, you must also supply the resultName argument.
- **S_resultName**: Results from an analysis.

Value Returned

- **n_list**: Returns a list of harmonic indices. For the ‘pss’, ‘pac’, and ‘pxf’ analyses, the index is an integer number. For the ‘pdisto’ analysis, the index is a list of integers that correspond with the frequency names listed in the ‘funds analysis parameter in the netlist.
  - **nil**: Returns nil if no harmonics are found in the data.

Example

For each of the following commands:

- harmonicList( ?result "pss-fd.pss" )
- harmonicList( ?result "pac-pac" )
- harmonicList( ?result "sweeppss_pss_fd-sweep" )
- harmonicList( ?result "sweeppss_pac-sweep" )

Each result is a list of integers.

For each of the following commands:

- harmonicList( ?result "pdisto-fi.pdisto" )
- harmonicList( ?result "sweepdisto_pdisto_pdisto_fb-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:

```plaintext
strcat(resultParam( 'largefundname ?result "pdisto-fi.pdisto" ) " "
resultParam( 'moderatefundnames ?result "pdisto-fi.pdisto" ))
```

Returns a string representing the order of the frequency names.
**histo**

```plaintext
histo( o_waveform x_bins n_min n_max )
   => o_histoWaveform/nil
```

**Description**

Returns a waveform that represents the statistical distribution of input data in the form of a histogram. The height of the bars (or bins) in the histogram represents the frequency of the occurrence of values within a specific period. Using the `histo` function, the range for capturing these frequencies can be specified through the `n_min` and `n_max` values.

**Arguments**

- **o_waveform**  
  Input waveform.
- **x_bins**  
  Number of bins to represent the input data.
- **n_min**  
  The first value on the horizontal axis of the histogram. By default, it assumes the minimum value of the input waveform.
- **n_max**  
  The last value on the horizontal axis of the histogram. By default, it assumes the maximum value of the input waveform.

**Value Returned**

- **o_histoWaveform**  
  Returns a waveform representing the statistical distribution of the input waveform `o_waveform`.
- **nil**  
  Returns `nil` in case of an error.

**Example**

```plaintext
histo( VT("/vin") 3 1.5 3.5)
   => out_wave
plot( out_wave )
```

Plots the output waveform `out_wave` as a histogram, which represents the statistical distribution of the input waveform `VT("/vin")`. 
**iinteg**

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

**Description**

Computes the indefinite integral of a waveform with respect to the X-axis variable.

**Arguments**

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

**Value Returned**

- `o_waveform` Returns a waveform representing the indefinite integral of the input waveform.
- `nil` Returns `nil` and an error message otherwise.

**Example**

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

Computes the indefinite integral of the waveform representing the voltage of "/net8".
imag

imag( {o_waveform | n_input} )
=> o_waveformImag/n_numberImag/nil

Description

Returns the imaginary part of a waveform representing a complex number or returns the imaginary part of a complex number.

Arguments

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

n_input Complex number.

Value Returned

o_waveformImag Returns a waveform when the input argument is a waveform.

n_numberImag Returns a number when the input argument is a number.

nil Returns nil and an error message otherwise.

Example

imag( v( "/net8" ) )

Returns a waveform representing the imaginary part of the voltage of "/net8". You also can use the vim alias to perform the same command, as in

vim( "net8" ).

x=complex( -1 -2 ) => complex(-1, -2)
imag( x ) => -2.0

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

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

  initial_limit_n    Initial limit for definite integration.

  final_limit_n    Final limit for definite integration.

Value Returned

  o_waveform    Returns a waveform representing the definite integral for a family of waveforms if the input argument is a family of waveforms.

  n_value    Returns a numerical value representing the definite integral of the input waveform if the input argument is a single waveform.

  nil    Returns nil and an error message otherwise.

Example

integ( v( "/out" ) )

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

integ( VT( "/out" ),12.5n,18n)
Returns the definite integral of the waveform representing the voltage of "/out" within a specified range.
**intersect**

intersect( o_waveform1 o_waveform2 )

=> o_wave/nil

**Description**

Returns a waveform containing the points of intersection for two waveforms passed as arguments.

**Arguments**

- **o_waveform1**
  Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave:XXXXX.)

- **o_waveform2**
  Additional waveform object.

**Value Returned**

- **o_wave**
  Returns a waveform containing the points of intersection for the two waveforms passed as arguments.

- **nil**
  Returns nil if the two waveforms are disjoint or overlap each other, and an error message, if the arguments to the function are not correct.

**Example**

intersect( VT("/inp1") VT("/inp2") )
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 nth-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

o_spurious  Waveform or number representing the spurious output power (in dBm).
o_reference Waveform or number representing the reference output power (in dBm).
f_ordspur   Order or slope of the spurious constant-slope power line. Default value: 3
f_ordref    Order or slope of the reference constant-slope power line. Default value: 1
f_epspur    Value (in dBm) used to indicate the point where the spurious constant-slope power line begins. If b_psweep is t, this value is the input power value of the point on the o_spurious waveform, otherwise this value is paired with the o_spurious value to define the point. This point should be in the linear region of operation. (If b_psweep is t, f_spspur defaults to the X coordinate of the first point of the o_spurious wave; if s_measure is ‘input, a number must be specified.)
**Ocean Reference**

Predefined and Waveform (Calculator) Functions

---

**f_epref**  
Value (in dBm) used to indicate the point where the reference constant-slope power line begins. If \( b_{psweep} \) is \( t \), this value is the input power value of the point on the \( o_{reference} \) waveform, otherwise this value is paired with the \( o_{reference} \) value to define the point. This point should be in the linear region of operation. (If \( b_{psweep} \) is \( t \), \( f_{epref} \) defaults to the X coordinate of the first point of the \( o_{reference} \) wave; if \( s_{measure} \) is \( 'input \), a number must be specified.)

**b_psweep**  
Boolean indicating that the input power to the circuit was a parametric sweep. The power sweep must both be in dBm and be performed at the lowest parametric level.  
Default value: \( t \)

**s_measure**  
Name indicating if measurement is to be input referred (\( 'input \)) or output referred (\( 'output \)).  
Default value: \( 'input \)

---

**Value Returned**

**o_waveform**  
Depending on setting of \( b_{psweep} \) and the dimension of the input waveforms, returns either a waveform or a family of waveforms.

**f_number**  
If \( o_{spurious} \) and \( o_{reference} \) are numbers or they are waveforms when \( b_{psweep} \) is \( t \), returns a number.

**nil**  
Returns \( nil \) and an error message otherwise.

---

**Example**

```
spurWave = db20(harmonic(wave signalHarmonic))
refWave = db20(harmonic(wave referenceHarmonic))
xloc = ipn( spurWave refWave 3.0 1.0 -25 -25 )
yloc = ipn( spurWave refWave 3.0 1.0 -25 -25 t "Output")
```

Computes the IP3 point for the given wave.

Each of the following examples returns an ip3 measurement.

```
ipn(db20(harmonic(v("/Pif" ?result "pss_fd") 9)))
  db20(harmonic(v("/Pif" ?result "pss_fd") 8)))
ipn(dbm(harmonic(spectralPower(v("/Pif" ?result "pss_fd")/50.0
  v("/Pif" ?result "pss_fd") 9)))
```

---
dbm(harmonic(spectralPower(v("/Pif" ?result "pss_fd") /50.0
v("/Pif" ?result "pss_fd")) 8)))

ipn(dbm(harmonic(spectralPower(v("/Pif" ?result "pss_fd")
/resultParam("rif:r" ?result "pss_td")
v("/Pif" ?result "pss_fd")) 9))
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))
ipnVRI

```latex
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 \( \text{dBm}(\text{spectralPower}(\text{v/r},v)) \) 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**

- `o_vport` Voltage across the output port. This argument must be a family of spectrum waveforms (1 point per harmonic), with the option of containing a parametric input power sweep (in dBm).

- `x_harmspur` Harmonic number of the spurious voltage contained in `o_vport`. When `o_iport` is specified, also applies to a current waveform contained in `o_iport`.

- `x_harmref` Harmonic index of the reference voltage contained in `o_vport`. When `o_iport` is specified, also applies to a current waveform contained in `o_iport`.

- `o_iport` Current into the output port. This argument must be a family of spectrum waveforms (1 point per harmonic), with the option of containing a parametric input power sweep (in dBm). When specified, power is calculated using voltage and current.

- `f_rport` Resistance into the output port. When specified and `o_iport` is nil, the output power is calculated using voltage and
resistance.
Default value: 50

\( f_{\text{ordspur}} \)
Order or slope of the spurious constant-slope power line.
Default value: 3

\( f_{\text{epoint}} \)
Value (in dBm) used to indicate the point where the spurious constant-slope power line begins. If \( b_{\text{psweep}} \) is \( t \), this value is the input power value of the point on the \( o_{\text{spurious}} \) waveform, otherwise this value is paired with the \( o_{\text{spurious}} \) value to define the point. This point should be in the linear region of operation.
Default value: If \( b_{\text{psweep}} \) is \( t \), the lowest input power value; if \( s_{\text{measure}} \) is ‘input’, a number must be specified.

\( b_{\text{psweep}} \)
Boolean indicating that the input power to the circuit was a parametric sweep. The power sweep must be in dBm and must be performed at the lowest parametric level.
Default value: \( t \)

\( f_{\text{epref}} \)
Value (in dBm) used to indicate the point where the reference constant-slope power line begins. If \( b_{\text{psweep}} \) is \( t \), this value is the input power value of the point on the \( o_{\text{reference}} \) waveform, otherwise this value is paired with the \( o_{\text{reference}} \) value to define the point. This point should be in the linear region of operation.
Default value: If \( f_{\text{epoint}} \) is not nil, \( f_{\text{epoint}} \); else if \( b_{\text{psweep}} \) is \( t \), the X coordinate of the first point of the \( o_{\text{reference}} \) wave; else if \( s_{\text{measure}} \) is ‘input’, a number must be specified.

\( f_{\text{ordref}} \)
Order or slope of the reference constant-slope power line.
Default value: 1

\( s_{\text{measure}} \)
Symbol indicating if measurement is to be input referred (‘input) or output referred (‘output).
Default value: ‘input

**Value Returned**

\( o_{\text{waveform}} \)
Depending on the setting of \( b_{\text{psweep}} \) and the dimension of input waveform(s), the ipnVRI function returns either a waveform or a family of waveforms.
**Ocean Reference**

Predefined and Waveform (Calculator) Functions

---

**f_number**

Depending on the setting of `b_psweep` and the dimension of input waveform(s), the `ipnVRI` function returns a number.

**nil**

Returns `nil` and an error message otherwise.

---

**Example**

Each of following examples returns an ip3 measurement:

```ruby
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"))
```
ipnVRICurves


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

- **o_vport**  
  Voltage across the output port. This argument must be a family of spectrum waveforms (1 point per harmonic), with the option of containing a parametric input power sweep (in dBm).

- **x_harmspur**  
  Harmonic index of the spurious voltage contained in o_vport. When o_iport is specified, also applies to a current waveform contained in o_iport.

- **x_harmref**  
  Harmonic index of the reference voltage contained in o_vport. When o_iport is specified, also applies to a current waveform contained in o_iport.

- **o_iport**  
  Current into the output port. This argument must be a family of spectrum waveforms (1 point per harmonic), with the option of containing a parametric input power sweep (in dBm). When specified, power is calculated using voltage and current.
**Ocean Reference**

**Predefined and Waveform (Calculator) Functions**

---

### Variables

**f_rport**

Resistance into the output port. When specified and `o_iport` is `nil`, the output power is calculated using voltage and resistance.

Default value: 50

**f_ordspur**

Order or slope of the spurious constant-slope power line.

Default value: 3

**f_epoint**

Value (in dBm) used to indicate the point where the spurious constant-slope power line begins. If `b_psweep` is `t`, this value is the input power value of the point on the `o_spurious` waveform, otherwise this value is paired with the `o_spurious` value to define the point. This point should be in the linear region of operation.

Default value: If `b_psweep` is `t`, the X coordinate of the first point of the `o_spurious` wave; otherwise a number must be specified.

**b_psweep**

Boolean indicating that the input power to the circuit was a parametric sweep. The power sweep must be in dBm and must be performed at the lowest parametric level.

Default value: `t`

**f_epref**

Value (in dBm) used to indicate the point where the reference constant-slope power line begins. If `b_psweep` is `t`, this value is the input power value of the point on the `o_reference` waveform, otherwise this value is paired with the `o_reference` value to define the point. This point should be in the linear region of operation.

Default value: If `f_epoint` is not `nil`, `f_epoint`; else if `b_psweep` is `t`, the X coordinate of the first point of the `o_reference` wave; else a number must be specified.

**f_ordref**

Order or slope of the reference constant-slope power line.

Default value: 1

---

### Value Returned

**o_waveform**

A family of waveforms that contains the spurious and reference tangent lines, and when `b_psweep` is `t`, contains the spurious and reference waveforms.

**nil**

Returns `nil` and an error message otherwise.
Example

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

```plaintext
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"))
```
**kf**

\[
\text{kf}(\ o\_s\_11\ o\_s\_12\ o\_s\_21\ o\_s\_22\ )
\]

\[
=>\ o\_\text{waveform}/\text{nil}
\]

**Description**

Returns the stability factor in terms of the supplied parameters.

**Arguments**

- **o\_s\_11**: Waveform object representing s11.
- **o\_s\_12**: Waveform object representing s12.
- **o\_s\_21**: Waveform object representing s21.
- **o\_s\_22**: Waveform object representing s22.

**Value Returned**

- **o\_\text{waveform}**: Waveform object representing the stability factor.
- **nil**: Returns nil if there is an error.

**Example**

\[
s11 = \text{sp}(1\ 1)
\]
\[
s12 = \text{sp}(1\ 2)
\]
\[
s21 = \text{sp}(2\ 1)
\]
\[
s22 = \text{sp}(2\ 2)
\]

\[
\text{plot(kf(s11\ s12\ s21\ s22))}
\]
\textbf{In}

\[ \ln( \{ o\_waveform \mid n\_number \} ) \]
\[ \Rightarrow o\_waveform/f\_number/\text{nil} \]

\textbf{Description}

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

\textbf{Arguments}

- \textit{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: \textit{srrWave:XXXXX}.)
- \textit{n\_number} Number.

\textbf{Value Returned}

- \textit{o\_waveform} Returns a waveform object representing the base-e (natural) logarithm of the input waveform if the input argument is a waveform object, or returns a family of waveforms if the input argument is a family of waveforms
- \textit{f\_number} Returns a number if the input argument is a number.
- \textit{nil} Returns \texttt{nil} and an error message otherwise.

\textbf{Example}

\[ \ln( v( \text{"/net9"} ) ) \]

Gets a waveform that is calculated as the natural logarithm of the input waveform.

\[ \ln(\text{ymax}(v(\text{"/net9"}))) \]

Gets a waveform that is calculated as the natural logarithm of the following: \texttt{ymax(v("/net9"))}.

\[ \ln(100) \]
\[ \Rightarrow 4.60517 \]

Gets the natural logarithm of 100.
**log10**

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

**Description**

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

**Arguments**

- **o_waveform**  
  Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: `srrWave:XXXXX`.)

- **n_number**  
  Number.

**Value Returned**

- **o_waveform**  
  Returns a waveform object if the input argument is a waveform object or returns a family of waveforms if the input argument is a family of waveforms.

- **n_number**  
  Returns a number that is calculated as the base-10 logarithm of the input number.

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

**Example**

```plaintext
log10( v( "/net9" ) )
```

Gets a waveform that is calculated as the base-10 logarithm of the input waveform.

```plaintext
log10( ymin( v( "/net9" ) ) )
```

Gets a waveform representing the base-10 logarithm of `ymin(v("/net9"))`.

```plaintext
log10( 100 )
=> 2.0
```

Gets the base-10 logarithm of 100, or 2.
**lsb**

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

**Description**

Computes the load 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. If it is specified as a linear range, the frequency is swept. The linear range is specified as a list with three values: the start of the range, the end of the range, and the increment. For example, `list(100M 1G 100M)` specifies a linear range with the following values:

  ```
  { 100M, 200M, 300M, 400M, 500M, 600M, 700M, 800M, 900M, 1G }
  ```

  In that case, a load stability circle is calculated at each one of the 10 frequencies.

**Value Returned**

- **o_waveform**: Waveform object representing the load stability circles.
- **nil**: Returns nil and an error message otherwise.

**Example**

```
plot(lsb(s11 s12 s21 s22 list(800M 1G 100M)))
```
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 \texttt{rshift} command.

Arguments

- \textit{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: \texttt{srrWave:XXXXX}.)
- \textit{n_delta} Value by which the waveform is to be shifted.

Value Returned

- \textit{o_waveform} Returns a waveform object representing the input waveform shifted to the left. Returns a family of waveforms if the input argument is a family of waveforms.
- \textit{nil} Returns \texttt{nil} and an error message otherwise.

Example

plot( lshift( v( "/net8" ) 30u ) )

Shifts the waveform representing the voltage of "/net8" to the left by 30u and plots the resulting waveform.
**mag**

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

**Description**

Gets the magnitude of a waveform or number.

**Arguments**

- **o_waveform**: Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: `srrWave:XXXXX`.)
- **n_number**: Number.

**Value Returned**

- **o_waveform**: Returns a waveform object if the input argument is a waveform object or returns a family of waveforms if the input argument is a family of waveforms.
- **n_number**: Returns a number if the input argument is a number.
- **nil**: Returns `nil` and an error message otherwise.

**Example**

`mag( v( "5" ) )`

Gets the magnitude of the waveform representing the voltage at net 5. You can also use the `vm` alias to perform the same command, as in `vm( "5" )`.

`mag( i( "VFB" ) )`

Gets the magnitude of the waveform representing current through the `VFB` component. You can also use the `im` alias to perform the same command, as in `im( "VFB" )`.

`mag( -10 ) => 10`

Returns the magnitude of `-10`. 
nc

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

Description

Computes the noise circles.

Arguments

<table>
<thead>
<tr>
<th>Argument</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>o_Fmin</td>
<td>Waveform object representing the minimum noise factor.</td>
</tr>
<tr>
<td>o_Gmin</td>
<td>Waveform object representing the optimum noise reflection.</td>
</tr>
<tr>
<td>o_rn</td>
<td>Waveform object representing the normalized equivalent noise resistance.</td>
</tr>
<tr>
<td>g_level</td>
<td>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 \texttt{linRg} function can be called to generate a linear range. For example, \texttt{linRg(-30, 30, 5)} is the same as \texttt{list(-30, -25, -20, -15, -10, -5, 0, 5, 10, 15, 20, 25, 30)} and the \texttt{g_level} argument can be specified as either of the above. In that case, a noise circle is calculated at each one of the 13 levels.</td>
</tr>
<tr>
<td>g_frequency</td>
<td>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, \texttt{list(100M, 1G, 100M)} specifies a linear range with the following values: {100M, 200M, 300M, 400M, 500M, 600M, 700M, 800M, 900M, 1G} In that case, a noise circle is calculated at each one of the 10 frequencies.</td>
</tr>
</tbody>
</table>

Value Returned

<table>
<thead>
<tr>
<th>Value</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>o_waveform</td>
<td>Waveform object representing the noise circles.</td>
</tr>
</tbody>
</table>
nil

Returns nil and an error message otherwise.

Example

Gopt = getData("Gopt")
Bopt = getData("Bopt")
Zref = zref(1 ?result "sp")
Gmin = gmin(Gopt Bopt Zref)
Fmin = getData("Fmin")
rn = getData("NRR")
NC = nc(Fmin Gmin rn 10 list(100M 1G 100M))
displayMode("smith")
smithType("impedance")
plot(NC)
overshoot

```plaintext
overshoot( o_waveform n_initVal g_initType n_finalVal g_finalType [b_multiple [s_Xname]] [b_histoDisplay] [x_noOfHistoBins] )
=> o_waveform/n_value/nil
```

**Description**

Computes the percentage by which an expression overshoots a step going from the initial value to the final value you enter.

This command returns the overshoot of `o_waveform` as a percentage of the difference between the initial value and the final value.

In the equation below, M represents Maximum Value of the peak wave, F represents Final Value of the settled wave, and I represents Initial Value of the wave.

![Overshoot Diagram](image)

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

**Arguments**

- **o_waveform**
  Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: `srrWave:XXXXX`.)

- **n_initVal**
  Initial X value at which to start the computation.

- **g_initType**
  Specifies how `initVal` functions. Valid values: a non-nil value specifies that the initial value is taken to be the value of the waveform, interpolated at `initVal`, and the waveform is clipped from below, as follows:

\[
\text{o_waveform} = \text{clip}( \text{o_waveform initVal nil} )
\]
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.)

\textbf{\textit{n\_finalVal}}

Final value at which to end the computation.

\textbf{\textit{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:

\begin{align*}
\text{o\_waveform} & = \text{clip} ( \text{o\_waveform nil finalVal} )
\end{align*}

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.)

\textbf{\textit{b\_multiple}}

An optional boolean argument that takes the value nil by default. If set to t, the function returns multiple occurrences of the overshoot event.

\textbf{\textit{s\_xName}}

An optional argument that is used only when \textit{b\_multiple} is set to t. It takes the value time by default. It controls the contents of the x vector of the waveform object returned by the function. Valid values: ‘time, ‘cycle

\textbf{\textit{b\_histoDisplay}}

When set to t, returns a waveform that represents the statistical distribution of the riseTime data in the form of a histogram. The height of the bars (bins) in the histogram represents the frequency of the occurrence of values within the range of riseTime data. Valid values: t nil

Default value: nil

\textbf{\textit{x\_noOfHistoBins}}

Denotes the number of bins represented in the histogram representation. Valid values: Any positive integer

Default value: nil

\textbf{Note:} \textit{b\_histoDisplay} and \textit{x\_noOfHistoBins} are added for backward compatibility only. It will be deprecated in future releases. Use the histo function for plotting the histogram of the resulting function.
Value Returned

**o_waveform**

Returns a waveform (or family of waveforms) representing the amount of overshoot in comparison to the whole signal if the input argument is a family of waveforms or if $b\_multiple$ is set to $t$.

**n_value**

Returns a value for the amount of overshoot in comparison to the whole signal if the input is a single waveform.

**nil**

Returns nil and an error message otherwise.

Example

overshoot( v( "/net8" ) 7n t 3.99u t )

Returns the value of the overshoot for the waveform representing the voltage of "/net8".

overshoot(VT("/out") 0.5 nil 4.95 nil 5 t 'time)

Returns multiple occurrences of overshoot specified against time-points at which each overshoot event occurs.

overshoot(VT("/out") 0.5 nil 4.95 nil 5 t 'cycle)

Returns multiple occurrences of overshoot specified against cycle numbers, where a cycle number refers to the n'th occurrence of the overshoot event in the input waveform.
**peak**

\[
\text{peak( } o\_\text{waveform } ?\text{from } f\_\text{from } ?\text{to } f\_\text{to } ?\text{xtol } f\_\text{xtol } ?\text{ytol } f\_\text{ytol } ) \Rightarrow o\_\text{waveform}/\text{nil}
\]

**Description**

Detects the peaks in the input waveform and returns the X and Y coordinates of these peak points in the form of a waveform.

**Note:** The function will not work for waveforms that comprise of complex numbers.

**Arguments**

- **o\_waveform**
  
  Input waveform.

- **f\_from**
  
  The initial point on the given waveform to start determining the peaks. By default, the first point of the waveform is the starting point.

- **f\_to**
  
  The final point on the given waveform up to which the peaks are to be determined. By default, the last point of the waveform is the end point.

- **f\_xtol**
  
  The distance on the X axis within which all peaks are to be filtered.
  
  Default: 0.0

- **f\_ytol**
  
  The distance on the Y axis within which all peaks are to be filtered.
  
  Default: 0.0

**Note:** If both \(f\_\text{xtol}\) and \(f\_\text{ytol}\) are specified, the filtering mechanism will operate as follows:

- The maximum peak is selected first.

- All adjacent peaks in the neighborhood of both \(f\_\text{xtol}\) in the X-axis direction and \(f\_\text{ytol}\) in the Y-axis direction are then filtered.

- Next, all the peaks in the rectangular window thus formed are filtered based on both \(f\_\text{xtol}\) and \(f\_\text{ytol}\).
If only one of $f_{xtol}$ or $f_{ytol}$ is specified, the peaks are filtered only in either the X-axis direction or the Y-axis direction, respectively.

**Value Returned**

- $o\_waveform$ Returns a waveform whose X and Y coordinates of the peaks are determined from the input waveform and the peaks are filtered based on the $f_{xtol}$ and $f_{ytol}$ criteria.
- nil Returns nil and an error message otherwise.

**Example**

```plaintext
peak( vt("/out") ?from 1n ?to 20n ?xtol 2n ?ytol 0.5)
```

Out of all the peaks in the region starting from 1n to 20n, the function returns a waveform comprising of some of these peaks that satisfy the criteria of $x\_tol\ (2n)$ and $y\_tol\ (0.5)$.
peakToPeak

peakToPeak( o_waveform ) => o_waveform/n_value/nil

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

Arguments

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

Value Returned

o_waveform Returns a waveform or a family of waveforms if the input argument is a family of waveforms.

n_value Returns the difference between the maximum and minimum values of a waveform if the input argument is a single waveform.

nil Returns nil and an error message otherwise.

Example

peakToPeak( v( "/net2" ) )

Returns the difference between the maximum and minimum values of the waveform representing the voltage of the "/net2" net.
**period_jitter**

`period_jitter( o_waveform t_crossType [?mode t_mode] [?threshold n_threshold] [binSize n_binSize] [?xName t_xName] [?outputType t_outputType] ) => o_waveform/f_val/nil`

**Description**

Computes the period jitter.

**Arguments**

- **o_waveform**
  Waveform, expression, or a family of waveforms.

- **t_crossType**
  The points at which the curves of the waveform intersect with the threshold. While intersecting, the curve may be either rising or falling.
  Valid values: *rising* and *falling*.
  Default value: *rising*.

- **t_mode**
  The mode for calculating the threshold.
  Valid values: *auto* and *user*.
  If set to *user*, an *n_threshold* value needs to be provided.
  If set to *auto*, *n_threshold* is calculated internally.
  Default value: *auto*.

- **n_threshold**
  The threshold value against which the frequency is to be calculated. It needs to be specified only when the *mode* selected is *user*.

- **n_binSize**
  The width of the moving average window. The deviation of value at the particular point from the average of this window is the jitter.

- **t_xName**
  The X-axis of the output waveform.
  Valid values: *time* and *cycle*.
  Default value: *time*.

- **t_outputType**
  Type of output.
  Valid values: *sd* and *plot*.
  If set to *sd*, the output is a standard deviation jitter.
  If set to *plot*, the output is a waveform.
  Default value: *plot*. 
Value Returned

\texttt{o\_waveform} \quad \text{Returns the period jitter values as a function of time or cycle when the \textit{outputType} is set to plot.}

\texttt{f\_val} \quad \text{Returns the standard deviation value when the \textit{outputType} is set to \texttt{sd}.}

\texttt{nil} \quad \text{Returns \texttt{nil} otherwise.}

Example

\begin{verbatim}
period_jitter( wave1 "rising" ?mode "user" ?threshold 1 ?binSize 2 ?xName "cycle" ?outputType "sd" )
=> 1.695467
\end{verbatim}

Returns the standard deviation for the period jitter of \texttt{wave1} with the threshold of 1 against the cycle on the x-axis.
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: srrWave:XXXXX.)

n_number
Number.

Value Returned

o_waveform
Returns a waveform object if the input argument is a waveform object or returns a family of waveforms if the input argument is a family of waveforms.

n_number
Returns a number if the input argument is a number.

nil
Returns nil and an error message otherwise.

Example

phase( v( "5" ) )

Gets the phase of the waveform representing the voltage at net 5. You can also use the vp alias to perform the same command, as in vp( "5" ).

phase( i( "VFB" ) )

Gets the phase of the waveform representing the current through the VFB component. You can also use the ip alias to perform the same command, as in ip( "VFB" ).

phase( -2.0 ) => 180.0

Gets the phase of −2.
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

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

n_number Number.

Value Returned

o_waveform Returns a waveform object representing the wrapped phase in degrees of the input waveform. Returns a family of waveforms if the input argument is a family of waveforms.

n_number Returns a number if the input argument is a number.

nil Returns nil and an error message otherwise.

Example

phaseDeg( v( "vout" ) )

Takes the input waveform, representing the voltage of the "vout" net, and returns the waveform object representing the wrapped phase in degrees.
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

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

n_number Number.

Value Returned

o_waveform Returns a waveform object representing the unwrapped phase in degrees of the input waveform. Returns a family of waveforms if the input argument is a family of waveforms.

n_number Returns a number if the input argument is a number.
nil Returns nil and an error message otherwise.

Example

phaseDegUnwrapped( v( "vout" ) )

Takes the input waveform, representing the voltage of the "vout" net, and returns the waveform object representing the unwrapped phase in degrees.
phaseMargin

phaseMargin( o_waveform )

=> o_waveform/n_value/nil

**Description**

Computes the phase margin of the loop gain of an amplifier.

You supply a waveform representing the loop gain of interest over a sufficiently large frequency range.

phaseMargin( gain ) = 180 + phase( value( gain f0 ) )

The phase margin is calculated as the difference between the phase of the gain in degrees at f0 and at -180 degrees. The frequency f0 is the lowest frequency where the gain is 1. For stability, the phase margin must be positive.

**Arguments**

- **o_waveform**
  Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave:XXXXX.)
Value Returned

\textbf{o\_waveform}  
Returns a waveform representing the phase margin of the loop gain of an amplifier for a family of waveforms if the input argument is a family of waveforms.

\textbf{n\_value}  
Returns the value (in degrees) equivalent to the phase margin of the input waveform.

\textbf{nil}  
Returns \texttt{nil} and an error message otherwise.

Example

\texttt{phaseMargin( v( "/OUT" ) )}

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

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

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

Arguments

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

  n_number          Number.

Value Returned

  o_waveform        Returns a waveform representing a discontinuous value (in radians) for the phase of the input waveform. Returns a family of waveforms if the input argument is a family of waveforms.

  n_number          Returns a number when the input argument is a number.

  nil               Returns nil and an error message otherwise.

Example

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

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

`phaseRadUnwrapped(o_waveform)`

=> o_waveform/nil

**Description**

Calculates the unwrapped (continuous) phase in radians of a waveform and returns a waveform.

**Arguments**

- **o_waveform**
  
  Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: `srrWave:XXXXX`.)

**Value Returned**

- **o_waveform**
  
  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.

- **nil**
  
  Returns nil and an error message otherwise.

**Example**

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

Returns the unwrapped phase of the waveform representing the voltage of the "OUT" net.
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

o_waveformBase          Waveform object to be used as the base for the expression.
o_waveformExpn          Waveform object to be used as the exponent for the expression.
n_numberBase            Number to be used as the base for the expression.
n_numberExpn            Number to used as the exponent for the expression.

Value Returned

o_waveform              Returns a family of waveforms if one of the input arguments is a family of waveforms or returns a waveform if one of the input arguments is a waveform (and none is a family).
n_result                Returns a number if both the input arguments are numbers.
nil                    Returns nil and an error message otherwise.

Example

pow( average( v( "/net9" ) ) 0.5 )

Gets the square root of the average value of the voltage at "/net9".

pow( 2 3 )
  => 8

Gets the value of 2 to the third power, or 8.

pow( -2 2 )
  => 4
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.
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

o_waveform
Time domain waveform object with units of volts or amps.

f_timeStart
Starting time for the spectral analysis interval. Use this parameter and f_timeEnd to exclude part of the interval. For example, you might set these values to discard initial transient data.

f_timeEnd
Ending time for the spectral analysis interval.

x_num
The number of time domain points to use. The maximum frequency in the Fourier analysis is proportional to x_num and inversely proportional to the difference between f_timeStart and f_timeEnd.
Default value: 512

t_windowName
The window to be used for applying the moving window FFT.
Default value: ‘Hanning

x_smooth
The Kaiser window smoothing parameter. The 0 value requests no smoothing.
Valid values: 0 <= x_smooth <= 15.
Default value: 1
**Ocean Reference**

Predefined and Waveform (Calculator) Functions

---

- **f_cohGain**
  A scaling parameter. A non-zero value scales the power spectral density by $1/(f_{\text{cohGain}})$.
  Valid values: $0 < f_{\text{cohGain}} < 1$ (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**

```plaintext
psd(VT("/net32" "/hm/test_bench/spectre/schematic"), 0, 16m, 12000,
    ?windowName ' Hanning, ?smooth 1, ?windowSize 256,
    ?detrending ' None, ?cohGain 1)
```
Consider applying this command to one of the waveforms in the following illustration.
The result is the following spectrum, which is displayed with a logarithmic vertical scale.
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

- **o_waveform1**: Time domain waveform object with units of volts or amps.
- **o_waveform2**: Time domain waveform object with units of volts or amps.
- **f_timeStart**: Starting time for the spectral analysis interval. Use this parameter and **f_timeEnd** to exclude part of the interval. For example, you might set these values to discard initial transient data.
- **f_timeEnd**: Ending time for the spectral analysis interval.
- **x_num**: The number of time domain points to use. The maximum frequency in the Fourier analysis is proportional to $x\_num$ and inversely proportional to the difference between $f\_timeStart$ and $f\_timeEnd$.
- **x_smooth**: The Kaiser window smoothing parameter. 0 requests no smoothing. Valid values: $0 \leq x\_smooth \leq 15$. Default value: 1
### f_cohGain

A scaling parameter. A non-zero value scales the power spectral density by $1/(f_{\text{cohGain}})$. 
Valid values: $0 < f_{\text{cohGain}} < 1$ (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

```scheme
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)
```
Consider applying this command to both of the waveforms in the following illustration.
The result is the following spectrum, which is displayed with a logarithmic vertical scale.
**pzbode**

\[
pzbode( f_c \ f_{\text{minf}} \ f_{\text{maxf}} \ x_{\text{npoints}} \ ?\text{poles} \ o_{\text{waveform1}} \ ?\text{zeros} \ o_{\text{wavefoem2}}) \\
=> \ o_{\text{waveform}}/\text{nil}
\]

**Description**

Calculates and plots the transfer function of a circuit from pole zero simulation data.

**Arguments**

- **f_c**
  - The transfer gain constant.
- **f_{minf}**
  - The minimum frequency for the bode plot.
- **f_{maxf}**
  - The maximum frequency for the bode plot.
- **x_{npoints}**
  - The frequency interval for the bode plot, in points per decade.
- **o_{waveform1}**
  - Poles from the dumped simulation data.
    - Default value: all
- **o_{waveform2}**
  - Zeros from the dumped simulation data.
    - Default value: all

**Value Returned**

- **o_{waveform}**
  - Waveform containing the x and y points of the transfer function.
  - The scale of the Y axis will be db20.
- **nil**
  - Returns nil and error message otherwise.

**Example**

\[
pzbode( 1.0 \ 1\text{M} \ 1\text{G} \ 20 \ ?\text{poles} \ \text{comple}xPole\text{Wave} \ ?\text{zeros} \ \text{comple}xZero\text{Wave} )
\]
pzfilter

pzfilter( [o_PoleWaveform] [o_ZeroWaveform] [?maxfreq t_maxfreq]
   [?reldist n_reldist] [?absdist n_absdist] [?minq n_minq] [?output_type
   o_output] )
=> o_waveform/nil

Description

Returns the filtered Pole and Zero waveforms.

**Note:** If you do not specify values for o_PoleWaveform and o_ZeroWaveform
arguments, you should have run pz analysis prior to using this function.

Arguments

- **o_PoleWaveform**  
  Input Pole waveform (complex points).
  Default value: Poles of the simulator pz-analysis dump

- **o_ZeroWaveform**  
  Input Zero waveform (complex points).
  Default value: Zeros of the simulator pz-analysis dump

- **t_maxfreq**  
  Maximum frequency.
  Default value: 1e10

- **n_reldist**  
  Relative distance to be considered while filtering.
  Default value: 0.05

- **n_absdist**  
  Absolute distance to be considered while filtering.
  Default value: 1e-6

- **n_minq**  
  Minimum q factor to be allowed while filtering.

- **o_output**  
  Specifies the type of the output. If this argument is not passed, the output is a family of waves with two child waveforms, representing poles and zeros respectively, with the real component of each waveform as the X values and the imaginary components as the Y values.
  Valid value: complexwave. The output is a family of waves with two child waves, both of which are complex and represent poles and zeros, respectively.
Value Returned

\textit{o\_waveform} \quad \text{Returns a family (waveform) of Pole and Zero waveforms.}

\textit{nil} \quad \text{Returns \textit{nil} otherwise.}

Example

\texttt{pzfilter( complexPoleWave complexZeroWave )}
\texttt{=> srrWave:175051584}

Returns a family of filtered Pole and Zero waveforms, which correspond to the sweep values of “Pole” and “Zero”, respectively.
real

real( {o_waveform | n_input} )
=> o_waveformReal/n_numberReal/nil

Description

Returns the real part of a waveform representing a complex number, or returns the real part of a complex number.

Arguments

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

-n_input: Complex number.

Value Returned

-o_waveformReal: Returns a waveform when the input argument is a waveform.

-n_numberReal: Returns a number when the input argument is a number.

-nil: Returns nil and an error message otherwise.

Example

real( v( "/net8" ) )

Returns a waveform representing the real part of the voltage of "/net8". You also can use the vr alias to perform the same command, as in vr( "net8").

x=complex( -1 -2 ) => complex(-1, -2)
real( x ) => -1.0

Creates a variable 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 [b_multiple [s_Xname][b_histoDisplay][x_noOfHistoBins] ] )
         => o_waveform/n_value/nil
```

### Description

Returns the rise time measured between \( \theta_1 \) (percent low) to \( \theta_2 \) (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`.

![Graph](image)

### Arguments

- **o_waveform**
  - Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: `srrWave:XXXXX`.)

- **n_initVal**
  - Initial value at which to start the computation.

- **g_initType**
  - Specifies how `n_initVal` functions.
  - Valid values: a non-nil value specifies that the initial value is taken to be the value of the waveform, interpolated at `n_initVal`, and the waveform is clipped from below as follows:
  - `o_waveform = clip( o_waveform g_initVal nil )`
where `nil` specifies that `n_initVal` is defined by the X value entered. (The command gets the Y value for the specified X value and uses that value for `n_initVal`.)

**n_finalVal**

Final value at which to end the computation.

**g_finalType**

Specifies how the `n_finalVal` argument functions.

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

\[
o_{\text{waveform}} = \text{clip}(o_{\text{waveform}} \text{ nil } n_{\text{finalVal}})
\]

where `nil` specifies that the `n_finalVal` argument is defined by the X value entered. (The command gets the Y value for the specified X value and uses that value for `n_finalVal`.)

**n_theta1**

Percent low.

**n_theta2**

Percent high.

**b_multiple**

An optional boolean argument that takes the value `nil` by default. If set to `t`, the function returns multiple occurrences of the riseTime event.

**s_xName**

An optional argument that is used only when `b_multiple` is set to `t`. It takes the value `time` by default. It controls the contents of the x vector of the waveform object returned by the function.

Valid values: `time`, `cycle`

**b_histoDisplay**

When set to `t`, returns a waveform that represents the statistical distribution of the riseTime data in the form of a histogram. The height of the bars (bins) in the histogram represents the frequency of the occurrence of values within the range of riseTime data.

Valid values: `t` `nil`

Default value: `nil`

**x_noOfHistoBins**

Denotes the number of bins represented in the histogram representation.

Valid values: Any positive integer

Default value: `nil`
Note: *b_histoDisplay* and *x_noOfHistoBins* are added for backward compatibility only. It will be deprecated in future releases. Use the *histo* function for plotting the histogram of the resulting function.

**Value Returned**

- **o_waveform**
  Returns a waveform representing the rise time for a family of waveforms if the input argument is a family of waveforms or if `b_multiple` is set to `t`.

- **n_value**
  Returns a value for the rise time if the input is a single waveform.

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

**Example**

```r
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:

```r
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.

```r
riseTime(VT("/out") 0.5 nil 4.5 nil 10 90 t "time") (s)
```

Returns multiple occurrences of `riseTime` specified against time-points at which each `riseTime` event occurs.

```r
riseTime(VT("/out") 0.5 nil 4.5 nil 10 90 t "cycle") (s)
```

Returns multiple occurrences of `riseTime` specified against cycle numbers, where a cycle number refers to the n'th occurrence of the `riseTime` event in the input waveform.
rms

rms( o_waveform )
    => o_waveform/n_value/nil

Description

Returns the root-mean-square value of a waveform.

Arguments

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

Value Returned

o_waveform Returns a waveform representing the root-mean-square value for a family of waveforms if the input argument is a family of waveforms.

n_value Returns a value for the root-mean-square value for the specified waveform if the input is a single waveform.

nil Returns nil and an error message otherwise.

Example

rms( v( "/out" ) )

Returns the root-mean-square value of the waveform representing the voltage of the "/out" net.
**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.
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

- \( o_{\text{waveform}} \)  
  Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: \( srrWave:XXXXX \).)

- \( n_{\text{rootVal}} \)  
  Y value of interest.

- \( x_n \)  
  Number that specifies which X value to return. If \( n \) equals 1, the first X value that crosses over the Y \( rootVal \) is returned. If \( n \) equals 2, the second X value that crosses over the Y \( rootVal \) is returned, and so on. If you specify a negative integer for \( n \), the X values that cross the \( rootVal \) are counted from right to left (from maximum to minimum). If you specify \( n \) as 0, the list of root values is returned.

Value Returned

- \( o_{\text{waveform}} \)  
  Returns a waveform if the input argument is a family of waveforms.

- \( n_{\text{value}} \)  
  Returns an X value when the input argument is a single waveform.

- \( l_{\text{value}} \)  
  Returns a list of all the root values when \( n \) is 0.

- nil  
  Returns nil and an error message otherwise.

Example

root( v( "vout" ), 1.0, 4 )
Returns the X value for the point at which the waveform curve crosses the 1.0 Y value for the fourth time.
rshift

rshift( o_waveform n_delta )
   => o_waveform/nil

Description

Shifts the waveform to the right by the \textit{n_delta} value.

This command is the inverse of the \texttt{lshift} command.

Arguments

\begin{itemize}
\item \texttt{o\_waveform} \hspace{2cm} Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: \texttt{srrWave:XXXXX}.)
\item \texttt{n\_delta} \hspace{2cm} Value by which the waveform is to be shifted.
\end{itemize}

Value Returned

\begin{itemize}
\item \texttt{o\_waveform} \hspace{2cm} Returns a waveform object. Returns a family of waveforms if the input argument is a family of waveforms.
\item \texttt{nil} \hspace{2cm} Returns \texttt{nil} and an error message otherwise.
\end{itemize}

Example

\texttt{rshift( v( "vout" ) 10n )}

Shifts the waveform representing the voltage through the \texttt{"vout"} net to the right by \texttt{10n}.

\begin{tikzpicture}
  \draw[->] (0,0) -- (5,0) node[right] {X};
  \draw[->] (0,-1) -- (0,1) node[above] {Voltage};
  \draw[very thick] (0,0) -- (1,0.5) -- (2,0) -- (3,-0.5) -- (4,0);
  \draw[very thick, dashed] (3,0) -- (4,0);
  \node at (1.5,0.25) {10n}; \node at (2.5,0.25) {20n};
  \node at (1.5,-0.25) {10n}; \node at (2.5,-0.25) {20n};
\end{tikzpicture}
sample

\[
sample( \text{o\_waveform}, \text{n\_from}, \text{n\_to}, \text{t\_type}, \text{n\_by} ) \\
\Rightarrow \text{o\_waveform}/\text{n\_number}/\text{nil}
\]

Description

Samples a waveform at the specified interval.

You can use this function to reduce the time it takes to plot waveforms that have many data points. If you sample a waveform beyond its range, you get the final value of the waveform. You can use this function to demodulate a signal. Consider an AM modulated sine wave. Assume the carrier frequency is 1 GHz, and the modulation frequency is 1 MHz. If the waveform is sampled every 1 ns, the resulting signal is cleanly demodulated (the 1 GHz carrier is completely eliminated by the sampling).

Note: The function can be used to sample both a waveform object as well as a family of waveforms. If the family is of dimension \( m \), the arguments \( n\_from, n\_to, \) and \( n\_by \) would be of dimension \( m-1 \).

Arguments

- **o\_waveform**
  - Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave:XXXXX.)

- **n\_from**
  - Starting value for the sampling.

- **n\_to**
  - Ending value for the sampling.

- **t\_type**
  - Type of the sampling.
  - Valid values: "linear" or "log"

- **n\_by**
  - Interval at which to sample.

Value Returned

- **o\_waveform**
  - Returns a waveform representing the sampling you specified.

- **n\_number**
  - Returns a number if the output contains only one point.

- **nil**
  - Returns nil and an error message otherwise.
Example

\texttt{sample( v( "vout" ) 0 50n "linear" 0.1n )}

Takes a linear sample of the waveform representing the voltage of the "vout" net.

\texttt{sample( v( "vout" ) 0 100m "log" 10 )}

Takes a logarithmic sample of the waveform representing the voltage of the "vout" net.
settlingTime

settlingTime( o_waveform n_initVal g_initType n_finalVal g_finalType n_theta
[b_multiple [s_Xname]] )
=> o_waveform/n_value,nil

Description

The settling time is the time by which the signal settles within the specified Percent of step (theta) of the difference between the Final Value and Initial Value from the Final Value.

Note: The above graph represents the Initial value of the signal as 0% and Final value as 100%. The Percent of Step is taken as 5%.

Arguments

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

- **n_initVal**
  Initial value at which to start the computation.

- **g_initType**
  Specifies whether the values entered are X values or Y values. Valid values: `t` specifies that `n_initVal` is defined by the X value entered; `nil` specifies that `n_initVal` is defined by the Y value entered.
**Ocean Reference**  
Predefined and Waveform (Calculator) Functions

---

### Final Value at Which to Start the Computation

- **n_finalVal**: Final value at which to start the computation.

### Specifies Whether the Values Entered Are X Values or Y Values

- **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.

### Percent of the Total Step

- **n_theta**: Percent of the total step.

### An Optional Boolean Argument That Takes the Value nil by Default

- **b_multiple**: An optional boolean argument that takes the value nil by default. If set to t, the function returns multiple occurrences of the settlingTime event.

### An Optional Argument That Is Used Only When b_multiple Is Set to t

- **s_xName**: An optional argument that is used only when `b_multiple` is set to t. It takes the value `time` by default. It controls the contents of the x vector of the waveform object returned by the function.  
  - Valid values: `time`, `cycle`

---

### Value Returned

- **o_waveform**: Returns a waveform representing the settling time for a family of waveforms if the input argument is a family of waveforms or if `b_multiple` is set to t.

- **n_value**: Returns a value for the settling time for the specified waveform if the input is a single waveform.

- **nil**: Returns nil and an error message otherwise.

---

### Example

- **settlingTime( v("/out") 0 t 2 t 90 )**

  Computes the time required for the waveform representing the voltage of the "/out" net to settle within 90 percent of the step from 0 to 2.

- **settlingTime(VT("/out") 0.5 nil 4.95 nil 5 t "time") (s)**

  Returns multiple occurrences of settlingTime specified against time-points at which each settlingTime event occurs.

- **settlingTime(VT("/out") 0.5 nil 4.95 nil 5 t "cycle") (s)**

---

August 2009  
421  
Product Version 6.1.3
Returns multiple occurrences of settlingTime specified against cycle numbers, where a cycle number refers to the n’th occurrence of the settlingTime event in the input waveform.
slewRate

\[
\text{slewRate( } o\_\text{waveform } n\_\text{initVal } g\_\text{initType } n\_\text{finalVal } g\_\text{finalType } n\_\theta_1 \\
\quad n\_\theta_2 \text{ [b\_multiple [s\_Xname]] [b\_histoDisplay] [x\_noOfHistoBins] } \\
\quad => o\_\text{waveform}/n\_value/nil
\]

Description

Computes the average rate at which an expression changes from \( \theta_1 \) (percent low) to \( \theta_2 \) (percent high) of the difference between the initial value and final value.

\[
slewRate = \frac{\Delta Y}{\Delta X}
\]

Arguments

- \( o\_\text{waveform} \): Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: \( \text{srrWave:XXXXX} \).)
- \( n\_\text{initVal} \): Initial X-axis value at which to start the computation.
- \( g\_\text{initType} \): Specifies whether the values entered are X values or Y values. Valid values: \( t \) specifies that \( \text{initVal} \) is defined by the X value entered; \( \text{nil} \) specifies that \( \text{initVal} \) is defined by the Y value entered
- \( n\_\text{finalVal} \): Final value at which to end the computation.
- \( g\_\text{finalType} \): Specifies whether the values entered are X values or Y values. Valid values: \( t \) specifies that \( \text{finalVal} \) is defined by the X value entered; \( \text{nil} \) specifies that \( \text{finalVal} \) is defined by the Y value entered
- \( n\_\theta_1 \): Percent low (percentage of the total step).
n_theta2

Percent high (percentage of the total step).

b_multiple

An optional boolean argument that takes the value nil by default. If set to t, the function returns multiple occurrences of the slewRate event.

s_xName

An optional argument that is used only when b_multiple is set to t. It takes the value time by default. It controls the contents of the x vector of the waveform object returned by the function.
Valid values: 'time, 'cycle

b_histoDisplay

When set to t, returns a waveform that represents the statistical distribution of the riseTime data in the form of a histogram. The height of the bars (bins) in the histogram represents the frequency of the occurrence of values within the range of riseTime data.
Valid values: t nil
Default value: nil

x_noOfHistoBins

Denotes the number of bins represented in the histogram representation.
Valid values: Any positive integer
Default value: nil

Note: b_histoDisplay and x_noOfHistoBins are added for backward compatibility only. It will be deprecated in future releases. Use the histo function for plotting the histogram of the resulting function.

Value Returned

o_waveform

Returns a waveform representing the slew rate for a family of waveforms if the input argument is a family of waveforms or if b_multiple is set to t.

n_value

Returns a value for the slew rate for the specified waveform if the input is a single waveform.

nil

Returns nil or an error message otherwise.

Example

slewRate( v( "vout" ) 10n t 30n t 10 90 )
Computes the slew rate for the waveform representing the voltage of the "vout" net from 10n to 30n.

\[ \text{slewRate}( v( "vout" ) 0 \text{ nil} 10 \text{ nil} 5 95 ) \]

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.

\[ \text{slewRate}(V T( "out" ) 0.5 \text{ nil} 4.5 \text{ nil} 10 90 \text{ t 'time}) \]

Return multiple occurrences of slewRate values, computed at different time-points.

\[ \text{slewRate}(V T( "out" ) 0.5 \text{ nil} 4.5 \text{ nil} 10 90 \text{ t 'cycle}) \]

Returns multiple occurrences of slewRate values specified against cycle numbers (where cycle number refers to the n'th occurrence of slewRate computation).
spectralPower

spectralPower( o_current o_voltage )
  => o_power/nil

Description

Returns the spectral power given the spectral current and voltage.

To obtain a list of the harmonic frequencies, use harmonicList.

Arguments

  o_current Waveform representing the current. The current can be obtained by calling the i data access function for the desired terminal.

  o_voltage Waveform representing the voltage. The voltage can be obtained by calling the v 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

  o_power Waveform representing the power of the net.

  nil Returns nil if there is an error.

Example

plot(db10(spectralPower(i("/PORT0/PLUS") v("/net28"))))

Plots power of the output "/net28". "/PORT0/PLUS" is a member of "/net28".
spectrum

spectrum( o_waveform x_numSamples x_noiseBins n_startFreq n_endFreq  
    t_windowName n_adcSpan t_measType )  
=> o_spectrumWaveform/g_value/nil

Description

Calculates Signal-to-Noise-and-Distortion Ratio (SINAD), Spurious Free Dynamic Range (SFDR), Effective Number of Bits (ENOB), and Signal-to-Noise Ratio (without distortion) by using discrete fourier transform of any given input signal.

The spectrum measure is used for characterizing A-to-D converters and is typically supported for transient simulation data.

Arguments

o_waveform  
Signal to measure.  
Valid values: Any transient signal or signal expression using calculator functions.  
A signal can be expressed as either a net or a terminal.

x_numSamples  
Optional number of sampled points used for the FFT.  
Valid values: Any integer power of two greater than zero. For a value that is not a power of two, the function rounds it up to the next closest power of two.  
Default value: Number of data points in the signal o_waveform.

x_noiseBins  
Optional number of noise bins, where the size of one bin is the reciprocal of the data window width. For example, 1 ms of transient data creates a bin size of 1 kHz.  
Valid values: Any integer power of two greater than or equal to zero.  
Default value: 0, implying that no signal is spilling into the bins.  
A frequency band of bin-size times the number of bins is calculated and adjusted as a function of the selected window.  
Frequency components in each band to the left and right of the fundamental or the harmonics are set to zero and do not contribute to any output result.
**n_startFreq**
Optional lower limit of frequency range for the spectrum measures.
Default value: First frequency point of the FFT.

**n_endFreq**
Optional upper limit of frequency range for the spectrum measures.
Default value: Last frequency point of the FFT.

**t_windowName**
Optional windowing function applied to o_waveform.
Default value: Rectangular.

**n_adcSpan**
Optional full-scale span, ignoring any DC offsets. This is used in ENOB calculation.
Valid values: Any floating point number.
Default value: If n_adcSpan is not specified or is nil, it is assumed to be 0 and is taken to be the peak-to-peak value of the fundamental.

**t_measType**
Result specifier.
Valid values: sinad, sfdr(db), enob, and snhr.

**Value Returned**

**o_spectrumWaveform**
Returns a waveform of spectrum measures if the input argument t_measType is specified as all.

**g_value**
Returns the spectrum measure specified by the t_measType argument.

**nil**
Returns nil and an error message otherwise.
Example

specMeas = spectrum( VT("/vcoOut") 1K nil 1K 10G "Rectangular" nil "all")
=> srrWave:180703344
value(specMeas "sinad")
=> -5.104-5.104
value(specMeas "sfdr")
=> 392.2m
value(specMeas "enob")
=> -1.14
value(specMeas "snhr")
=> -4.948

The variable specMeas represents the waveform returned by the spectrum function. The value function retrieves the value corresponding to a particular spectrum measure from specMeas.

specMeas = spectrum( VT("/vcoOut") 1K nil 1K 10G "Rectangular" nil "snhr")
=> -4.948

Returns the value of the spectrum measure snhr, as specified by the spectrum function.
**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 \( y = f(x) \)

\[
\int_{\text{from}}^{\text{to}} (y - \text{average}(y))^2
\]

\[
\text{stddev}(y) = \sqrt{\int_{\text{from}}^{\text{to}} (y - \text{average}(y))^2}
\]

Arguments

\( o\_waveform \)

Waveform object or family of waveforms representing simulation results that can be displayed as a series of points. (A waveform object identifier looks like this: srrWave:XXXXX)

Value Returned

\( n\_stddev \)

Returns a number representing the standard deviation value of the input waveform.

\( o\_waveformStddev \)

Returns a waveform representing the average value if the input is a family of waveforms.

\( \text{nil} \)

Returns \( \text{nil} \) or an error message.

Example

`stddev( v( "/net9" ) )`

Gets the standard deviation of the voltage (Y-axis value) of \( /\text{net9} \) over the entire time range specified in the simulation analysis.
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

\(o\_waveform\)  Waveform object representing the wave.

\(n_x\)  X coordinate of the point. The default value is the X coordinate of the first point on the wave.

\(n_y\)  Y coordinate of the point. The default value is the Y coordinate at the given or default X coordinate.

\(n\_slope\)  Slope of the line.
Default value: 1.0

Value Returned

\(o\_waveform\)  Wave object representing the line.

\(nil\)  Returns \(nil\) if there is an error.

Example

refLine
=> tangent(refWave ?x -25 ?slope 1.0)
thd

thd( o_waveform n_from n_to x_num n_fund)
   => o_waveform/n_thdValue/nil

Description

The thd function computes the percentage of total harmonic content of a signal with respect to the fundamental frequency expressed as a voltage percentage.

The computation uses the dft function. Assume that the dft function returns complex coefficients $A_0, A_1, ..., A_f, ...$. Please note that fundamental frequency $f$ is the frequency contributing to the largest power in the signal. $A_0$ is the complex coefficient for the DC component and $A_i$ is the complex coefficient for the $i$th harmonic where $i \neq 0, f$. Then, total harmonic distortion is computed as:

$$\sqrt{\frac{\sum_{i=1, i \neq 0, f} |A_i|^2}{|A_f|}} \times 100\%$$

Arguments

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

- **n_from**
  Starting value for the computation.

- **n_to**
  Ending value for the computation.

- **x_num**
  Number of timepoints. If $x_{num}$ is not a power of 2, it is forced to be the next higher power of 2.

- **n_fund**
  Fundamental Frequency of the signal. If it is nil or zero then the non-zero frequency contributing to the largest power in the signal is used as the fundamental frequency. Otherwise, the harmonic frequency nearest to its value is used as the fundamental frequency.
Value Returned

\texttt{o\_waveform} \quad Returns a waveform representing the absolute value of the total harmonic distortion if the input argument is a family of waveforms.

\texttt{n\_thdValue} \quad Returns the absolute value of the total harmonic distortion of the input waveform.

\texttt{nil} \quad Returns \texttt{nil} and an error message otherwise.

Example

\texttt{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.

\texttt{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.


### unityGainFreq

unityGainFreq( o_gainFreqWaveform )

=> n_frequency/nil

**Description**

Computes and reports the frequency at which the gain is unity.

**Arguments**

- **o_gainFreqWaveform**  
  Gain frequency waveform.

**Value Returned**

- **n_frequency**  
  Returns a scalar value representing the frequency at which the gain of the input waveform is unity.

- **nil**  
  Returns nil otherwise.

**Example**

unityGainFrequency( VF("/out") )
value

value( o_waveform [s_name] g_value ?period n_period [b_multiple [s_Xname]]
[b_histoDisplay][x_noOfHistoBins])
=> o_waveform/g_value/nil

Description

Returns the Y value of a waveform for a given X value.

Arguments

- **o_waveform**: Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave:XXXXX.)
- **s_name**: The name of the innermost or outermost sweep variable. If the sweep variable name is not supplied, the innermost sweep variable is used.
- **g_value**: Value (X value) at which to provide the Y value. If a string has been defined for a value or set of values, the string may be used instead of the value.
- **n_period**: The interval or period after which the value needs to be computed.
- **b_multiple**: An optional boolean argument that takes the value nil by default. If set to t, the function returns multiple occurrences of the interpolated value.
- **s_xName**: An optional argument that is used only when **b_multiple** is set to t. It takes the value time by default. It controls the contents of the x vector of the waveform object returned by the function.
  - **Valid values**: 'time, 'cycle
- **b_histoDisplay**: When set to t, returns a waveform that represents the statistical distribution of the riseTime data in the form of a histogram. The height of the bars (bins) in the histogram represents the frequency of the occurrence of values within the range of riseTime data.
Valid values: t nil  
Default value: nil

\texttt{x_noOfHistoBins} \hspace{1cm} Denotes the number of bins represented in the histogram representation.  
Valid values: Any positive integer  
Default value: nil

**Note:** \texttt{b_histoDisplay} and \texttt{x_noOfHistoBins} are added for backward compatibility only. It will be deprecated in future releases. Use the histo function for plotting the histogram of the resulting function.

For the simplest calls to the function, which specify only the given waveform (\texttt{o_waveform}) and the X value (\texttt{g_value}), the given waveform can be a family of waveforms. If the family is of dimension \(m\), \texttt{g_value} can be either of dimension \(m-1\) or a scalar. If \texttt{g_value} is scalar, the function returns the Y value of all the components of the family at the specified \texttt{g_value}.

**Value Returned**

\texttt{o_waveform} \hspace{1cm} Returns a waveform or a family of waveforms if the input argument is a family of waveforms or if values are expected at multiple points.

\texttt{g_value} \hspace{1cm} Returns the Y value if the input argument is a single waveform. For parametric sweeps, the value might be a single waveform that can be printed with the \texttt{ocnPrint} command.

\texttt{nil} \hspace{1cm} Returns \texttt{nil} and an error message if the value cannot be printed.

**Example**

\texttt{value( v( "/net18" ) 4.428e-05 )}

Prints the value of \texttt{"/net18"} at \texttt{time=4.428e-05}. This is a parametric sweep of temperature over time.

\texttt{value( v( "/OUT" )'TEMPDC 20.0 )}

Returns \texttt{srrWave:XXXXX}, indicating that the result is a waveform.

\texttt{print( value( v( "/OUT" )'TEMPDC 20.0 ) )}

Prints the value of \texttt{v( "/OUT" )} at every time point for \texttt{TEMPDC=20}. 
**Ocean Reference**  
Predefined and Waveform (Calculator) Functions

- **print( value( v("/OUT") 200n ?period 100n ) )**
  
  Prints the value of \( v("/OUT") \) at 200n, 300n and so on at intervals of 100n until the end of the waveform.

- **value(VT("/out") 2e-07 ?period 2e-07 ?xName "time") (V)**
  
  Returns multiple occurrences of the value specified against time-points at which each interpolated value occurs.

- **value(VT("/out") 2e-07 ?period 2e-07 ?xName "cycle") (V)**
  
  Returns multiple occurrences of value specified against cycle numbers, where a cycle number refers to the n'th occurrence of the value event in the input waveform.
xmax

xmax( o_waveform x_numberOfPeaks )
  => o_waveform/g_value/l_value/nil

Description

Computes the value of the independent variable (X) at which the Y value attains its maximum value.

Arguments

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

x_numberOfPeaks Specifies the nth X value corresponding to the maximum Y value. For example, if x_numberOfPeaks is 3, the X value corresponding to the third maximum Y value is returned. If you specify a negative integer for x_numberOfPeaks, the X values are counted from right to left (from maximum to minimum). If x_numberOfPeaks is 0, xmax returns a list of X locations.

Value Returned

o_waveform Returns a waveform (or a family of waveforms) if the input argument is a family of waveforms.

g_value Returns the X value corresponding to the peak specified with x_numberOfPeaks if the input argument is a single waveform.

l_value Returns a list of X locations when x_numberOfPeaks is 0 and the input argument is a single waveform.

nil Returns nil and an error message otherwise.

Example

xmax( v( "/net9" ) 1 )
Ocean Reference
Predefined and Waveform (Calculator) Functions

 Gets the time value (X-axis value) at which the voltage of "/net9" attains its first peak value.
\[ \text{xmax}( v( "/net9" ) 0 ) \]

 Gets the list of time values (X-axis values) at which the voltage of "/net9" attains each of its peak values.
x\text{min} \quad \text{xmin}(\text{o\_waveform \ x\_numberOfValleys}) \\
\Rightarrow \text{o\_waveform/g\_value/l\_value/nil}

\textbf{Description}

Computes the value of the independent variable (X) at which the Y value attains its minimum value.

\textbf{Arguments}

- \text{o\_waveform} \quad \text{Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: srrWave:XXXX.)}

- \text{x\_numberOfValleys} \quad \text{Specifies the n\text{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.}

\textbf{Value Returned}

- \text{o\_waveform} \quad \text{Returns a waveform (or a family of waveforms) if the input argument is a family of waveforms.}

- \text{g\_value} \quad \text{Returns the X value corresponding to the valley specified with x\_numberOfValleys if the input argument is a single waveform.}

- \text{l\_value} \quad \text{Returns a list of X locations when x\_numberOfValleys is 0 and the input argument is a single waveform.}

- \text{nil} \quad \text{Returns nil and an error message otherwise.}

\textbf{Example}

\text{xmin( v( "/net9" ) 1 )}
Gets the time value (X axis) at which the voltage of "/net 9" has its first low point or valley.

```
xmin( v( "/net 9" ) 0 )
```

Gets the list of time values (X axis) at which the voltage of "/net 9" has low points or valleys.
xval

\[ xval( o\_waveform ) -> o\_waveform/nil \]

**Description**

Returns a waveform whose X vector and Y vector are equal to the input waveform’s X vector.

**Arguments**

- \( o\_waveform \)  
  Waveform object representing simulation results that can be displayed as a series of points on a grid. (A waveform object identifier looks like this: \texttt{srrWave:XXXXX}.)

**Value Returned**

- \( o\_waveform \)  
  Returns a waveform if the input argument is a single waveform. Returns a family of waveforms if the input argument is a family of waveforms.

- nil  
  Returns \texttt{nil} and an error message otherwise.

**Example**

\[ xval( v( "/net8" ) ) \]

Returns a waveform in which the X vector for the voltage of "/net8" is also used for the Y vector.
**ymax**

```
ymax( o_waveform )
=> n_max/o_waveformMax/nil
```

**Description**

Computes the maximum value of the waveform’s Y vector.

A waveform consists of an independent-variable X vector and a corresponding Y vector.

**Arguments**

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

**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".
ymin

\[ \text{ymin}( \text{o\_waveform} ) \]
\[ \Rightarrow n\_min/o\_waveformMin/\text{nil} \]

**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**

\text{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: \text{srrWave:XXXXX}.)

**Value Returned**

\text{n\_min}  
Returns a number representing the minimum value of Y if the input argument is a single waveform.

\text{o\_waveformMin}  
Returns a waveform (or family of waveforms) representing the minimum value of Y if the input argument is a family of waveforms.

\text{nil}  
Returns \text{nil} and an error message otherwise.

**Example**

\text{ymin( v( "/net9" ) )}

Gets the minimum voltage (Y value) of "/net9".
Parametric Analysis Commands

These commands set up a parametric analysis. When you run a parametric analysis, you can plot the resulting data as a family of curves.

This chapter contains information on the following commands:

- paramAnalysis on page 448
- paramRun on page 452
paramAnalysis

paramAnalysis(t_desVar [?start n_start] [?stop n_stop] [?center n_center]
    [?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.

\begin{align*}
  n_{\text{start}}, n_{\text{stop}} & \quad \text{highest precedence} \\
  n_{\text{center}}, n_{\text{span}} & \quad \text{precedence}
\end{align*}

Similarly, if you specify more than one step control, the OCEAN environment uses the
following precedence.

\begin{align*}
  f_{\text{step}} & \quad \text{highest precedence} \\
  n_{\text{lin}} & \quad \text{precedence} \\
  n_{\text{dec}} & \\
  n_{\text{log}} & \\
  n_{\text{oct}} & \\
  n_{\text{times}} & \quad \text{lowest precedence}
\end{align*}

To run the analysis, use the \texttt{paramRun} command described in “paramRun” on page 452.

Arguments

\begin{align*}
  t_{\text{desVar}} & \quad \text{Name of the design variable to be swept.} \\
  n_{\text{start}} & \quad \text{Beginning value for the design variable.}
\end{align*}
**n_stop**
Final value for the design variable.

**n_center**
Center point for a range of values that you want to sweep.

**n_span**
Range of values that you want to sweep around the center point. For example, if *n_center* is 100 and *n_span* is 20 then the sweep range extends from 90 to 110.

**f_step**
Increment by which the value of the design variable changes. For example, if *n_start* is 1.0, *n_stop* is 2.1, and *f_step* is 0.2, the parametric analyzer simulates at values 1.0, 1.2, 1.4, 1.6, 1.8, and 2.0.

**n_lin**
The number of steps in the analysis. The parametric analyzer automatically assigns equal intervals between the steps. With this option, there is always a simulation at both *n_start* and *n_stop*. The value for the *n_lin* argument must be an integer greater than 0.

For example, if *n_start* is 0.5, *n_stop* is 2.0, and *n_lin* is 4, the parametric analyzer simulates at values 0.5, 1.0, 1.5, and 2.0.

**n_log**
The number of steps between the starting and stopping points at equal-ratio intervals using the following formula:

\[
\text{log multiplier} = (n\text{-stop}/n\text{-start})^{(n\text{-log}-1)}
\]

The number of steps can be any positive number, such as 0.5, 2, or 6.25.

Default value: 5

For example, if *n_start* is 3, *n_stop* is 15, and *n_log* is 5, the parametric analyzer simulates at values 3, 4.48605, 6.7082, 10.0311, and 15.

The ratios of consecutive values are equal, as shown below.

\[
\]

**n_dec**
The number of steps between the starting and stopping points calculated using the following formula:
\[ \text{decade multiplier} = 10^{1/n_{\text{dec}}} \]

The number of steps can be any positive number, such as 0.5, 2, or 6.25.
Default value: 5

For example, if \( n_{\text{start}} \) is 1, \( n_{\text{stop}} \) is 10, and \( n_{\text{dec}} \) 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 \( 10^1 \).

\( n_{\text{oct}} \)

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 \( n_{\text{start}} \) is 2, \( n_{\text{stop}} \) is 4, and \( n_{\text{oct}} \) 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 \( 2^2 \).

\[ \text{octave multiplier} = 2^{1/(n_{\text{oct}})} \]

\( n_{\text{times}} \)

A multiplier. The parametric analyzer simulates at the points between \( n_{\text{start}} \) and \( n_{\text{stop}} \) that are consecutive multiples of \( n_{\text{times}} \).

For example, if \( n_{\text{start}} \) is 1, \( n_{\text{stop}} \) is 1000, and \( n_{\text{times}} \) is 2, the parametric analyzer simulates at values 1, 2, 4, 8, 16, 32, 64, 128, 256, and 512.

\( n_{\text{spanPercent}} \)

Range specified as a percentage of the center value. For example, if \( n_{\text{center}} \) is 100 and \( n_{\text{spanPercent}} \) is 40, the sweep range extends from 80 to 120.
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.

Example
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("rl" ?start 200 ?stop 600 ?step 200
   paramAnalysis("rs" ?start 300 ?stop 700 ?step 200
  )
)

Sets up a nested parametric analysis for the rl design variable.

paramAnalysis("temp" ?start -50 ?stop 100 ?step 50)

Sets up a parametric analysis for temperature.
paramRun

```
paramRun( [s_paramAnalysis] )
=> t / nil
paramRun( [?jobName t_jobName] [?drmsCmd t_drmsCmd] )
=> s_jobName/nil
paramRun( [?jobName t_jobName] t_jobName )
=> s_jobName/nil
```

**Description**

 Runs the specified parametric analysis.

If you do not specify a parametric analysis, all specified analyses are run. Distributed processing must be enabled using the `hostmode` command before parametric analyses can be run in distributed mode.

When the `paramRun` command finishes, the PSF directory contains a file named `runObjFile` that points to a family of data. To plot the family, use a normal `plot` command. For example, you might use `plot(v("/out")).`

For information about specifying a parametric analysis, see the `paramAnalysis` command described in “paramAnalysis” on page 448.

**Arguments**

- **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.
### OCEAN Reference

#### Parametric Analysis Commands

<table>
<thead>
<tr>
<th>Variable</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td><code>t_termTime</code></td>
<td>Termination time for job. If the job is not completed by <code>t_termTime</code>, the job is terminated.</td>
</tr>
<tr>
<td><code>t_dependentOn</code></td>
<td>List of jobs on which the specified analysis is dependent. The analysis is not started until after dependent jobs are complete.</td>
</tr>
<tr>
<td><code>t_mailingList</code></td>
<td>List of users to be notified by e-mail when the analysis is complete.</td>
</tr>
<tr>
<td><code>s_block</code></td>
<td>When <code>s_block</code> is not nil, the OCEAN script halts until the job is complete.</td>
</tr>
<tr>
<td></td>
<td>Default value: nil</td>
</tr>
<tr>
<td><code>s_notifyFlag</code></td>
<td>When <code>notifyFlag</code> is not nil, a job completion message is echoed to the OCEAN interactive window.</td>
</tr>
<tr>
<td></td>
<td>Default value: <code>t</code></td>
</tr>
<tr>
<td><code>s_paramAnalysis</code></td>
<td>Parametric analysis.</td>
</tr>
<tr>
<td><code>t_drmsCmd</code></td>
<td>A DRMS (Distributed Resource Management System) command, such as a bsub command for LSF or a qsub command for SGE (Sun Grid Engine) used to submit a job. When this argument is used, all other arguments, except <code>?jobName</code> will be ignored. Moreover, it will not be possible to call the OCEAN function <code>wait</code> on the jobs submitted using this argument.</td>
</tr>
<tr>
<td></td>
<td>To know more about the command option, refer to the section Submitting a Job in the chapter <em>Using the Distributed Processing Option in the Analog Design Environment</em> of the <em>Virtuoso Analog Distributed Processing Option User Guide</em>.</td>
</tr>
<tr>
<td><code>s_lsfResourceStr</code></td>
<td>Specifies an LSF Resource Requirement string to submit a job. It is effective only in the LSF mode.</td>
</tr>
</tbody>
</table>

#### Value Returned

<table>
<thead>
<tr>
<th>Value</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td><code>t</code></td>
<td>Returned if successful.</td>
</tr>
<tr>
<td><code>nil</code></td>
<td>Returns <code>nil</code> and prints an error message if unsuccessful.</td>
</tr>
</tbody>
</table>
Example

paramRun() => t

Runs all specified parametric analyses.

rsAnalysis = paramAnalysis("CAP" ?values '(10 20))
paramRun('rsAnalysis)

OR

rsAnalysis = paramAnalysis("CAP" ?values '(10 20) paramAnalysis("RES" ?values '(10 20 )))
paramRun('rsAnalysis)

Runs the rs parametric analysis.

paramRun(?queue "background" ?lsfResourceStr "mem>500")

Runs the analysis in the queue named background on a machine, if it has at least 500 MB of RAM memory.
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 456
- digitalHostMode on page 457
- digitalHostName on page 458
- hostMode on page 459
- hostName on page 460
- killJob on page 461
- monitor on page 462
- remoteDir on page 463
- resumeJob on page 464
- suspendJob on page 465
- wait on page 466

This chapter also provides sample OCEAN scripts that optimally use these commands. See the section Sample Scripts on page 467.

For detailed information on distributed processing, refer to Virtuoso Analog Distributed Processing Option User Guide.
**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**

```sh
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**

- `'local'`: Sets the simulation to run locally on the user’s machine.
- `'remote'`: Sets the simulation to run on a remote host. If you use this argument, you must specify the host name by using the `digitalHostName` command.

**Value Returned**

- `t`: Returns `t` if successful.
- `nil`: Returns `nil` and prints an error message if unsuccessful.

**Example**

digitalHostMode( 'local' )

Sets the digital simulator to run locally on the user’s host.
**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

'local
Sets the simulation to run locally on the user's machine.

'remote
Sets the simulation to run on a remote host queue. For this release, the remote host is specified in the .cdsenv file.

'distributed
Sets the simulation to run using the distributed processing software.

Value Returned

t
Returns t if successful.

nil
Returns nil and prints an error message if unsuccessful.

Example

hostMode( 'distributed )
=> t

Enables distributed processing on the current host.
**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**

*t_name*  
Name used to identify the remote host.

**Value Returned**

*t*  
Returns *t* if successful.

*nil*  
Returns *nil* and prints an error message if unsuccessful.

**Example**

hostName( "remotehost" )

*Specifies that the host called remotehost is to be used for remote simulation.*
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.
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

\textit{t}\_path \hspace{1cm} Specifies the path to the project directory on the remote host to be used for remote simulation.

Value Returned

t \hspace{1cm} Returns t if successful.

\textbf{nil} \hspace{1cm} Returns nil and prints an error message if unsuccessful.

Example
remoteDir( "/simulation" )

Specifies that the project directory is "/simulation."
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.

Additional jobs that you want to resume

t_jobName2...t_jobNameN

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.
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.

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.
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

  t_jobName   Name used to identify the job. The job name is user defined or
               system generated, depending on how the user submitted the job.

  t_jobName2...t_jobnameN  Additional jobs that you want to postpone.

Value Returned

  t         Returns t if successful.

  nil       Returns nil and prints an error message if unsuccessful.

Example

wait( 'myckt1 )
 => t

Postpones execution of all subsequent OCEAN commands until the job myckt1 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 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.

```plaintext
; set up the environment for the jobs
simulator( 'spectre )
hostMode( 'distributed )
design( "//home/simulation/test2/spectre/schematic/netlist/netlist" )
resultsDir( "//home/simulation/test2/spectre/schematic" )
analysis('tran ?stop "5u" )
temp( 27 )

jobList = nil

; starting first job
jobList = append1( jobList run( ?queue "test" ?host "menaka" ) )

analysis('tran ?stop "50u"

; starting second job
jobList = append1( jobList run(?jobName "job_2" ?queue "test" ?host "menaka")

analysis('tran ?stop "10u"

; starting third job, which is dependent on job_2
jobList= append1(jobList run(?jobName "disable" ?queue "test" ?dependentOn symbolToString(car(last(jobList))))

; wait for all the jobs to complete
```
wait((append1 last(jobList) nil))

; open and plot the results of the jobs
openResults( car(last(jobList)))
selectResult( 'tran )
newWindow()
plot(getData("/net61") )

openResults( nth(1 jobList))
selcetResult('tran)
newWindow()
plot(getData("/net61") )

**To set up and run a simple analysis in blocking mode and select results**

; set up the environment for Simple Analysis
simulator( 'spectre )
hostMode( 'distributed )
design( 
  "/home/amit/Artist446/simulation/ampTest/spectre/schematic/netlist/netlist" )
resultsDir( "/home/Artist446/simulation/ampTest/spectre/schematic"")
modelFile( 
  "/home/Artist446/Models/myModels.scs"")
analysis( 'tran ?stop "3u"")
desVar( "CAP" 0.8p )
temp( 27 )

; submit the job in blocking mode, to the queue test and machine menaka
run( ?queue "test" ?host "menaka" ?block t)

; select and plot the results
selectResult( 'tran )
plot(getData( "/out") )

**To set up and run a parametric analysis in blocking mode and select results**

; set up the environment for parametric analysis.
simulator( 'spectre )
hostMode( 'distributed )
design( }
"/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 submit multiple jobs without using wait or selecting results

; set up the environment for the jobs
simulator( 'spectre )
hostMode( 'distributed )
design( ":/home/Artist446/simulation/ampTest/spectre/schematic/netlist/netlist")
resultsDir( ":/home/Artist446/simulation/ampTest/spectre/schematic"
)
modelFile(  
   ("/home/Artist446/Models/myModels.scs"
)
)

; setup and submit first job
analysis(’tran ?stop "3u"  )
desVar( "CAP" 0.8p    )
temp( 27 )
run(?queue "SUN5_5032" ?host "menaka")

; setup and submit second job
analysis(‘ac ?start "1M" ?stop "2M"  )
analysis(’tran ?stop "3u"  )
desVar( "CAP" 0.8p    )
temp( 27 )
run(?queue "SUN5_5032" ?host "menaka")

To submit multiple jobs using wait and selection of results

; set up the environment for the jobs
simulator( 'spectre )
hostMode( 'distributed )
design(
"/home/Artist446/simulation/ampTest/spectre/schematic/netlist/netlist")
resultsDir( "/home/Artist446/simulation/ampTest/spectre/schematic" )
modelFile(
   "'/home/Artist446/Models/myModels.scs""
)

; initialize jobList to nil
jobList = nil

; setup and submit first job
analysis( 'tran ?stop "3u" )
desVar( "CAP" 0.8p )
temp( 27 )
jobList = append1( jobList run(?queue "SUN5_5032" ?host "menaka") )

; setup and submit second job
analysis( 'ac ?start "1M" ?stop "2M" )
analysis( 'tran ?stop "3u" )
desVar( "CAP" 0.8p )
temp( 27 )
jobList = append1( jobList run(?queue "SUN5_5032" ?host "menaka") )

; wait for both the jobs to finish
wait( (append1 jobList nil) )

; open and plot the result of first job
openResults( (car jobList))
selectResult( 'tran )
plot(getData("/out") )

; open and plot the result of second job
openResults( (cadr jobList))
selectResult( 'tran )
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 483
- cond on page 485
- for on page 478
- foreach on page 480
- if on page 474
- unless on page 476
- when on page 477
- while on page 482
if

\[ \text{if}( \ g\_condition \ g\_thenExpression \ [g\_elseExpression] \ ) \]
\[ \Rightarrow \ g\_result \]

**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 \text{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.

**Example**

\[ x = 2 \]
\[ \text{if}( \ x > 5 \) \]
\[ \Rightarrow 0 \]

Returns 0 because \( x \) is less than 5.

\[ a = "npn" \]
\[ \text{if}( \ a == "npn" \) \]
\[ \text{print}( \ a \) \]
\[ \Rightarrow "npn" \]

Prints the string \( npn \) and returns the result of \text{print}.

\[ x = 5 \]
\[ \text{if}( \ x "non-nil" "nil" \) \]
\[ \Rightarrow "non-nil" \]
Returns "non-nil" because \( x \) was not nil. If \( x \) was nil, "nil" would be returned.

\[
x = 7 \\
if( x > 5 ) \\
=> 1
\]

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.

Example

  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

- \textit{g\_condition} Any SKILL expression.
- \textit{g\_expr1...} Any SKILL expression.

Value Returned

- \textit{g\_result} Returns the value of the last expression of the sequence \textit{g\_expr1...} if \textit{g\_condition} evaluates to non-nil.
- \textit{nil} returns \textit{nil} if the \textit{g\_condition} expression evaluates to \textit{nil}.

Example

- \texttt{x = -123}
- \texttt{when( x < 0 println( "x is negative" ) -x ) => 123}

Prints "x is negative" as a side effect.

- \texttt{when( x >= 0 println( "x is positive" ) x) => nil}

Returns \textit{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

s_loopVar Name of the local loop variable that must not be changed inside the loop.

x_initialValue Integer expression setting the initial value for the local loop variable.

x_finalValue Integer expression giving final limit value for the loop.

g_expr1 Expression to evaluate inside loop.

g_expr2 ... Additional expressions to evaluate inside loop.

Value Returned

t This construct always returns t.
Example

```c
sum = 0
for( i 1 10
    sum = sum + i
    printf( "%d" sum ))
=> t
```

Prints 10 numbers and returns t.

```c
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

$s_{formalVar}$ Name of the variable.

$g_{exprList}$ Expression whose value is a list of elements to assign to the
formal variable $s_{formalVar}$.

$g_{expr1}$ $g_{expr2}$ ... Expressions to execute.

$g_{exprTable}$ Association table whose elements are to be processed.

Value Returned

$l_{valueList}$ Returns the value of the second argument, $g_{exprList}$.

$o_{valueTable}$ Returns the value of $g_{exprTable}$.

Example

```plaintext
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.

```plaintext
foreach( key myTable printf( "%L : %L" key myTable[key] ) )
```

Accesses an association table and prints each key and its associated data.

```plaintext
( 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 \textit{g\_condition} and the sequence of expressions \textit{g\_expr1} ... if the condition is true.

This process is repeated until \textit{g\_condition} evaluates to false (\textit{nil}). Note that because this form always returns \textit{t}, it is principally used for its side effects.

\textbf{Note:} Everything that can be done with a \textit{for} loop can also be done with a \textit{while} loop.

Arguments

\begin{itemize}
\item \textit{g\_condition} \hspace{1em} Any SKILL expression.
\item \textit{g\_expr1} \hspace{1em} Any SKILL expression.
\end{itemize}

Value Returned

\textit{t} \hspace{1em} Always returns \textit{t}.

Example

\begin{verbatim}
i = 0
while( (i <= 10) printf(\%d\ i++) )
=> t
\end{verbatim}

Prints the digits 0 through 10.
**case**

```latex
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 \textit{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 \textit{g\_selectionExpr} expression is evaluated and the resulting selector value is matched sequentially against comparators defined in \textit{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 \textit{case} is then the value of the last expression evaluated (that is, the last expression in the clause selected). If there is no match, \textit{case} returns \textit{nil}.

The symbol \texttt{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**

\begin{itemize}
    
    \begin{itemize}
        \item \textit{g\_selectionExpr} \hspace{1cm} 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.
        
        \item \textit{l\_clause1} \hspace{1cm} An expression whose first element is an atom or list of atoms to be compared against the value of \textit{g\_selectionExpr}. The remainder of the \textit{l\_clause} is evaluated if a match is found.
        
        \item \textit{l\_clause2...} \hspace{1cm} Zero or more clauses of the same form as \textit{l\_clause1}.
    \end{itemize}
\end{itemize}

**Value Returned**

\begin{itemize}
    \item \textit{g\_result} \hspace{1cm} Returns the value of the last expression evaluated in the matched clause.
\end{itemize}
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".
Tests each of the arguments according to the conditions specified with cond.
This chapter contains information on the following commands:

close on page 488
fscanf on page 489
gets on page 491
infile on page 492
load on page 493
newline on page 495
outfile on page 496
pfile on page 498
printf on page 499
println on page 500
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**

\[
\text{fscanf}(\ p_{\text{inputPort}} \ t_{\text{formatString}} \ [s_{\text{var1}} \ ...] ) \\
=> x_{\text{items}}/\text{nil}
\]

**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**

<table>
<thead>
<tr>
<th>Format Specification</th>
<th>Types of Argument</th>
<th>Scans for</th>
</tr>
</thead>
<tbody>
<tr>
<td>%d</td>
<td>fixnum</td>
<td>An integer</td>
</tr>
<tr>
<td>%f</td>
<td>flonum</td>
<td>A floating-point number</td>
</tr>
<tr>
<td>%s</td>
<td>string</td>
<td>A string (delimited by spaces) in the input</td>
</tr>
</tbody>
</table>

**Arguments**

\[
p_{\text{inputPort}} \text{ Input port to read from.}
\]

\[
t_{\text{formatString}} \text{ Format string to match against in the reading.}
\]

\[
s_{\text{var1}}... \text{ Name of the variable in which to store results.}
\]
Value Returned

\textit{x\_items} \quad \text{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.}

\texttt{nil} \quad \text{Returns \texttt{nil} if an end of file is encountered}

Example

\begin{verbatim}
\texttt{fscanf( p \"%d %f\" i d )}
\end{verbatim}

Scans for an integer and a floating-point number from the input port \texttt{p} and stores the values read in the variables \texttt{i} and \texttt{d}, respectively.

Assume a file \texttt{testcase} with one line:

\begin{verbatim}
hello 2 3 world
\end{verbatim}

\begin{verbatim}
\texttt{x = infile("testcase")}
\texttt{\rightarrow \texttt{port:}\texttt{"testcase"}}
\texttt{fscanf( x \"%s %d %d %s\" a b c d )}
\texttt{\rightarrow 4}
\texttt{(list a b c d) \rightarrow ("hello" 2 3 "world")}
\end{verbatim}
gets
gets( s_variableName [p_inputPort] ) => t_string/nil

Description
Reads a line from the input port and stores the line as a string in the variable. This is a macro.
The string is also returned as the value of gets. The terminating newline character of the line becomes the last character in the string.

Arguments
s_variableName Variable in which to store the input string.
p_inputPort Name of input port.
  Default value: piport

Value Returned
t_string Returns the input string when successful.
nil Returns nil when the end of file is reached.
  (s_variableName maintains its last value.)

Example
Assume the test1.data file has the following first two lines:
#This is the data for test1
0001 1100 1011 0111
p = infile("test1.data") => port:"test1.data"
gets(s p) => "#This is the data for test1"
gets(s p) => "0001 1100 1011 0111"
s => "0001 1100 1011 0111"

Gets a line from the test1.data file and stores it in the variable s. The s variable contains the last string stored in it by the gets function.
**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.

**Example**

in = infile( "~/test/input.il" ) => port:"~/test/input.il"
close( in )
=> t

Closes the /test/input.il port.

Opens the input port /test/input.il.

infile("myFile") => nil

Returns nil if myFile does not exist according to the current setting of the SKILL path or exists but is not readable.
load

load( _fileName [ _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

_fileName File to be loaded. Uses the file name extension to determine the language mode to use.

Valid values:

- .il Means the file contains SKILL code.
- .ils Means the file contains SKILL++ code.
- .ocn Means the file contains OCEAN commands (with SKILL or SKILL++ commands)

_password Password, if _fileName is an encrypted file.
Value Returned

t

Returns t if the file is successfully loaded.

Example

load( "test.ocn" )

Loads the test.ocn file.

procedure( trLoadSystem()  
   load( "test.il" );;; SKILL code  
   load( "test.ils" );;; SKILL++ code  
); procedure

You might have an application partitioned into two files. Assume that test.il contains SKILL code and test.ils contains SKILL/SKILL++ code. This example loads both files.
newline

newline( [p_outputPort] )
=> nil

Description
Prints a newline (backslash \n) character and then flushes the output port.

Arguments

p_outputPort Output port.
Defaults value: poport

Value Returned
nil Prints a newline and then returns nil.

Example
print( "Hello" ) newline() print( "World!" )
"Hello"
"World!"
=> nil

Prints a newline character after Hello.
outfile

outfile( S_fileName [t_mode] )
=> p_outport/nil

Description

Opens an output port ready to write to a file.

Various print commands can write to this file. Commands write first to a character buffer, which writes to the file when the character buffer is full. If the character buffer is not full, the contents are not written to the file until the output port is closed or the drain command is entered. Use the close or drain command to write the contents of the character buffer to the file. The file can be specified with either an absolute path or a relative path. If a relative path is given and the current SKILL path setting is not nil, all directory paths from SKILL path are checked in order, for that file. If found, the system overwrites the first updatable file in the list. If no updatable file is found, it places a new file of that name in the first writable directory.

Arguments

S_fileName       Name of the file to open or create.

t_mode       Mode in which to open the file. If a, the file is opened in append mode; If w, a new file is created for writing (any existing file is overwritten).
            Default value: w

Value Returned

p_outport       An output port ready to write to the specified file.

returns nil if the named file cannot be opened for writing. An error is signaled if an illegal mode string is supplied.

Example

p = outfile( "~/tmp/out.il" "w" )
=> port:"~/tmp/out.il"

Opens the /tmp/out.il port.

outfile( "~/bin/ls" )
=> nil
Returns \texttt{nil}, indicating that the specified port could not be opened.
pfile

pfile( [S_fileName | p_port] )
   => p_port/nil

Description

Opens an output port ready to write to a file or returns the name of an existing port indicating that it is available.

This command is similar to the outfile command when a valid S_fileName is specified. When p_port is specified, it returns the file port that is currently being used by p_port. When no argument is specified, it opens the stdout port.

Arguments

S_fileName            Name of the file to open or create.
P_port                Retrieves the name of the file port that is being used.

Value Returned

P_port                The ID of the port that was opened, or stdout.
nil                    Returns nil if the named file cannot be opened for writing.

Example

p = pfile( "/tmp/out.il" "w" )
=> port:"/tmp/out.il"

Opens the /tmp/out.il port.
pfile( "/bin/ls" )
=> nil

Returns nil, indicating that the specified port could not be opened.
p = pfile()
=> port:"*stdout*"

Returns stdout as the file port indicating that stdout has been opened.
pfile( p )
=> port:"/tmp/out.il"

Returns the file port.
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 Cadence SKILL Language User Guide.

printf is identical to fprintf except that it does not take the p_port argument and the output is written to poport.

Arguments

  t_formatString
      Characters to be printed verbatim, intermixed with format specifications prefixed by the “%” sign.

  g_arg1...
      Arguments following the format string are printed according to their corresponding format specifications.

Value Returned

  t
      Prints the formatted output and returns t.

Example

  x = 197.9687 => 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 \textit{g\_value}. The \textit{println} function flushes the output port after printing each newline character.

Arguments
\begin{itemize}
  \item \textit{g\_value} \hspace{1cm} Any SKILL value.
  \item \textit{p\_outputPort} \hspace{1cm} Port to be used for output.
      \hspace{1cm} Default value: \textit{poport}
\end{itemize}

Value Returned
\begin{itemize}
  \item \textit{nil} \hspace{1cm} Prints the given object and returns \textit{nil}.
\end{itemize}

Example
\begin{verbatim}
for( i 1 3 println( "hello" ))
"hello"
"hello"
"hello"
=> t
\end{verbatim}

Prints hello three times. \texttt{for} always returns \texttt{t}.
OCEAN Commands in XL Mode

The following OCEAN XL commands provide for multi-test ADE XL support in OCEAN.

- ocnSetXLMode on page 504
- ocnxlBeginTest on page 505
- ocnxlEndTest on page 506
- ocnxlEndXLMode on page 507
- ocnxlSelectTest on page 508
- ocnxlSweepVar on page 509
- ocnxlSweepParam on page 510
- ocnxlCorner on page 511
- ocnxlDisableTest on page 512
- ocnxlDisableSweepVar on page 513
- ocnxlDisableSweepParam on page 514
- ocnxlDisableCornerForTest on page 515
- ocnxlGlobalOptimizationOptions on page 516
- ocnxlJobSetup on page 518
- ocnxlLocalOptimizationOptions on page 519
- ocnxlModelGroup on page 520
- ocnxlOutputOceanScript on page 521
- ocnxlOutputMatlabScript on page 522
ocnxIMonteCarloOptions on page 523
ocnxIPutToleranceSpec on page 525
ocnxIPutMinSpec on page 526
ocnxIPutMaxSpec on page 527
ocnxIPutGreaterthanSpec on page 528
ocnxIPutLessthanSpec on page 529
ocnxIPutRangeSpec on page 530
ocnxIPutTargetSpec on page 531
ocnxIResultsLocation on page 532
ocnxIRun on page 533
ocnxIRunSetupSummary on page 534
ocnxISamplingOptions on page 535
ocnxISetupLocation on page 536
ocnxIOutputExpr on page 537
ocnxIOutputSignal on page 538
ocnxIOutputTerminal on page 539
ocnxIOutputSummary on page 540
ocnxITargetCellView on page 542
ocnxIYieldImprovementOptions on page 543
ocnxIEnableCornerForTest on page 546
ocnxIEnableSweepParam on page 547
ocnxIEnableSweepVar on page 548
ocnxIEnableTest on page 549
ocnxIGetBestPointParams on page 550
ocnxlGetCorners on page 551
ocnxlGetCurrentHistory on page 552
ocnxlGetCurrentHistoryId on page 553
ocnxlGetSession on page 554
ocnxlGetSpecs on page 555
ocnxlGetTests on page 556
ocnxlRemoveSpec on page 557
ocnxlRenameCurrentHistory on page 558
ocnxlRun on page 559
ocnxlHistoryPrefix on page 561
ocnxlLoadSetupState on page 562
ocnxlStartingPoint on page 565
ocnxlOutputAreaGoal on page 566
ocnxlConjugateGradientOptions on page 567
ocnxlMTSEnable on page 568
ocnxlMTSBloc k on page 569
ocnxlProjectDir on page 571
ocnxlSimResultsLocation on page 572
ocnxlDisableCorner on page 572
ocnxlEnableCorner on page 574
ocnxlSaveSetupAs on page 575
ocnxlSetAllParametersDisabled on page 576
ocnxlSetAllVarsDisabled on page 577
ocnSetXLMode

ocnSetXLMode()
=> t / nil

Description

Sets OCEAN mode to XL. In this mode, multi-tests can be created and run. Also Corners, MonteCarlo commands can be given. Once mode is set to XL, it cannot be reverted.

Arguments

None

Value Returned

<table>
<thead>
<tr>
<th>Value</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>t</td>
<td>Returns t if the mode is set to XL.</td>
</tr>
<tr>
<td>nil</td>
<td>Returns nil otherwise.</td>
</tr>
</tbody>
</table>

Example

ocnSetXLMode()
ocnxlBeginTest

ocnxlBeginTest(t_testName)
   => t / nil

Description

This command indicates the beginning of the test specified by testName. Subsequent commands populate this test. The test specification ends when ocnxlEndTest() command is given. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

t_testName The name of the test.

Value Returned

t Returns t if its able to begin the test.
nil Returns nil otherwise.

Example

ocnxlBeginTest("test_one")
ocnxlEndTest

ocnxlEndTest()
=> t / nil

Description

This command indicates the end of the current test’s specification. See help on ocnxlBeginTest(). This command works only in XL mode. See help on ocnSetXLMode().

Arguments

None.

Value Returned

<table>
<thead>
<tr>
<th>t</th>
<th>Returns t if the test setup completes.</th>
</tr>
</thead>
<tbody>
<tr>
<td>nil</td>
<td>Returns nil otherwise.</td>
</tr>
</tbody>
</table>

Example

ocnxlBeginTest("test_one")
design("solutions" "ampTest" "schematic")
simulator('spectre)
ocnxlEndTest()
ocnxIEndXLMode

ocnxIEndXLMode()
=> t

Description

This command ends the XL mode. Also see help on ocnxlSetXLMode().

Arguments

None.

Value Returned

t

Returns t if it exits the XL mode.

Example

ocnxIEndXLMode()
ocnxlSelectTest

ocnxlSelectTest (t_testName)
    => t / nil

Description

Lets you select a test. List of test names can be obtained by ocnxlGetTests() command. See help on ocnxlGetTests(). This command works only in XL mode. See help on ocnSetXLMode().

Arguments

$t_testName$ The name of the test.

Value Returned

  t Returns t if the test is selected.

  nil Returns nil otherwise.

Example

ocnxlSelectTest("test_two") => t
Sets "test_two" as the currently selected test.
ocnxlSweepVar

ocnxlSweepVar(t_varName t_varValue)
    => t / nil

Description

Lets you define a sweep variable along with its value. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

  t_varName             Name of the variable.
  t_varValue            Value of the variable and the specification for the sweep.

Value Returned

  t                   Returns t if the sweep is set.
  nil                 Returns nil otherwise.

Example

  ocnxlSweepVar("CAP" "5p") => t
ocnxlSweepParam

ocnxlSweepParam(t_paramName t_paramValue)
   => t / nil

Description

Lets you define a sweep parameter along with its value. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

\textit{t\_paramName} \quad \text{Name of the parameter.}
\textit{t\_paramValue} \quad \text{Value of the parameter.}

Value Returned

\textit{t} \quad \text{Returns } t \text{ if sweep for the parameter is set.}
\textit{nil} \quad \text{Returns } \text{nil otherwise.}

Example

ocnxlSweepParam("solutions/ampTest/schematic/R1/r" "10K") => t
ocnxlCorner

ocnxlCorner(t_cornerName l_cornerDetails)
   => t / nil

Description

Lets you define a corner. cornerDetails is a list of elements where each element is (t_type t_varName t_value). Available types are variable, parameter, and model. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

t_cornerName  Name of the corner.

l_cornerDetails  Details of the corner. Details is a list of items where each item has a tag, name, and a value. The tag can be of 3 types — variable, parameter, and model.

Value Returned

t  Returns t if the corner is defined.

nil  Returns nil otherwise.

Example

ocnxlCorner("C0" '( ("variable" "CAP" "2p") ("variable" "T" "78"))) => t
**ocnxlDisableTest**

`ocnxlDisableTest(t_testName)`  
=> t / nil

**Description**

Lets you disable a test. A disabled test will not be run when `ocnxlRun()` command is fired. See help on `ocnxlRun()`. This command works only in XL mode. See help on `ocnSetXLMode()`.

**Arguments**

`t_testName` Name of the test.

**Value Returned**

<table>
<thead>
<tr>
<th>t</th>
<th>Returns t if the test is disabled.</th>
</tr>
</thead>
<tbody>
<tr>
<td>nil</td>
<td>Returns nil otherwise.</td>
</tr>
</tbody>
</table>

**Example**

`ocnxlDisableTest("test_two")` => t
ocnxlDisableSweepVar

ocnxlDisableSweepVar(t_varName)
 => t / nil

Description

Lets you disable a sweep variable. A disabled sweep is not run when `ocnxlRun()` command is fired. See help on `ocnxlRun()`. This command works only in XL mode. See help on `ocnSetXLMode()`.

Arguments

`t_varName`   Name of the variable.

Value Returned

`t`          Returns `t` if the sweep variable is disabled.

`nil`     Returns `nil` otherwise.

Example

`ocnxlDisableSweepVar("CAP") => t`
ocnxlDisableSweepParam

ocnxlDisableSweepParam(t_paramName)
    => t / nil

Description

Lets you disable a sweep parameter. A disabled parameter is not run when ocnxlRun() command is fired. See help on ocnxlRun(). This command works only in XL mode. See help on ocnSetXLMode().

Arguments

t_paramName                  Name of the parameter.

Value Returned

t                  Returns t if the sweep parameter is disabled.
nil                Returns nil otherwise.

Example

ocnxlDisableSweepParam("solutions/ampTest/schematic/R1/r") => t
ocnxlDisableCornerForTest

ocnxlDisableCornerForTest(t_cornerName t_testName)
   => t / nil

Description

Lets you disable a corner for a test. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

  t_cornerName       Name of the corner.
  t_testName         Name of the test.

Value Returned

  t               Returns t if the corner of the test is disabled.
  nil             Returns nil otherwise.

Example

ocnxlDisableCornerForTest("C0" "test_one")
ocnxIGlobalOptimizationOptions

ocnxIGlobalOptimizationOptions( [ ?effort t_effort ]
   [ ?runFullEvaluation t_runFullEvaluation ]
   [ ?refPoint t_refPoint ]
   [ ?meetAllGoals t_meetAllGoals ]
   [ ?timeLimit t_timeLimit ]
   [ ?numPoints t_numPoints ]
   [ ?noImprvPoints t_noImprvPoints ]
   [ ?pointsAfterAllSpecsSatisfied t_pointsAfterAllSpecsSatisfied ]
)
=> t / nil

Description

Lets you specify options for global optimization run. See help on ocnxlRun() for run modes. This command works only in XL mode. See help on ocnSetXLMode()

Arguments

\textbf{t\_effort} \hspace{1cm} \text{The value for effort. Default for \texttt{effort} is \texttt{coarse}. Possible values are \texttt{fine} and \texttt{coarse}.}

\textbf{t\_runFullEvaluation} \hspace{1cm} \text{Sets the program to run full optimization. Default for \texttt{runFullEvaluation} is \texttt{0}. Possible values are \texttt{0} and \texttt{1}.}

\textbf{t\_refPoint} \hspace{1cm} \text{Sets the current point as a starting place for sizing. Default for \texttt{refPoint} is \texttt{0}. Possible values are \texttt{0} and \texttt{1}.}

\textbf{t\_meetAllGoals} \hspace{1cm} \text{Sets the program to run optimization only until all specifications are met. Default for \texttt{meetAllGoals} is \texttt{0}. Possible values are \texttt{0} and \texttt{1}.}

\textbf{t\_timeLimit} \hspace{1cm} \text{Sets the time limit (in minutes) for the optimization run. Default for \texttt{timeLimit} is "".}

\textbf{t\_numPoints} \hspace{1cm} \text{Sets the maximum number of points for the optimization run. Default for \texttt{numPoints} is "".}

\textbf{t\_noImprvPoints} \hspace{1cm} \text{Default for \texttt{noImprvPoints} is "".}

\textbf{t\_pointsAfterAllSpecsSatisfied} \hspace{1cm} \text{Sets the number of points to be run after all specifications are satisfied. Default for \texttt{t\_pointsAfterAllSpecsSatisfied} is "".}
Value Returned

\( t \) Returns \( t \) if options are specified for global optimization run.

\( \text{nil} \) Returns \( \text{nil} \) otherwise.

Example

\[ \text{ocnxlGlobalOptimizationOptions(?runFullEvaluation "1")} \]

\# Sets global optimization to be run only until all specifications are met.
ocnxlJobSetup

ocnxlJobSetup (l_setupOptions)
   => t / nil

Description

 Specifies job setup information. This information guides how ocean xl distributed jobs. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

l_setupOptions The list of setup options. This depends on the underlying distribution method used.

Value Returned

 t Returns t if the job setup information is specified.
 nil Returns nil otherwise.

Example

ocnxlJobSetup (’( "configuretimeout" "300" "distributionmethod" "Local" "lingertimeout" "300" "maxjobs" "1" "runtimeout" "-1" "starttimeout" "300") )
ocnx1LocalOptimizationOptions

ocnx1LocalOptimizationOptions( [ ?effort t_effort ]
   [ ?runFullEvaluation t_runFullEvaluation ] [ ?meetAllGoals t_meetAllGoals ]
   [ ?timeLimit t_timeLimit ] [ ?numPoints t_numPoints ] )
=> t / nil

Description

Lets you specify options for local optimization run. See help on ocnxlRun() for run modes.

Arguments

\( t_{\text{effort}} \)

Value for effort. Default for effort is coarse. Possible values are fine and coarse.

\( t_{\text{runFullEvaluation}} \)

Sets the program to run full optimization. Default for runFullEvaluation is 0. Possible values are 0 and 1.

\( t_{\text{meetAllGoals}} \)

Sets the program to run optimization only until all specifications are met. Default for meetAllGoals is 0. Possible values are 0 and 1.

\( t_{\text{timeLimit}} \)

Sets the time limit (in minutes) for the optimization run. Default for timeLimit is "".

\( t_{\text{numPoints}} \)

Sets the maximum number of points for the optimization run. Default for numPoints is "".

Value Returned

\( t \)

Returns \( t \) if options are specified for local optimization run.

\( \text{nil} \)

Returns \text{nil} otherwise.

Example

ocnx1LocalOptimizationOptions(?effort "coarse" )
# Sets coarse as the effort for local optimization run.
ocnxlModelGroup

ocnxlModelGroup( t_modelGroupName l_modelFileSetup) => t/nil

Description

Lets you add and define a new model group. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

<table>
<thead>
<tr>
<th>t_modelGroupName</th>
<th>Name of the new model group.</th>
</tr>
</thead>
<tbody>
<tr>
<td>l_modelFileSetup</td>
<td>List of model file spec.</td>
</tr>
</tbody>
</table>

Model file spec:

\[
( t_modelFilePath
  [?section t_section]
  [?enabled g_enabled]
  [?test t_test]
  [?block t_block]
)
\]

Value Returned

<table>
<thead>
<tr>
<th>t</th>
<th>Returns t if a new model is defined.</th>
</tr>
</thead>
<tbody>
<tr>
<td>nil</td>
<td>Returns nil otherwise.</td>
</tr>
</tbody>
</table>

Example

ocnxlModelGroup( "F2"

'(  
  
( "/myModels/Models/model1.scs" ?enabled nil ?section "")
( "/myModels/Models/model2.scs" ?section "")
( "/myModels/Models/model3.scs" ?enabled nil ?section "")
))
ocnxlOutputOceanScript

ocnxlOutputOceanScript(t_script [ ?name t_outputName ] [ ?plot plot ] [ ?save save ])
=> t/nil

Description

Adds an OCEAN script based output in the current test being specified. A test's specification begins with ocnxlBeginTest(). See help on ocnxlBeginTest().

This command works only in XL mode. See help on ocnSetXLMode().

Arguments

- **t_script** Name and location of the script file.
- **t_outputName** Name of the output file.
- **plot** Specifies if the values are to be plotted.
- **save** Specifies if the output are to be saved.

Value Returned

- **t** Returns t if the output is generated.
- **nil** Returns nil otherwise.

Example

ocnxlOutputOceanScript( "/tmp/my_measure.ocn" ?name "MAX" ?plot t ?save t ) => t
Adds "MAX" in the outputs.
ocnxlOutputMatlabScript

ocnxlOutputMatlabScript( t_script [ ?name t_outputName ] [ ?plot plot ] [ ?save save ] ) => t/nil

Description

Adds a MATLAB script based output in the current test being specified. A test’s specification begins with ocnxlBeginTest(). See help on ocnxlBeginTest().

This command works only in XL mode. See help on ocnSetXLMode().

Arguments

t_script Name and location of the script file.

t_outputName Name of the output file.

plot Specifies if the values are to be plotted.

save Specifies if the output are to be saved.

Value Returned

t Returns t if the output is generated.

nil Returns nil otherwise.

Example

ocnxlOutputMatlabScript( "/tmp/my_measure.m" ?name "MAX" ?plot t ?save t ) => t

Adds "MAX" in the outputs.
ocnxI Monte Carlo Options

ocnxI Monte Carlo Options( [ ?mcmethod t_mcmethod ] [ ?mcNumPoints t_mcNumPoints ] [ ?samplingMode t_samplingMode ] [ ?saveAllPlots t_saveAllPlots ] [ ?saveProcess t_saveProcess ] [ ?saveMismatch t_saveMismatch ] [ ?dutSummary t_dutSummary ] [ ?useReference t_useReference ] [ ?donominal t_donominal ] [ ?monteCarloSeed t_monteCarloSeed ] [ ?mcStartingRunNumber t_mcStartingRunNumber ] [ ?designUnderTest t_designUnderTest ] [ ?dutInstances t_dutInstances ] )
=> t / nil

Description

Lets you specify options for Monte Carlo runs. See help on ocnxI Run() for run modes.

Arguments

\textbf{t\_mcmethod} \quad Sets the statistical variation method for Monte Carlo runs. Default for mcmethod is all. Possible values are global, mismatch and all.

\textbf{t\_mcNumPoints} \quad Sets the number of points you want to simulate for Monte Carlo runs. Default for mcNumPoints is 100.

\textbf{t\_samplingMode} \quad Sets the default statistical sampling method for Monte Carlo runs. Default for samplingMode is random. Possible values are random and lhs.

\textbf{t\_saveAllPlots} \quad Saves raw data (psf files) for every Monte Carlo iteration so that you can plot a family of curves. Default for saveAllPlots is 0. Possible values are 0 and 1.

\textbf{t\_saveProcess} \quad Controls whether 'process' parameters need to be saved to the results database. Default value is “1”. Possible values are “0” and “1”.

\textbf{t\_saveMismatch} \quad Controls whether ‘mismatch’ parameters need to be saved to the results database. Default value is “0”. Possible values are “0” and “1”.

\textbf{t\_dutSummary} \quad Optional argument for a user to specify DUT instances in a Monte Carlo run.
Value Returned

t
Returns t if options for montecarlo run are specified.

nil
Returns nil otherwise.

Example


\textit{t} \textbf{useReference}

Specifies whether to use a schematic point or a reference point that you have created as a starting place for sizing.

Possible values are “0” and “1”. The default value is “0”.

\textit{t} \textbf{donominal}

Specifies whether to run a simulation at the reference point prior to beginning the Monte Carlo process. Possible values are “0” and “1”. If set to “1”, Spectre will run a simulation at the reference point, and, if this fails, then the sampling process is not initiated and the simulation stops.

The default value is “1”.

\textit{t} \textbf{monteCarloSeed}

Specifies a different seed for Monte Carlo runs. Default for \texttt{monteCarloSeed} is “12345”.

\textit{t} \textbf{mcStartingRunNumber}

Specifies a starting run number for Monte Carlo runs. Default for \texttt{mcStartingRunNumber} is “1”.

\textit{t} \textbf{designUnderTest}

Specifies the design under test (DUT) for Monte Carlo runs. Default for \texttt{designUnderTest} is "". Specify the DUT using the format:

\texttt{"libname:cellname:viewname"}

\textit{t} \textbf{dutInstances}

Specifies the the DUT instances for Monte Carlo runs. Default for \texttt{dutInstances} is "". Specify the DUT instances using the characters %/ to separate two DUT instances. For example, specify:

\texttt{"libname:cellname:instname1%/}
\texttt{libname:cellname:instname2"}
ocnx1PutToleranceSpec

ocnx1PutToleranceSpec(t_testName t_resultName t_value
    s_type t_tolerance g_weight)
=> t / nil

Description

Lets you specify a tolerance spec for a result. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

- **t_testName**: The name of the test.
- **t_resultName**: The name of the result.
- **t_value**: The target value.
- **s_type**: The type of tolerance.
- **t_value**: The tolerance value.
- **g_weight**: The weighting factor for the spec.

Value Returned

- **t**: Returns `t` if the specifications are specified.
- **nil**: Returns `nil` otherwise.

Example

```
ocnx1PutToleranceSpec("test_one" "VT('/out')" "5.0" 'percent" "10" 4) => t
# Spec is defined that transient voltage for /out signal must be 5.0 volts with
tolerance 10%. The weighting factor for the spec is 4.
```
ocnxIPutMinSpec

ocnxIPutMinSpec(t_testName t_resultName t_minValue g_weight)
 => t / nil

Description

Lets you specify a minimum spec for a result. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

$t_testName$ Name of the test.
$t_resultName$ Name of the result on which you specify the test.
$t_minValue$ The minimum value.
$g_weight$ The weighting factor for the spec.

Value Returned

$t$ Returns $t$ if the specification is set.
$nil$ Returns $nil$ otherwise.

Example

ocnxIPutMinSpec("test_one" "VT('/out')" "3.5" 4) => t
Spec is defined that minimum transient voltage for /out signal must be 3.5 volts. The weighting factor for the spec is 4.
ocnxlPutMaxSpec

ocnxlPutMaxSpec(t_testName t_resultName t_maxValue g_weight)
=> t / nil

Description

Lets you specify a maximum spec for a result. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

- t_testName: Name of the test.
- t_resultName: Name of the result on which you specify the test.
- t_maxValue: The maximum value.
- g_weight: The weighting factor for the spec.

Value Returned

- t: Returns t if the specification is set.
- nil: Returns nil otherwise.

Example

ocnxlPutMaxSpec("test_one" "VT('/out')" "6.5" 4) => t
Spec is defined that maximum transient voltage for /out signal must be 6.5 volts. The weighting factor for the spec is 4.
ocnxlPutGreaterthanSpec

ocnxlPutGreaterthanSpec(t_testName t_resultName t_Value g_weight)
=> t / nil

Description

Lets you specify that value for a result must be greater than a threshold value. This command
works only in XL mode. See help on ocnSetXLMode().

Arguments

<table>
<thead>
<tr>
<th>Argument</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>t_testName</td>
<td>Name of the test.</td>
</tr>
<tr>
<td>t_resultName</td>
<td>Name of the result on which you specify the test.</td>
</tr>
<tr>
<td>t_Value</td>
<td>The threshold value.</td>
</tr>
<tr>
<td>g_weight</td>
<td>The weighting factor for the spec.</td>
</tr>
</tbody>
</table>

Value Returned

<table>
<thead>
<tr>
<th>Value</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>t</td>
<td>Returns t if the specification is set.</td>
</tr>
<tr>
<td>nil</td>
<td>Returns nil otherwise.</td>
</tr>
</tbody>
</table>

Example

ocnxlPutGreaterthanSpec("test_one" "VT('/out')" "3.5" 4) => t
Spec is defined that transient voltage for /out signal
must always be greater than 3.5 volts. The weighting factor for the spec is 4.
ocnxlPutLess-than-Spec

ocnxlPutLess-than-Spec(t_testName t_resultName t_Value g_weight)
  => t / nil

Description

Lets you specify that value for a result must be less than a threshold value. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

- **t_testName**
  - Name of the test.
- **t_resultName**
  - Name of the result on which you specify the test.
- **t_Value**
  - The threshold value.
- **g_weight**
  - The weighting factor for the spec.

Value Returned

- **t**
  - Returns t if the specification is set.
- **nil**
  - Returns nil otherwise.

Example

```plaintext
ocnxlPutLess-than-Spec("test_one" "VT('/out')" "6.5" 4) => t
Spec is defined that the transient voltage for /out signal must always be less than 6.5 volts. The weighting factor for the spec is 4.
```
ocnxlPutRangeSpec

ocnxlPutRangeSpec(t_testName t_resultName t_maxValue t_minValue g_weight)
  => t / nil

Description

Lets you specify a range spec for a result. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

  t_testName Name of the test.
  t_resultName Name of the result on which you specify the test.
  t_maxValue The maximum value.
  t_minValue The minimum value.
  g_weight The weighting factor for the spec.

Value Returned

  t Returns t if the specification is set.
  nil Returns nil otherwise.

Example

ocnxlPutRangeSpec("test_one" "VT('/out')" "6.5" "3.5" 4) => t
Spec is defined that maximum transient voltage for /out signal
must be 6.5 volts and minimum must be 3.5 volts. The weighting factor for the spec
is 4.
**ocnxlPutTargetSpec**

`ocnxlPutTargetSpec(t_testName t_resultName t_value)`

=> t / nil

**Description**

Lets you specify a target spec for a result. This command works only in XL mode. See help on `ocnSetXLMode()`.

**Arguments**

- `t_testName` Name of the test.
- `t_resultName` Name of the result on which you specify the test.
- `t_value` The target value.

**Value Returned**

- `t` Returns `t` if the specification is set.
- `nil` Returns `nil` otherwise.

**Example**

`ocnxlPutTargetSpec("test_one" "VT(\'/out\')" "5.0") => t`

Spec is defined that transient voltage for /out signal must be 5.0 volts.
ocnxlResultsLocation

ocnxlResultsLocation(t_resultsDir)
    => t / nil

Description

Sets the results directory to the specified location. All results data goes into this location. By
default data goes into the target cell view. See help on ocnxlTargetCellView(). This
command works only in XL mode. See help on ocnSetXLMode().

Arguments

t_resultsDir Location of the results directory.

Value Returned

t Returns t if the results location is set.
nil Returns nil otherwise.

Example

ocnxlResultsLocation("/home/ocnuser")
ocnxlRun

ocnxlRun( [ ?nominalCornerEnabled b_nominalCornerEnabled ]
          [ ?allCornersEnabled b_allCornersEnabled ]
          [ ?allSweepsEnabled b_allSweepsEnabled ]
          => t / nil

Description

Lets you run the setup. It also lets you control what you want to run. Even if you have added many sweeps and corners, in this command, you can control whether to run the nominal corner, all corners and all sweeps.

By default, the entire setup runs, that is, all corners, all sweeps and nominal corner run. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

b_nominalCornerEnabled  Specifies if the nominal corner is enabled for run. This is an optional argument.

b_allCornersEnabled     Specifies if all corners are enabled for run. This is an optional argument.

b_allSweepsEnabled      Specifies if all sweeps are enabled for run. This is an optional argument.

Value Returned

t           Returns t if the run is scheduled.

nil          Returns nil otherwise.

Example

ocnxlRun(?allCornersEnabled nil)
No corner will be run but rest of the setup will be run.
ocnxlRunSetupSummary

ocnxlRunSetupSummary()
=> t / nil

Description

Generates the run setup summary. It shows how many tests, sweeps and corners are there and whether they are enabled. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

None.

Value Returned

t Returns t if the summary is generated.
nil Returns nil otherwise.

Example

ocnxlRunSetupSummary()
**ocnxlSamplingOptions**

```ocnxlSamplingOptions ( [ ?points t_numberOfPoints ])
  => t / nil
```

**Description**

Lets you specify options for sampling run. See help on `ocnxlRun()` for run modes. This command works only in XL mode. See help on `ocnSetXLMode()`.

**Arguments**

- `t_numberOfPoints` Specifies the number of points. The default value for points is 200.

**Value Returned**

- `t` Returns `t` if the options for the run are specified.
- `nil` Returns `nil` otherwise.

**Example**

```ocnxlSamplingOptions(?points "500")
  # Sets 500 as the number of points for sampling run.
```
ocnxlSetupLocation

ocnxlSetupLocation(t_setupDir)
    => t / nil

Description

Sets the setup directory to the specified location. All setup data goes into this location. By default data goes into the target cell view. See help on ocnxlTargetCellView(). This command works only in XL mode. See help on ocnSetXLMode().

Arguments

\texttt{t\_setupDir} \quad \text{Location of the setup directory.}

Value Returned

\texttt{t} \quad \text{Returns } t \text{ if the location of the setup directory is set.}

\texttt{nil} \quad \text{Returns } \text{nil} \text{ otherwise.}

Example

ocnxlSetupLocation("/home/ocnuser")
ocnxOutputExpr

ocnxOutputExpr( t_expr
  [ ?name t_outputName ]
  [ ?plot plot ]
  [ ?save save ]
)
=> t / nil

Description

This command adds an output expression in the current test being specified. A test's specification begins with ocnxBeginTest(). See help on ocnxBeginTest(). This command works only in XL mode. See help on ocnSetXLMode().

Arguments

- **t_expr**: The expression that you want to add.
- **t_outputName**: The name of the expression. This is an optional argument.
- **plot**: Whether to plot or not. This is an optional argument.
- **save**: Whether to save or not. This is an optional argument.

Value Returned

- **t**: Returns \( t \) if the output expression is set.
- **nil**: Returns nil otherwise.

Example

ocnxOutputExpr( "ymax(VT("/out"))" ?name "MAX" ?plot t ?save t )
# Adds "/out" in the outputs.
ocnxlOutputSignal

ocnxlOutputSignal( t_signal [ ?plot plot ] [ ?save save ] )
=> t / nil

Description

This command adds an output signal in the current test being specified. A test’s specification begins with ocnxlBeginTest(). See help on ocnxlBeginTest(). This command works only in XL mode. See help on ocnSetXLMode().

Arguments

t_signal The name of the signal.
plot Whether to plot or not. This is an optional argument.
save Whether to save or not. This is an optional argument.

Value Returned

t Returns t if the output signal is set.
nil Returns nil otherwise.

Example

ocnxlOutputSignal( "/out" ?plot t ?save t )
# Adds "/out" in the outputs.
ocnxlOutputTerminal

ocnxlOutputTerminal( t_term [ ?plot plot ] [ ?save save ] )
=> t / nil

Description

This command adds an output terminal in the current test being specified. A test's specification begins with ocnxlBeginTest(). See help on ocnxlBeginTest(). This command works only in XL mode. See help on ocnSetXLMode().

Arguments

- t_term: The name of the terminal.
- plot: Whether to plot or not. This is an optional argument.
- save: Whether to save or not. This is an optional argument.

Value Returned

- t: Returns t if the output terminal is set.
- nil: Returns nil otherwise.

Example

ocnxlOutputTerminal( "/I8/inp" ?plot t ?save t )
# Adds "/I8/inp" in the outputs.
ocnxlOutputSummary

ocnxlOutputSummary(?exprSummary b_exprSummary ?specSummary b_specSummary ?detailed b_detailed ?wave b_wave)
=> t/nil

Description

Generates output summary. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

b_exprSummary Default value for exprSummary is t. When exprSummary is t, then expressions summary is printed. If detailed is also set to t, then detailed expression summary is also printed. If you do not want to see the expressions summary, give nil for exprSummary.

b_specSummary Default value for specSummary is t. When specSummary is t, then spec summary is printed. If detailed is also set to t, then detailed spec summary is also printed. If you do not want to see the spec summary, give nil for specSummary.

b_detailed Default value for detailed is t. When detailed is t, then details of expr/spec are printed. Otherwise only summary is printed.

b_wave Default value of wave is t. When wave is t, then the value of expressions evaluating to a waveform is printed as "wave". If you do not want to see the expressions that are evaluated to waveforms in the output, set wave to nil.

Value Returned

t Returns t if the summary is generated.

nil Returns nil otherwise.

Example

ocnxlOutputSummary()
# This will print the details of expressions and specs for each sweep point and each corners. It will also print the summary of expressions (minimum and maximum values) and for specs (pass/fail, minimum and maximum values) for each sweep point across all corners.

`ocnxlOutputSummary(?exprSummary nil)`

# This will print the details of specs for each sweep point and each corners. It will also print the summary of specs (pass/fail, minimum and maximum values) for each sweep point across all corners. This will not print any expression details/summary.

`ocnxlOutputSummary(?specSummary nil ?detailed nil)`

# This will print only the summary of expressions (minimum and maximum values) for each sweep point across all corners. This will not print any details for expressions. This will also not print any spec details/summary.

`ocnxlOutputSummary(?wave nil)`

# This will print the details of expressions and specs for each sweep point and each corners. It will also print the summary of expressions (minimum and maximum values) and for specs (pass/fail, minimum and maximum values) for each sweep point across all corners. This will skip printing the outputs that evaluate to waveforms.
ocnxlTargetCellView

ocnxlTargetCellView(t_lib t_cell t_view)
  => t / nil

Description

Specifies target cellview where ADE XL data will be created. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

  t_lib          Name of the library.
  t_cell         Name of the cell.
  t_view         Name of the view.

Value Returned

  t             Returns t if its able to use the lib:cell:view as the target.
  nil           Returns nil otherwise.

Example

  ocnxlTargetCellView("opamplib" "ampTest" "schematic")
ocnxlYieldImprovementOptions

ocnxlYieldImprovementOptions([?iymethod t_iymethod]

[?refPoint t_refPoint] [?algorithm t_algorithm] [?timeLimit t_timeLimit]

[?iterations t_iterations] [?numPoints t_numPoints]

[?sigmaTarget t_sigmaTarget] [?stopIfNoImprovement t_stopIfNoImprovement]

[?runFullEvaluation t_runFullEvaluation]

[?optimizationMethod t_optimizationMethod] [?effort t_effort]

[?iysamplingmethod t_iysamplingmethod]

[?iymontecarlodomoninal t_iymontecarlodomoninal]

[?iymontecarloseed t_iymontecarloseed] [?iymcnumpoints t_iymcnumpoints]

[?iymontecarlostartingrun t_iymontecarlostartingrun]

[?WYCmethod t_WYCmethod] [?designUnderTest t_designUnderTest]

[?dutInstances t_dutInstances])

=> t / nil

Description

Lets you specify options for improve yield runs. See help on ocnxlRun() for run modes. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

**t_iymethod**

The yield improvement method to be used. The default value is "all". The possible values are "global", "mismatch" and "all".

**t_refPoint**

Specifies whether to use a schematic point or a reference point that you have created as a starting place for sizing. Possible values are "0" and "1". The default value is "0".

**t_algorithm**

The default value is "0". The possible values are "0" and "1".

**t_timeLimit**

Sets a time limit for the improve yield run. The default value is "". The timeLimit is in minutes.

**t_iterations**

Specifies the number of sizing/Monte Carlo iterations, The default value is "3".
**t_numPoints**
Specifies the maximum number of points processed per iteration.
The default value is "3000".

**t_sigmaTarget**
Allows you to increase the mean of the goal distribution to target (of goal) value even after achieving 100% yield. ADE GXL allows you to achieve 4, 5, or even 6 sigma designs.
The default value is "6". The possible values are "4", "5", and "6".

**t_stopIfNoImprovement**
Specifies if the yield improvement run must be stopped if there is no yield improvement. The default value is "0". The possible values are "0" and "1".

**t_runFullEvaluation**
Sets the program to run full optimization. Default for runFullEvaluation is "0". Possible values are "0" and "1".

**t_optimizationMethod**
Sets the optimization method. Default for optimizationMethod is "global". Possible values are "global" and "local".

**t_effort**
Specifies the effort level if you are using local optimization. Default for effort is "fine". Possible values are "fine" and "coarse".

**t_iysamplingmethod**
Sets the default statistical sampling method for improve yield runs. The default value is "random". Possible values are "random" and "lhs".

**t_iymontecarlonominal**
Specifies whether to run a simulation at the reference point prior to beginning the improve yield process. Possible values are "0" and "1". If set to "1", Spectre will run a simulation at the reference point, and, if this fails, then the sampling process is not initiated and the simulation stops.
The default value is "1".

**t_iymontecarlosseed**
Specifies a different seed for Monte Carlo runs. Default for monteCarloSeed is "12345".

**t_iymcnumpoints**
Sets the number of Monte Carlo points you want to simulate. The default value is nil.

**t_iymontecarlostaringrun**
Specifies the run that Monte Carlo begins with. The default value is "1".
**t_designUnderTest**  
Specifies the design under test (DUT) for improve yield runs. Default for `designUnderTest` is "". Specify the DUT using the format:  

```
"libname:cellname:viewname"
```

**t_dutInstances**  
Specifies the DUT instances for improve yield runs. Default for `dutInstances` is "". Specify the DUT instances using the characters %/ to separate two DUT instances. For example, specify:  

```
"libname:cellname:instname1%/libname:cellname:instname2"
```

**Value Returned**

- **t**: Returns `t` if the options are specified.
- **nil**: Returns `nil` otherwise.

**Example**

```ruby
ocnxlYieldImprovementOptions(?iymethod "mismatch")  
# Sets mismatch as the method for yield improvement run.
```
ocnxlEnableCornerForTest

ocnxlEnableCornerForTest(t_cornerName t_testName)
 => t / nil

Description

Enables a corner for a test. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

<table>
<thead>
<tr>
<th>t_cornerName</th>
<th>Name of the corner.</th>
</tr>
</thead>
<tbody>
<tr>
<td>t_testName</td>
<td>Name of the test.</td>
</tr>
</tbody>
</table>

Value Returned

<table>
<thead>
<tr>
<th>t</th>
<th>Returns t if successful.</th>
</tr>
</thead>
<tbody>
<tr>
<td>nil</td>
<td>Returns nil otherwise.</td>
</tr>
</tbody>
</table>

Example

ocnxlEnableCornerForTest("C0" "test_one") => t
ocnlEnableSweepParam

ocnlEnableSweepParam(t_paramPath)
   => t / nil

Description

Enables a sweep parameter. A disabled sweep parameter is not run when the ocnlRun() command is run. This command works only in XL mode. See help on ocnlSetXLMode().

Arguments

t_paramPath 
   Name of the sweep parameter.

Value Returned

   t 
   Returns t if successful.

   nil 
   Returns nil otherwise.

Example

ocnlEnableSweepParam("solutions/ampTest/schematic/R1/r") => t
ocnxlEnableSweepVar

ocnxlEnableSweepVar(t_varName)
   => t / nil

Description

Enables a sweep variable. A disabled sweep variable is not run when the ocnxlRun() command is run. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

t_varName  Name of the sweep variable.

Value Returned

<table>
<thead>
<tr>
<th>t</th>
<th>Returns t if successful.</th>
</tr>
</thead>
<tbody>
<tr>
<td>nil</td>
<td>Returns nil otherwise.</td>
</tr>
</tbody>
</table>

Example

ocnxlEnableSweepVar("CAP") => t
ocnxlEnableTest

ocnxlEnableTest(t_testName)
    => t / nil

Description

Enables a test. A disabled test will not be run when the ocnxlRun() command is run. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

t_testName                Name of the test.

Value Returned

t Returns t if successful.
nil Returns nil otherwise.

Example

ocnxlEnableTest("test_two") => t
ocnxlGetBestPointParams

ocnxlGetBestPointParams()
   => t / nil

**Description**

Returns a list of best design points. This command works only in XL mode. See help on ocnSetXLMode().

**Arguments**

None

**Value Returned**

<table>
<thead>
<tr>
<th>Value</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>t</td>
<td>Returns t if successful.</td>
</tr>
<tr>
<td>nil</td>
<td>Returns nil otherwise.</td>
</tr>
</tbody>
</table>

**Example**

ocnxlGetBestPointParams()
ocnxlGetCorners

ocnxlGetCorners() => t / nil

Description

Returns a list of corners names. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

None

Value Returned

- **t**: Returns *t* if successful.
- **nil**: Returns *nil* otherwise.

Example

ocnxlGetCorners()
ocnxlGetCurrentHistory

ocnxlGetCurrentHistory()
   => t / nil

Description

Returns the current history (checkpoint) name. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

None

Value Returned

t Returns t if successful.

nil Returns nil otherwise.

Example

ocnxlGetCurrentHistory()
ocnxlGetCurrentHistoryId

ocnxlGetCurrentHistoryId()
=> t / nil

Description

Returns the ID of the current history (checkpoint). This command works only in XL mode. See help on ocnSetXLMode().

Arguments

None

Value Returned

<table>
<thead>
<tr>
<th>t</th>
<th>Returns t if successful.</th>
</tr>
</thead>
<tbody>
<tr>
<td>nil</td>
<td>Returns nil otherwise.</td>
</tr>
</tbody>
</table>

Example

ocnxlGetCurrentHistoryId()
ocnxlGetSession

ocnxlGetSession()

=> t / nil

Description

Returns the session name. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

None

Value Returned

<table>
<thead>
<tr>
<th>Value</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>t</td>
<td>Returns t if successful.</td>
</tr>
<tr>
<td>nil</td>
<td>Returns nil otherwise.</td>
</tr>
</tbody>
</table>

Example

ocnxlGetSession()
ocnxlGetSpecs

ocnxlGetSpecs()
 => t / nil

Description

Returns a list of parameter specification names. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

None

Value Returned

<table>
<thead>
<tr>
<th>t</th>
<th>Returns t if successful.</th>
</tr>
</thead>
<tbody>
<tr>
<td>nil</td>
<td>Returns nil otherwise.</td>
</tr>
</tbody>
</table>

Example

ocnxlGetSpecs()
ocnxlGetTests

ocnxlGetTests()
  => t / nil

Description

Returns a list of test names. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

None

Value Returned

- t: Returns t if successful.
- nil: Returns nil otherwise.

Example

ocnxlGetTests()
ocnxlRemoveSpec

ocnxlRemoveSpec(t_specName)
  => t / nil

Description

Removes the specified parameter specification. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

  t_specName                Name of the spec.

Value Returned

  t                          Returns t if successful.
  nil                        Returns nil otherwise.

Example

  ocnxlRemoveSpec("MAX")
ocnxlRenameCurrentHistory

ocnxlRenameCurrentHistory(t_newNameForHistory)
  => t / nil

Description

Renames the current history (checkpoint). This command works only in XL mode. See help on ocnSetXLMode().

Arguments

  t_newNameForHistory  New name for the current history.

Value Returned

  t  Returns t if successful.
  nil  Returns nil otherwise.

Example

  ocnxlRenameCurrentHistory("myHistory")
ocnxlRun

ocnxlRun( [ ?mode s_mode ]
    [ ?nominalCornerEnabled b_nominalCornerEnabled ]
    [ ?allCornersEnabled b_allCornersEnabled ]
    [ ?allSweepsEnabled b_allSweepsEnabled ]
    [ ?verboseMode b_verboseMode ]
)
=> t/nil

Description

Specifies the run mode for simulation and whether to run the nominal corner, corners and
sweeps during simulation. Also specifies whether to report completion of points during
simulation. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

\textit{s\_mode} \quad \text{Lets you run simulations in one of following modes:}
\begin{itemize}
  \item 'sweepsAndCorners
  \item 'localOptimization
  \item 'globalOptimization
  \item 'monteCarlo
  \item 'yieldImprovement
  \item 'sampling
  \item 'sensitivity
\end{itemize}

\textit{b\_nominalCornerEnabled} \quad \text{Accepts boolean values \texttt{t} or \texttt{nil}. The default value is \texttt{t}. If set to \texttt{nil}, ADE XL excludes nominal corners from the simulation run.}

\textit{b\_allCornersEnabled} \quad \text{Accepts boolean values \texttt{t} or \texttt{nil}. The default value is \texttt{t}. If set to \texttt{nil}, ADE XL excludes all corners from the simulation run.}

\textit{b\_allSweepsEnabled} \quad \text{Accepts boolean values \texttt{t} or \texttt{nil}. The default value is \texttt{t}. If set to \texttt{nil}, ADE XL excludes all sweeps from the simulation run.}
b_verboseMode

Accepts boolean values t or nil. The default value is t. If set to nil, ADE XL does not report the progress in the simulation of points in the simulation run.

**Note:** Cadence recommends that you specify the value nil if you have setup a large number of points.

**Value Returned**

<table>
<thead>
<tr>
<th>Value</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>t</td>
<td>Returns t if successful.</td>
</tr>
<tr>
<td>nil</td>
<td>Returns nil otherwise.</td>
</tr>
</tbody>
</table>

**Example**

```ocnxlRun(?allCornersEnabled nil)
No corner will be run but rest of the ADE XL setup will be run.ocnxlHistoryPrefix```
ocnxlHistoryPrefix

ocnxlHistoryPrefix(t_prefixName)
  => t / nil

Description

Sets the prefix used in the names of history items created by OCEAN XL runs. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

  t_prefixName  The prefix to be used in the names of history items.

Value Returned

  t  Returns t if successful.
  nil Returns nil otherwise.

Example

  ocnxlHistoryPrefix("check")

  Creates history items with names like check.0, check.1, and so on.
ocnxILoadSetupState

ocnxILoadSetupState( t_stateName t_mode [ ?tests t_tests ]
[ ?vars t_vars ] [ ?parameters t_parameters ] [ ?currentMode t_currentMode ]
[ ?allSweepsEnabled t_allSweepsEnabled ]
[ ?allCornersEnabled t_allCornersEnabled ]
[ ?defaultCornerEnabled t_defaultCornerEnabled ]
[ ?runOptions t_runOptions ] [ ?specs t_specs ] [?corners t_corners ]
[ ?modelGroups t_modelGroups ] [?extensions t_extensions]
)
=> t / nil

Description

Restores the settings in the specified setup state to the active setup. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

\textbf{t\_stateName} \hspace{2cm} The name of the setup state to be restored.

\textbf{t\_mode} \hspace{2cm} Specifies the mode for restoring the settings in the setup state to the active setup.

Valid values: \textquoteleft retain, \textquoteleft merge, \textquoteleft overwrite

\textbf{t\_tests} \hspace{2cm} Specifies whether the tests in the setup state should be restored to the active setup.

Valid values: \textit{t, nil}

Default Value: \textit{t}

\textbf{t\_vars} \hspace{2cm} Specifies whether the global variables in the setup state should be restored to the active setup.

Valid values: \textit{t, nil}

Default Value: \textit{t}

\textbf{t\_parameters} \hspace{2cm} Specifies whether the parameters in the setup state should be restored to the active setup.

Valid values: \textit{t, nil}

Default Value: \textit{t}
### t_currentMode
Specifies whether the run mode in the setup state should be restored to the active setup

- **Valid values:** t, nil
- **Default Value:** t

### t_allSweepsEnabled
Specifies whether the run mode settings in the setup state should be restored to the active setup

- **Valid values:** t, nil
- **Default Value:** t

### t_allCornersEnabled
Specifies whether the run mode settings in the setup state should be restored to the active setup

- **Valid values:** t, nil
- **Default Value:** t

### t_defaultCornerEnabled
Specifies whether the run mode settings in the setup state should be restored to the active setup

- **Valid values:** t, nil
- **Default Value:** t

### t_runOptions
Specifies whether the run options in the setup state should be restored to the active setup

- **Valid values:** t, nil
- **Default Value:** t

### t_specs
Specifies whether the parameter specifications in the setup state should be restored to the active setup

- **Valid values:** t, nil
- **Default Value:** t

### t_corners
Specifies whether the corners in the setup state should be restored to the active setup

- **Valid values:** t, nil
- **Default Value:** t
**t_modelGroups**

Specifies whether the model groups in the setup state should be restored to the active setup.

Valid values: t, nil

Default Value: t

**t_extensions**

Specifies whether the extensions in the setup state should be restored to the active setup.

Valid values: t, nil

Default Value: t

---

**Value Returned**

- **t**
  - Returns t if successful.

- **nil**
  - Returns nil otherwise.

---

**Example**

```
ocnxlLoadSetupState("optimize")
```

Loads the setup state named optimize.
ocnxlStartingPoint

ocnxlStartingPoint(l_startingPointDetails)
    => t / nil

Description

Lets you specify a reference point—a starting place for sizing—for Improve Yield, Global Optimization or Monte Carlo runs. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

l_startingPointDetails A list of elements where each element is:

    (t_type t_varName t_value)

Where:

■ t_type can be a variable or parameter.
■ t_varName is the name of the variable or parameter
■ t_value is the value of the variable or parameter.

Value Returned

t Returns t if successful.
nil Returns nil otherwise.

Example

ocnxlStartingPoint('(("variable" "CAP" "2p")
    ("parameter" "ether_adcflash/adc_cascode_opamp/schematic/M2/fw" "16.3u"))}
ocnxlOutputAreaGoal

ocnxlOutputAreaGoal( t_expr
    [ ?name t_outputName ]
    [ ?plot plot ]
    [ ?save save ]
)
=> t / nil

Description

Adds an area goal output expression in the current test being specified. A test’s specification begins with `ocnxlBeginTest()`. See help on `ocnxlBeginTest()`. This command works only in XL mode. See help on `ocnSetXLMode()`.

Arguments

- **t_expr**
  The expression that you want to add.
- **t_outputName**
  The name of the expression. This is an optional argument.
- **plot**
  Whether to plot or not. This is an optional argument.
- **save**
  Whether to save or not. This is an optional argument.

Value Returned

- **t**
  Returns t if the output expression is set.
- **nil**
  Returns nil otherwise.

Example

```
ocnxlOutputAreaGoal( "((I8/R4) ('res' 'w*l' 'default' 'enabled'))...)" ?name "MAX" ?plot t ?save t ) => t
```

Adds an area goal output expression named `MAX` for the current test.
ocnxlConjugateGradientOptions

ocnxlConjugateGradientOptions (  
    [?runFullEvaluation t_runFullEvaluation]  
    [?meetAllGoals t_meetAllGoals]  
    [?timeLimit t_timeLimit]  
    [?numPoints t_numPoints]  
)  
=> t | nil

Description

Sets options for conjugate gradient runs. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

  t_runFullEvaluation  Sets to run full evaluation.  
                      Possible Values: “1” and “0”.  
                      Default Value: “1”
  t_meetAllGoals  Sets to run only until all goals are met.  
                 Possible Values: “1” and “0”.  
                 Default Value: “1”
  t_timeLimit  Sets a time limit (in seconds) for the run.
  t_numPoints  Sets the limit in the number of points to be run.

Value Returned

  t  Returns t if options are specified for conjugate gradient run.
  nil  Returns nil otherwise.

Example

  ocnxlConjugateGradientOptions(?numPoints "3000")
  Sets to run for 3000 points.
ocnxlMTSEnable

ocnxlMTSEnable( b_enable)

=> t / nil

Description

Enables or disables multi-technology simulation (MTS) mode for the current test. The current
test’s specification begins with ocnxlBeginTest(). See help on ocnxlBeginTest().
This command works only in XL mode. See help on ocnSetXLMode().

Arguments

b_enable

Enables or disables MTS mode for the current test.

Valid values: t, nil

Default Value: nil

Value Returned

t

Returns t if successful.

nil

Returns nil otherwise.

Example

ocnxlMTSEnable(t)

Enables MTS mode for the current test.
ocnxIMTSBlock

ocnxIMTSBlock( s_blockName
            [?isMtsBlock  b_isMtsBlock]
            [?includeFile  l_includeFile]
            [?modelFiles  l_modelFiles]
         )
     => t/nil

Description

Enables a block for multi-technology simulation (MTS) and specifies the include files and model files associated with the block. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

\textit{s\_blockName} \hspace{1cm} \text{Specifies the name of the block that needs to be enabled for multi-technology simulation.}
\textit{Valid values: a string}

\textit{b\_isMtsBlock} \hspace{1cm} \text{Specifies whether the block is enabled or disabled for multi-technology simulation.}
\textit{Valid values: t, nil}
\textit{Default Value: nil}

\textit{l\_includeFile} \hspace{1cm} \text{Specifies the include files associated with the block.}
\textit{Valid values: a list of strings or nil}
\textit{Default value: nil}

\textit{l\_modelFiles} \hspace{1cm} \text{Specifies the model files associated with the block.}
\textit{Valid values: a list of strings or nil}
\textit{Default value: nil}

Value Returned

\textit{t} \hspace{1cm} \text{Returns t if successful.}
\textit{nil} \hspace{1cm} \text{Returns nil otherwise.}
Example

ocnx1MTSBlock('digLib\ inv_usim
?isMtsBlock t
?modelFiles '(['Models/myModels.scs" "ss"
("Models/spectre_cl0131v.scs" "aa"))

Enables the inv_usim cell in the digLib library for multi-technology simulation and specifies the model files (and the sections of the model files) associated with the block.
ocnxlProjectDir

ocnxlProjectDir(t_projectDir)
    => t / nil

Description

Sets the project directory to the specified location. All simulation data goes into this location by default. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

    t_projectDir      Sets the location of the project directory.

Value Returned

    t                Returns t if successful.
    nil              Returns nil otherwise.

Example

    ocnxlProjectDir("/home/ocnuser/simulation")
ocnxlSimResultsLocation

ocnxlSimResultsLocation(t_simResultsDir)
   => t / nil

Description

Sets the simulation results directory to the specified location. All simulation data goes into this location. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

t_simResultsDir  Sets the location of the simulation results directory.

Value Returned

t  Returns t if successful.
nil  Returns nil otherwise.

Example

ocnxlSimResultsLocation("/home/ocnuser")

ocnxlDisableCorner

ocnxlDisableCorner(t_cornerName)
   => t/nil

Description

Lets you disable a corner. A disabled corner will not be run when the ocnxlRun() command is run. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

t_cornerName  The name of the corner to be disabled.
Value Returned

\[
\begin{align*}
t & \quad \text{Returns } t \text{ if the corner is disabled.} \\
nil & \quad \text{Returns } \text{nil} \text{ otherwise.}
\end{align*}
\]

Example

\[
\text{ocexlDisableCorner("C0") } \Rightarrow t
\]
ocnxlEnableCorner

ocnxlEnableCorner(t_cornerName)
    => t/nil

Description

Lets you enable a corner. An enabled corner will be run when ocnxlRun() command is run. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

t_cornerName The name of the corner to be enabled.

Value Returned

t Returns t if the corner is disabled.
nil Returns nil otherwise.

Example

ocnxlEnableCorner("C0") => t
**ocnxlSaveSetupAs**

`ocnxlSaveSetupAs(t_lib t_cell t_view)`

=> `t/nil`

**Description**

Saves the current setup to a different adexl view. This command works only in XL mode. See help on `ocnSetXLMode()`.

**Arguments**

- `t_lib`: The name of the library in which the new adexl view is to be saved.
- `t_cell`: The name of the cell in which the new adexl view is to be saved.
- `t_view`: The name of the the new adexl view.

**Value Returned**

- `t`: Returns `t` if the save is successful.
- `nil`: Returns `nil` otherwise.

**Example**

`ocnxlSaveSetupAs("solution" "ampTest" "newView")`
ocnxlSetAllParametersDisabled

ocnxlSetAllParametersDisabled(g_disabled)
   => t/nil

Description

Lets you enable or disable all parameters. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

  g_disabled  Specify t to disable all parameters, and nil to enable all parameters.

Value Returned

  t  Returns t if all parameters are enabled or disabled.
  nil  Returns nil otherwise.

Example

  ocnxlSetAllParametersDisabled(t) => t
ocnxlSetAllVarsDisabled

ocnxlSetAllVarsDisabled(g_disabled)
  => t/nil

Description

Lets you enable or disable all global variables. This command works only in XL mode. See help on ocnSetXLMode().

Arguments

  g_disabled

  Specify t to disable all variables, and nil to enable all variables.

Value Returned

  t

  Returns t if all variables are enabled or disabled.

  nil

  Returns nil otherwise.

Example

  ocnxlSetAllVarsDisabled(t) => t
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.
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:

```plaintext
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.
# Index

## Symbols

<table>
<thead>
<tr>
<th>Symbol</th>
<th>Page</th>
</tr>
</thead>
<tbody>
<tr>
<td>&amp;&amp; (and) operator</td>
<td>55</td>
</tr>
<tr>
<td>... in syntax</td>
<td>20</td>
</tr>
<tr>
<td>... in syntax</td>
<td>20</td>
</tr>
<tr>
<td>/ in syntax</td>
<td>20</td>
</tr>
<tr>
<td>[] in syntax</td>
<td>19</td>
</tr>
<tr>
<td>{} in syntax</td>
<td>19</td>
</tr>
<tr>
<td></td>
<td>in syntax</td>
</tr>
<tr>
<td></td>
<td></td>
</tr>
</tbody>
</table>

## Numerics

<table>
<thead>
<tr>
<th>Number</th>
<th>Page</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>20</td>
</tr>
<tr>
<td>2</td>
<td>21</td>
</tr>
</tbody>
</table>

## A

<table>
<thead>
<tr>
<th>Function/Expression</th>
<th>Page</th>
</tr>
</thead>
<tbody>
<tr>
<td>abs</td>
<td>256</td>
</tr>
<tr>
<td>abs function</td>
<td>256</td>
</tr>
<tr>
<td>ac</td>
<td>83</td>
</tr>
<tr>
<td>acos</td>
<td>257</td>
</tr>
<tr>
<td>add1</td>
<td>258</td>
</tr>
<tr>
<td>addSubwindow</td>
<td>193</td>
</tr>
<tr>
<td>addSubwindowTitle</td>
<td>194</td>
</tr>
<tr>
<td>addTitle</td>
<td>195</td>
</tr>
<tr>
<td>addWaveLabel</td>
<td>196</td>
</tr>
<tr>
<td>addWindowLabel</td>
<td>198</td>
</tr>
<tr>
<td>aliases</td>
<td>247</td>
</tr>
<tr>
<td>Allocating an Array of a Given Size</td>
<td>66</td>
</tr>
<tr>
<td>alphalessp function</td>
<td>67</td>
</tr>
<tr>
<td>alphaNumCmp function</td>
<td>68</td>
</tr>
<tr>
<td>analysis</td>
<td>85</td>
</tr>
<tr>
<td>Appending a maximum number of characters from two input strings (strncat)</td>
<td>67</td>
</tr>
<tr>
<td>appendString</td>
<td>74</td>
</tr>
<tr>
<td>arithmetic</td>
<td></td>
</tr>
<tr>
<td>operators</td>
<td>52</td>
</tr>
<tr>
<td>predefined functions</td>
<td>254</td>
</tr>
<tr>
<td>Arithmetic and Logical Expressions</td>
<td>59</td>
</tr>
<tr>
<td>Arithmetic Operators</td>
<td>52</td>
</tr>
<tr>
<td>Arrays</td>
<td>66</td>
</tr>
<tr>
<td>arrays</td>
<td></td>
</tr>
</tbody>
</table>

## B

<table>
<thead>
<tr>
<th>Function/Expression</th>
<th>Page</th>
</tr>
</thead>
<tbody>
<tr>
<td>b1f</td>
<td>284</td>
</tr>
<tr>
<td>bandwidth</td>
<td>285</td>
</tr>
<tr>
<td>binary minus operator</td>
<td>57</td>
</tr>
<tr>
<td>Blocking and Nonblocking Modes</td>
<td>35</td>
</tr>
<tr>
<td>Blocking Mode</td>
<td>35</td>
</tr>
<tr>
<td>braces in syntax</td>
<td>19</td>
</tr>
<tr>
<td>brackets in syntax</td>
<td>19</td>
</tr>
<tr>
<td>buildString function</td>
<td>66</td>
</tr>
</tbody>
</table>

## C

<table>
<thead>
<tr>
<th>Function/Expression</th>
<th>Page</th>
</tr>
</thead>
<tbody>
<tr>
<td>C language comparison</td>
<td></td>
</tr>
<tr>
<td>escape characters</td>
<td>66</td>
</tr>
<tr>
<td>parentheses</td>
<td>58</td>
</tr>
<tr>
<td>strings</td>
<td>65</td>
</tr>
<tr>
<td>case</td>
<td>483</td>
</tr>
<tr>
<td>case statement</td>
<td>483</td>
</tr>
<tr>
<td>clearAll</td>
<td>199</td>
</tr>
<tr>
<td>clearSubwindow</td>
<td>200</td>
</tr>
<tr>
<td>clip</td>
<td>286</td>
</tr>
<tr>
<td>clip function</td>
<td>286</td>
</tr>
<tr>
<td>close</td>
<td>488</td>
</tr>
<tr>
<td>close function</td>
<td>488</td>
</tr>
<tr>
<td>command types</td>
<td>26</td>
</tr>
<tr>
<td>commands</td>
<td></td>
</tr>
<tr>
<td>data access</td>
<td></td>
</tr>
<tr>
<td>DataTypes</td>
<td>153</td>
</tr>
<tr>
<td>getData</td>
<td>154</td>
</tr>
<tr>
<td>i</td>
<td>157</td>
</tr>
<tr>
<td>noiseSummary</td>
<td>219</td>
</tr>
<tr>
<td>ocnHelp</td>
<td>159</td>
</tr>
<tr>
<td>ocnPrint</td>
<td>223, 226, 228</td>
</tr>
<tr>
<td>ocnResetResults</td>
<td>161</td>
</tr>
<tr>
<td>Function</td>
<td>Page Numbers</td>
</tr>
<tr>
<td>----------</td>
<td>--------------</td>
</tr>
<tr>
<td>openResults</td>
<td>34, 162</td>
</tr>
<tr>
<td>outputParams</td>
<td>164</td>
</tr>
<tr>
<td>outputs</td>
<td>166</td>
</tr>
<tr>
<td>pv</td>
<td>170</td>
</tr>
<tr>
<td>report</td>
<td>239</td>
</tr>
<tr>
<td>results</td>
<td>174</td>
</tr>
<tr>
<td>selectResult</td>
<td>175</td>
</tr>
<tr>
<td>sweepNames</td>
<td>179</td>
</tr>
<tr>
<td>sweepValues</td>
<td>181</td>
</tr>
<tr>
<td>v</td>
<td>184</td>
</tr>
<tr>
<td>plotting</td>
<td></td>
</tr>
<tr>
<td>addSubwindow</td>
<td>193</td>
</tr>
<tr>
<td>addSubwindowTitle</td>
<td>194</td>
</tr>
<tr>
<td>addTitle</td>
<td>195</td>
</tr>
<tr>
<td>addWaveLabel</td>
<td>196</td>
</tr>
<tr>
<td>addWindowLabel</td>
<td>198</td>
</tr>
<tr>
<td>clearAll</td>
<td>199</td>
</tr>
<tr>
<td>clearSubwindow</td>
<td>200</td>
</tr>
<tr>
<td>currentSubwindow</td>
<td>201</td>
</tr>
<tr>
<td>currentWindow</td>
<td>202</td>
</tr>
<tr>
<td>dbCompressionPlot</td>
<td>203</td>
</tr>
<tr>
<td>deleteSubwindow</td>
<td>204</td>
</tr>
<tr>
<td>deleteWaveform</td>
<td>209</td>
</tr>
<tr>
<td>displayMode</td>
<td>210, 211</td>
</tr>
<tr>
<td>graphicsOff</td>
<td>212</td>
</tr>
<tr>
<td>graphicsOn</td>
<td>213</td>
</tr>
<tr>
<td>hardCopy</td>
<td>214</td>
</tr>
<tr>
<td>hardCopyOptions</td>
<td>215</td>
</tr>
<tr>
<td>ip3Plot</td>
<td>217</td>
</tr>
<tr>
<td>newWindow</td>
<td>218</td>
</tr>
<tr>
<td>plot</td>
<td>230</td>
</tr>
<tr>
<td>plotStyle</td>
<td>233</td>
</tr>
<tr>
<td>removeLabel</td>
<td>238</td>
</tr>
<tr>
<td>xLimit</td>
<td>242</td>
</tr>
<tr>
<td>yLimit</td>
<td>243</td>
</tr>
<tr>
<td>return values</td>
<td>31</td>
</tr>
<tr>
<td>simulation</td>
<td></td>
</tr>
<tr>
<td>ac</td>
<td>83</td>
</tr>
<tr>
<td>analysis</td>
<td>85</td>
</tr>
<tr>
<td>appendPath</td>
<td>74</td>
</tr>
<tr>
<td>createFinalNetlist</td>
<td>89, 93, 105, 107</td>
</tr>
<tr>
<td>dc</td>
<td>96</td>
</tr>
<tr>
<td>delete</td>
<td>99</td>
</tr>
<tr>
<td>design</td>
<td>101</td>
</tr>
<tr>
<td>desVar</td>
<td>103</td>
</tr>
<tr>
<td>envOption</td>
<td>108</td>
</tr>
<tr>
<td>forcenode</td>
<td>112, 113, 114</td>
</tr>
<tr>
<td>ic</td>
<td>116, 125</td>
</tr>
<tr>
<td>includeFile</td>
<td>117</td>
</tr>
<tr>
<td>nodeset</td>
<td>119</td>
</tr>
<tr>
<td>noise</td>
<td>120</td>
</tr>
<tr>
<td>ocnDisplay</td>
<td>122</td>
</tr>
<tr>
<td>off</td>
<td>126</td>
</tr>
<tr>
<td>option</td>
<td>127</td>
</tr>
<tr>
<td>paramAnalysis</td>
<td>33, 448</td>
</tr>
<tr>
<td>paramRun</td>
<td>452</td>
</tr>
<tr>
<td>path</td>
<td>75, 79, 80</td>
</tr>
<tr>
<td>prependPath</td>
<td>76</td>
</tr>
<tr>
<td>restore</td>
<td>129</td>
</tr>
<tr>
<td>resultsDir</td>
<td>130</td>
</tr>
<tr>
<td>run</td>
<td>131</td>
</tr>
<tr>
<td>save</td>
<td>135</td>
</tr>
<tr>
<td>simulator</td>
<td>139, 140</td>
</tr>
<tr>
<td>store</td>
<td>143</td>
</tr>
<tr>
<td>temp</td>
<td>144</td>
</tr>
<tr>
<td>tran</td>
<td>145</td>
</tr>
<tr>
<td>commenting code</td>
<td>57</td>
</tr>
<tr>
<td>Comments</td>
<td>57</td>
</tr>
<tr>
<td>Common SKILL Syntax Characters Used In OCEAN</td>
<td>27</td>
</tr>
<tr>
<td>compare</td>
<td>288</td>
</tr>
<tr>
<td>Comparing Strings</td>
<td>67</td>
</tr>
<tr>
<td>Comparing Two String or Symbol Names Alphanumerically or Numerically (alphaNumCmp)</td>
<td>68</td>
</tr>
<tr>
<td>Comparing Two Strings Alphabetically (strcmp)</td>
<td>68</td>
</tr>
<tr>
<td>Comparing Two Strings or Symbol Names Alphabetically (alphalessp)</td>
<td>67</td>
</tr>
<tr>
<td>complex</td>
<td>296</td>
</tr>
<tr>
<td>complexp</td>
<td>297</td>
</tr>
<tr>
<td>compression</td>
<td>290</td>
</tr>
<tr>
<td>compressionVRI</td>
<td>292</td>
</tr>
<tr>
<td>compressionVRICurves</td>
<td>294</td>
</tr>
<tr>
<td>Concatenating a list of strings with separation characters (buildString)</td>
<td>66</td>
</tr>
<tr>
<td>Concatenating Strings (Lists)</td>
<td>66</td>
</tr>
<tr>
<td>Concatenating two or more input strings (strcat)</td>
<td>67</td>
</tr>
<tr>
<td>cond</td>
<td>485</td>
</tr>
<tr>
<td>cond statement</td>
<td>485</td>
</tr>
<tr>
<td>conjugate</td>
<td>298</td>
</tr>
<tr>
<td>conjugate function</td>
<td>298</td>
</tr>
<tr>
<td>Constants</td>
<td>60</td>
</tr>
<tr>
<td>constants</td>
<td>60</td>
</tr>
<tr>
<td>Constants and Variables</td>
<td>65</td>
</tr>
<tr>
<td>Convention</td>
<td>28, 29, 30</td>
</tr>
<tr>
<td>conventions for user-defined arguments</td>
<td>19</td>
</tr>
<tr>
<td>for user-entered text</td>
<td>19</td>
</tr>
<tr>
<td>convolve</td>
<td>299</td>
</tr>
</tbody>
</table>
convolve function 299
cos 261
cPwrContour 301
createFinalNetlist 89, 93, 105, 107
createNetlist 94
Creating Arithmetic and Logical Expressions 61
Creating OCEAN Scripts 42
Creating Scripts from Analog Artist 42
Creating Scripts from the Analog Design Environment 42
Creating Scripts Using Sample Script Files 42
cRefContour 303
cross 305
currentSubwindow 201
currentWindow 202

data access commands. See commands, data access
Data Access Without Running a Simulation 34
Data Types 63
data types
   SKILL 30
   supported 63
Data Types Used in OCEAN 30
dataTypes 153
db10 307
db20 308
dbCompressionPlot 203
dbm 309
dc 96
declare function 66
Declaring a SKILL Function 68
Defining Function Parameters 69
Defining Local Variables (let) 69
definitionFile 98
delay 310
delete 99
deleteJob 456
deleteSubwindow 204, 208
deleteWaveform 209
deriv 314
design 101
design variables 31
Design Variables in OCEAN 31
desVar 103
dft 315, 317
dftbb 317
discipline 105
displayMode 210, 211
displayNetlist 107
Distributed Processing 34
dnl 319
double quotes 28

E

envOption 108
Errors and Warnings 481
evcdFile 110
evcdInfoFile 111
evmQAM 323
evmQpsk 325
exp 262, 263
expressions, nested 58
eyeDigram 327

F

file commands and functions
   See functions, file
flip 329
floating-point numbers 30, 53, 64
for 478
for statement 478
forcenode 112, 113, 114
foreach 480
fourEval 330
freq 332
frequency 336
From a UNIX Shell 45
From the CIW 45
cscanf 489
function body 71
functions
   file
      close 488
      fscanf 489
      gets 491
      inline 492
      load 493
      newline 495
      outfile 496
      pfile 498
      SKILL
abs 256
acos 257
add 1 258
asin 259
atan 260
cos 261
dep 262
int 263
max 267
min 268
mod 269
phaseNoise 168
random 270
resultParam 172
round 271
sin 272
sp 177
sqrt 273
srandom 274
sub1 275
tan 276
vswr 186
xor 277
zm 188
zref 190
waveform
average 279
b1f 284
bandwidth 285
clip 286
compare 288
compression 290
conjugate 298
convolve 299
cross 305
db10 307
db20 308
dbm 309
delay 310
deriv 314
dft 315, 317
dnl 319
evmQAM 323
evmQpsk 325
flip 329
fourEval 330
frequency 332, 336
ga 337
gac 338
gainBwProd 340
gainMargin 342
gmax 343
gmin 344
gmux 346
gpc 348
groupDelay 350
gsmg 345
gt 351
Harmonic 352
harmonicList 356
histo 358
iinteg 359
imag 360
integ 361
ipn 364
kf 373
ln 374
log10 375
lsb 376
lshift 377
mag 378
nc 379
overshoot 381
peak 384
peakToPeak 386, 387
phase 389
phaseDeg 390
phaseDegUnwrapped 391
phaseMargin 392
phaseRad 394
phaseRadUnwrapped 395
pow 396
psd 398
psdBB 402
pzbode 406
pzfilter 407
real 409
riseTime 410
rms 413
rmsNoise 414
root 415
rshift 417
sample 418
settingTime 420
slewRate 423
spectralPower 426, 427
ssb 430, 431
tangent 432
thd 433, 435
value 436
xmax 439
xmin 441
Interactive Session Demonstrating the OCEAN Use Model

if statement 474
im alias 248
integ 359
im alias 248
imag 360
includeFile 117
infile 492
infix arithmetic operators 56
infix operators 59, 61
int 263
integ 361
integer 63

keywords 19
kf 373
killJob 461

let 69
Line Continuation 59
linRg 264
literal characters 19
ln 374
load 493
Loading OCEAN Scripts 45
local variables 69
log 265
log10 375
Logical Operators 55
logical operators 55
logRg 266
lsb 376
lshift 377

i 157
ic 116, 125
if 474
**Mixed-Signal in OCEAN 4.4.6** 579

- `mag` 378
- `max` 267
- `min` 268

**Naming Conventions** 52
- `nc` 379
- `nesting, expressions` 58
- `newline` 495
- `newWindow` 218
- `NF` 245
- `NFmin` 245
- `NNR` 245
- `nodeset` 119
- `noise` 120
- `noiseSummary` 219

**Nonblocking Mode** 35
- `Numbers` 64
- `numbers`
  - `floating-point` 30, 53, 64
  - `integer` 63
- `numbers, scaling factors` 52

**OCEAN**
- `aliases` 247
- `definition` 25
  - `design variables` 31
  - `OCEAN in Non-Graphical Mode` 38
- `OCEAN Online Help` 26
- `OCEAN Return Values` 31
- `OCEAN Syntax Overview` 27
- `OCEAN Tips` 48
- `OCEAN Use Models` 37
- `ocnCloseSession` 121
- `ocnDisplay` 122
- `ocnGetAdjustedPath` 124
- `ocnHelp` 159
- `ocnPrint` 223, 226, 228
- `ocnResetResults` 161
  - `ocnSetAttrib` 226
  - `ocnSetSilentMode` 80
  - `ocnWaveformTool` 125
  - `ocnxlConjugateGradientOptions` 567
  - `ocnxlDisableCorner` 572
  - `ocnxlEnableCorner` 574
  - `ocnxlEnableCornerForTest` 546
  - `ocnxlEnableSweepParam` 547
  - `ocnxlEnableSweepVar` 548
  - `ocnxlEnableTest` 549
  - `ocnxlGetBestPointParams` 550
  - `ocnxlGetcorners` 551
  - `ocnxlGetCurrentHistory` 552
  - `ocnxlGetCurrentHistoryId` 553
  - `ocnxlgetSession` 554
  - `ocnxlGetSpecs` 555
  - `ocnxlGetTests` 556
  - `ocnxlHistoryPrefix` 561
  - `ocnxlLoadSetupState` 562
  - `ocnxlMTSBlock` 569
  - `ocnxlMTSEnable` 568
  - `ocnxlOutputAreaGoal` 566
  - `ocnxlProjectDir` 571
  - `ocnxlRemoveSpec` 557
  - `ocnxlRenameCurrentHistory` 558
  - `ocnxlRun` 559
  - `ocnxlSaveSetupAs` 575
  - `ocnxlSetAllParametersDisabled` 576
  - `ocnxlSetAllVarsDisabled` 577
  - `ocnxlSimResultsLocation` 572
  - `ocnYvsYplot` 228
  - `off` 126
  - `online help` 26
  - `openResults` 34, 162
  - `operators`
    - `arithmetic` 52
    - `binary minus` 57
    - `infix` 56, 61
    - `introduction` 52
    - `logical` 55
    - `relational` 54
    - `unary minus` 57
    - `option` 127
    - `Or-bars in syntax` 19
    - `order of evaluation` 58
    - `outfile` 496, 498
    - `outputParams` 164
  - `outputs` 166
  - `outputs() in OCEAN` 32
  - `overshoot` 381
P

paramAnalysis 448
parameters, definition 70
Parametric Analysis 33
parametric analysis 33
paramRun 452
paramValPair Format 78
Parentheses 27
parentheses 27, 58
Parentheses in C 58
Parentheses in SKILL 58
path 75, 79, 80
peak 384
peakToPeak 386, 387
period_jitter 387
pfile 498
phase 389
phaseDeg 390
phaseDegUnwrapped 391
phaseMargin 392
phaseNoise 168
phaseRad 394
phaseRadUnwrapped 395
plot 230
plotStyle 233
Plotting and Printing SpectreRF Functions in
OCEAN 245
plotting commands. See commands, plotting and printing
pow 396
Predefined Arithmetics 254
prependPath 76
primitives 59
print 499
println 500
procedure 70
procedures, definition 70
See also SKILL functions
psd 398
psdbb 402
pv 170
pzbode 406
pzfilter 407
pzSummary 236

Q

Question Mark 29

question mark 29

R

random 270
real 409
Recovering from an Omitted Double
Quote 28
Related Documents 18
Relational and Logical Operators 54
relational operators 54
removeLabel 238
report 239
restore 129
resultParam 172
results 174
resultsDir 130
resumeJob 464
return value (=>) 70
return values 31
right arrow in syntax 20
riseTime 410
rms 413
rmsNoise 414
RN 245
Role of Parentheses 58
root 415
round 271
rshift 417
run 131
Running Multiple Simulators 48

S

sample 418
save 135
saveOption 137
Scaling Factors 52
scaling factors 52
Selectively Creating Scripts 42
selectResult 175
settingTime 420
settlingTime 420
setup 77
simulation commands
See commands, simulation
simulator 139, 140
sin 272
Single Quotes 29
single quotes 29
SKILL
   commenting code 57
   primitives 59
   white space in code 57
SKILL data types 30
Skill Function Return Values 70
SKILL functions
   arguments 70
   declaring 68
   defining parameters 69
   definition 70
   parameters 70
   terminology 70
Skill Functions 63
SKILL functions, syntax conventions 20
SKILL Syntax 56
SKILL syntax 27
SKILL Syntax Examples 20
slash in syntax 20
slewRate 423
solver 140
sp 177
Special Characters 56
spectralPower 426, 427
spectrum 427
sqrt 273
srandom 274
ssb 430, 431
stddev 431
stimulusFile 141
store 143
strcat function 67
strcmp function 68
Strings 65
strings
   comparing 67
   concatenating 66
   definition 65
strncat function 67
sub1 275
sub1 function 275
suspendJob 465
sweepNames 179
sweepValues 181
sweepVarValues 182
syntax 56
   double quotes 28
   functions 70
   overview 27

parentheses 27
question mark 29
single quotes 29
syntax conventions 19
Syntax Functions for Defining Functions 70

T
tan 276
tan function 276, 277
tangent 432
temp 144
Terms and Definitions 70
thd 433, 435
The Advantages of SKILL 51
tran 145
types of commands 26
Types of OCEAN Commands 26
Typographic and Syntax Conventions 19

U
unary minus operator 57
unbound variables 65
unityGainFreq 435
unless 476
unless statement 476
Using && 55
Using || 56
Using OCEAN from a UNIX Shell 38
Using OCEAN from the CIW 39
Using OCEAN Interactively 38
Using Variables 60

V
v 184
value 436
value function 436
Variables 60
variables
   defining local 69
   definition 60
   introduction 60
   unbound 65
vcdFile 146
vcdInfoFile 147
vdb alias 247
vecFile 148
vertical bars in syntax 19
vim alias 248
vm alias 247
vp alias 247
vr alias 248
vswr 186

W
wait 466
Waveform (Calculator) Functions 278
when 477
when statement 477
while 482
while statement 482
White Space 57
white space 57

X
xLimit 242
xmax 439
xmin 441
xor 277
xval 443

Y
yLimit 243
ymax 444
ymin 445

Z
zm 188
zref 190