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# **Sage Reference Manual: Interpreter Interfaces**

***Release 6.9***

**The Sage Development Team**

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<b>1</b>	<b>Common Interface Functionality</b>	<b>3</b>
<b>2</b>	<b>Common Interface Functionality through Pexpect</b>	<b>9</b>
<b>3</b>	<b>Sage wrapper around pexpect's spawn class</b>	<b>13</b>
<b>4</b>	<b>Interface to Axiom</b>	<b>15</b>
<b>5</b>	<b>The Elliptic Curve Factorization Method</b>	<b>21</b>
<b>6</b>	<b>Interface to 4ti2</b>	<b>27</b>
<b>7</b>	<b>Interface to FriCAS</b>	<b>33</b>
<b>8</b>	<b>Interface to Frobbly for fast computations on monomial ideals.</b>	<b>37</b>
<b>9</b>	<b>Interface to GAP</b>	<b>41</b>
9.1	First Examples . . . . .	41
9.2	GAP and Singular . . . . .	41
9.3	Saving and loading objects . . . . .	42
9.4	Long Input . . . . .	42
9.5	Changing which GAP is used . . . . .	43
<b>10</b>	<b>Interface to GAP3</b>	<b>53</b>
10.1	Obtaining GAP3 . . . . .	53
10.2	Changing which GAP3 is used . . . . .	53
10.3	Functionality and Examples . . . . .	54
10.4	Common Pitfalls . . . . .	55
10.5	Examples . . . . .	55
<b>11</b>	<b>Interface to Groebner Fan</b>	<b>63</b>
<b>12</b>	<b>Interface to Giac</b>	<b>65</b>
12.1	Tutorial . . . . .	65
<b>13</b>	<b>Interface to the Gnuplot interpreter</b>	<b>73</b>
<b>14</b>	<b>Interface to the GP calculator of PARI/GP</b>	<b>75</b>
<b>15</b>	<b>Interface for extracting data and generating images from Jmol readable files.</b>	<b>85</b>
<b>16</b>	<b>Interface to KASH</b>	<b>87</b>
16.1	Issues . . . . .	87

16.2	Tutorial . . . . .	87
16.3	Long Input . . . . .	93
<b>17</b>	<b>Interface to LiE</b>	<b>95</b>
17.1	Tutorial . . . . .	95
<b>18</b>	<b>Lisp Interface</b>	<b>103</b>
<b>19</b>	<b>Interface to Macaulay2</b>	<b>107</b>
<b>20</b>	<b>Interface to Magma</b>	<b>117</b>
20.1	Parameters . . . . .	117
20.2	Multiple Return Values . . . . .	117
20.3	Long Input . . . . .	118
20.4	Garbage Collection . . . . .	118
20.5	Other Examples . . . . .	119
<b>21</b>	<b>Interface to the free online MAGMA calculator</b>	<b>135</b>
<b>22</b>	<b>Interface to Maple</b>	<b>137</b>
22.1	Tutorial . . . . .	137
<b>23</b>	<b>Interface to Mathematica</b>	<b>145</b>
23.1	Tutorial . . . . .	146
23.2	Long Input . . . . .	148
23.3	Loading and saving . . . . .	149
23.4	Complicated translations . . . . .	149
<b>24</b>	<b>Interface to MATLAB</b>	<b>153</b>
24.1	Tutorial . . . . .	153
<b>25</b>	<b>Pexpect interface to Maxima</b>	<b>157</b>
25.1	Tutorial . . . . .	158
25.2	Examples involving matrices . . . . .	160
25.3	Laplace Transforms . . . . .	160
25.4	Continued Fractions . . . . .	161
25.5	Special examples . . . . .	161
25.6	Miscellaneous . . . . .	162
25.7	Interactivity . . . . .	162
25.8	Latex Output . . . . .	162
25.9	Long Input . . . . .	163
<b>26</b>	<b>Abstract interface to Maxima</b>	<b>167</b>
<b>27</b>	<b>Library interface to Maxima</b>	<b>183</b>
<b>28</b>	<b>Interface to MuPAD</b>	<b>197</b>
<b>29</b>	<b>Interface to mwrnk</b>	<b>201</b>
<b>30</b>	<b>Interface to GNU Octave</b>	<b>205</b>
30.1	Computation of Special Functions . . . . .	205
30.2	Tutorial . . . . .	206
<b>31</b>	<b>Interface to PHC.</b>	<b>211</b>
<b>32</b>	<b>POV-Ray, The Persistence of Vision Ray Tracer</b>	<b>217</b>

<b>33</b>	<b>Parallel Interface to the Sage interpreter</b>	<b>219</b>
<b>34</b>	<b>Interface to QEPCAD</b>	<b>221</b>
<b>35</b>	<b>Interface to Bill Hart’s Quadratic Sieve</b>	<b>247</b>
<b>36</b>	<b>Interface to R</b>	<b>251</b>
<b>37</b>	<b>Interface to several Rubik’s cube solvers.</b>	<b>263</b>
<b>38</b>	<b>Interface to Sage</b>	<b>265</b>
<b>39</b>	<b>Interface to Scilab</b>	<b>269</b>
<b>40</b>	<b>Interface to Singular</b>	<b>275</b>
40.1	Introduction . . . . .	275
40.2	Tutorial . . . . .	275
40.3	Computing the Genus . . . . .	278
40.4	An Important Concept . . . . .	278
40.5	Long Input . . . . .	279
<b>41</b>	<b>The Tachyon Ray Tracer</b>	<b>295</b>
<b>42</b>	<b>Interface to TIDES</b>	<b>297</b>
<b>43</b>	<b>Interface to the Sage cleaner</b>	<b>303</b>
<b>44</b>	<b>Quitting interfaces</b>	<b>305</b>
<b>45</b>	<b>An interface to read data files</b>	<b>307</b>
<b>46</b>	<b>Indices and Tables</b>	<b>309</b>
	<b>Bibliography</b>	<b>311</b>



Sage provides a unified interface to the best computational software. This is accomplished using both C-libraries (see C/C++ Library Interfaces) and interpreter interfaces, which are implemented using pseudo-tty's, system files, etc. This chapter is about these interpreter interfaces.

---

**Note:** Each interface requires that the corresponding software is installed on your computer. Sage includes GAP, PARI, Singular, and Maxima, but does not include Octave (very easy to install), MAGMA (non-free), Maple (non-free), or Mathematica (non-free).

There is overhead associated with each call to one of these systems. For example, computing  $2+2$  thousands of times using the GAP interface will be slower than doing it directly in Sage. In contrast, the C-library interfaces of C/C++ Library Interfaces incur less overhead.

---

In addition to the commands described for each of the interfaces below, you can also type e.g., `%gap`, `%magma`, etc., to directly interact with a given interface in its state. Alternatively, if `X` is an interface object, typing `X.interact()` allows you to interact with it. This is completely different than `X.console()` which starts a complete new copy of whatever program `X` interacts with. Note that the input for `X.interact()` is handled by Sage, so the history buffer is the same as for Sage, tab completion is as for Sage (unfortunately!), and input that spans multiple lines must be indicated using a backslash at the end of each line. You can pull data into an interactive session with `X` using `sage(expression)`.

The console and interact methods of an interface do very different things. For example, using gap as an example:

1. `gap.console()`: You are completely using another program, e.g., `gap/magma/gp`. Here Sage is serving as nothing more than a convenient program launcher, similar to `bash`.
2. `gap.interact()`: This is a convenient way to interact with a running gap instance that may be “full of” Sage objects. You can import Sage objects into this gap (even from the interactive interface), etc.

The console function is very useful on occasion, since you get the exact actual program available (especially useful for tab completion and testing to make sure nothing funny is going on).





## COMMON INTERFACE FUNCTIONALITY

See the examples in the other sections for how to use specific interfaces. The interface classes all derive from the generic interface that is described in this section.

AUTHORS:

- William Stein (2005): initial version
- William Stein (2006-03-01): got rid of infinite loop on startup if client system missing
- Felix Lawrence (2009-08-21): edited `._sage_()` to support lists and float exponents in foreign notation.
- Simon King (2010-09-25): `Expect._local_tmpfile()` depends on `Expect.pid()` and is cached; `Expect.quit()` clears that cache, which is important for forking.
- Jean-Pierre Flori (2010,2011): Split non Pexpect stuff into a parent class.
- Simon King (2015): Improve pickling for `InterfaceElement`

```
class sage.interfaces.interface.AsciiArtString
    Bases: str
```

```
class sage.interfaces.interface.Interface(name)
    Bases: sage.structure.parent_base.ParentWithBase
```

Interface interface object.

**call** (*function\_name*, \**args*, \*\**kwds*)

**clear** (*var*)

Clear the variable named *var*.

**console** ()

**cputime** ()

CPU time since this process started running.

**eval** (*code*, \*\**kwds*)

**execute** (\**args*, \*\**kwds*)

**function\_call** (*function*, *args*=None, *kwds*=None)

EXAMPLES:

```
sage: maxima.quad_qags(x, x, 0, 1, epsrel=1e-4)
```

```
[0.5, 5.5511151231257...e-15, 21, 0]
```

```
sage: maxima.function_call('quad_qags', [x, x, 0, 1], {'epsrel': '1e-4'})
```

```
[0.5, 5.5511151231257...e-15, 21, 0]
```

**get** (*var*)

Get the value of the variable *var*.

**get\_seed()**

Returns the seed used to set the random number generator in this interface. The seed is initialized as None but should be set when the interface starts.

## EXAMPLES:

```
sage: s = Singular()
sage: s.set_seed(107)
107
sage: s.get_seed()
107
```

**get\_using\_file(var)**

Return the string representation of the variable var in self, possibly using a file. Use this if var has a huge string representation, since it may be way faster.

**Warning:** In fact unless a special derived class implements this, it will *not* be any faster. This is the case for this class if you're reading it through introspection and seeing this.

**help(s)****interact()**

This allows you to interactively interact with the child interpreter. Press Ctrl-D or type 'quit' or 'exit' to exit and return to Sage.

---

**Note:** This is completely different than the console() member function. The console function opens a new copy of the child interpreter, whereas the interact function gives you interactive access to the interpreter that is being used by Sage. Use sage(xxx) or interpretername(xxx) to pull objects in from sage to the interpreter.

---

**name(new\_name=None)****new(code)****rand\_seed()**

Returns a random seed that can be put into set\_seed function for any interpreter. This should be overridden if the particular interface needs something other than a small positive integer.

## EXAMPLES:

```
from sage.misc.random_testing import random_testing
sage: from sage.interfaces.interface import Interface
sage: i = Interface("")
sage: i.rand_seed() # random
318491487L

sage: s = Singular()
sage: s.rand_seed() # random
365260051L
```

**read(filename)**

## EXAMPLES:

```
sage: filename = tmp_filename()
sage: f = open(filename, 'w')
sage: f.write('x = 2\n')
sage: f.close()
sage: octave.read(filename) # optional - octave
sage: octave.get('x')       # optional - octave
' 2'
```

```
sage: import os
sage: os.unlink(filename)
```

**set** (*var*, *value*)

Set the variable *var* to the given value.

**set\_seed** (*seed=**None*)

Sets the random seed for the interpreter and returns the new value of the seed. This is dependent on which interpreter so must be implemented in each separately. For examples see `gap.py` or `singular.py`. If *seed* is *None* then should generate a random seed.

EXAMPLES:

```
sage: s = Singular()
sage: s.set_seed(1)
1
sage: [s.random(1,10) for i in range(5)]
[8, 10, 4, 9, 1]

sage: from sage.interfaces.interface import Interface
sage: i = Interface("")
sage: i.set_seed()
Traceback (most recent call last):
...
NotImplementedError: This interpreter did not implement a set_seed function
```

```
class sage.interfaces.interface.InterfaceElement (parent, value, is_name=False,
                                                    name=None)
```

Bases: `sage.structure.element.RingElement`

Interface element.

**attribute** (*attrname*)

If this wraps the object *x* in the system, this returns the object *x.attrname*. This is useful for some systems that have object oriented attribute access notation.

EXAMPLES:

```
sage: g = gap('SO(1,4,7)')
sage: k = g.InvariantQuadraticForm()
sage: k.attribute('matrix')
[[ [ 0*Z(7), Z(7)^0, 0*Z(7), 0*Z(7) ], [ 0*Z(7), 0*Z(7), 0*Z(7), 0*Z(7) ],
  [ 0*Z(7), 0*Z(7), Z(7), 0*Z(7) ], [ 0*Z(7), 0*Z(7), 0*Z(7), Z(7)^0 ] ]

sage: e = gp('ellinit([0,-1,1,-10,-20])')
sage: e.attribute('j')
-122023936/161051
```

**bool** ()

**gen** (*n*)

**get\_using\_file** ()

Return this element's string representation using a file. Use this if self has a huge string representation. It'll be way faster.

EXAMPLES:

```
sage: a = maxima(str(2^1000))
sage: a.get_using_file()
'1071508607186267320948425049060001810561404811705533607443750388370351051124936122493198378'
```

**hasattr** (*attrname*)

Returns whether the given attribute is already defined by this object, and in particular is not dynamically generated.

EXAMPLES:

```
sage: m = maxima('2')
sage: m.hasattr('integral')
True
sage: m.hasattr('gcd')
False
```

**is\_string** ()

Tell whether this element is a string.

By default, the answer is negative.

**name** (*new\_name=None*)

Returns the name of self. If *new\_name* is passed in, then this function returns a new object identical to self whose name is *new\_name*.

Note that this can overwrite existing variables in the system.

EXAMPLES:

```
sage: x = r([1,2,3]); x
[1] 1 2 3
sage: x.name()
'sage3'
sage: x = r([1,2,3]).name('x'); x
[1] 1 2 3
sage: x.name()
'x'

sage: s5 = gap.SymmetricGroup(5).name('s5')
sage: s5
SymmetricGroup( [ 1 .. 5 ] )
sage: s5.name()
's5'
```

**sage** (*\*args, \*\*kws*)

Attempt to return a Sage version of this object.

This method does nothing more than calling `_sage_()`, simply forwarding any additional arguments.

EXAMPLES:

```
sage: gp(1/2).sage()
1/2
sage: _.parent()
Rational Field
sage: singular.lib("matrix")
sage: R = singular.ring(0, '(x,y,z)', 'dp')
sage: singular.matrix(2,2).sage()
[0 0]
[0 0]
```

**class** sage.interfaces.interface.**InterfaceFunction** (*parent, name*)

Bases: sage.structure.sage\_object.SageObject

Interface function.

```
class sage.interfaces.interface.InterfaceFunctionElement (obj, name)  
    Bases: sage.structure.sage_object.SageObject  
    Interface function element.  
    help()  
sage.interfaces.interface.is_InterfaceElement (x)
```



## COMMON INTERFACE FUNCTIONALITY THROUGH PEXPECT

See the examples in the other sections for how to use specific interfaces. The interface classes all derive from the generic interface that is described in this section.

AUTHORS:

- William Stein (2005): initial version
- William Stein (2006-03-01): got rid of infinite loop on startup if client system missing
- Felix Lawrence (2009-08-21): edited `._sage_()` to support lists and float exponents in foreign notation.
- Simon King (2010-09-25): `Expect._local_tmpfile()` depends on `Expect.pid()` and is cached; `Expect.quit()` clears that cache, which is important for forking.
- Jean-Pierre Flori (2010,2011): Split non Pexpect stuff into a parent class.
- Simon King (2010-11-23): Ensure that the interface is started again after a crash, when a command is executed in `_eval_line`. Allow synchronisation of the GAP interface.

```
class sage.interfaces.expect.Expect(name, prompt, command=None, server=None,
                                     server_tmpdir=None, ulimit=None, maxread=100000,
                                     script_subdirectory=None, restart_on_ctrlc=False, ver-
                                     bose_start=False, init_code=[], max_startup_time=None,
                                     logfile=None, eval_using_file_cutoff=0, do_cleaner=True,
                                     remote_cleaner=False, path=None, terminal_echo=True)
```

Bases: `sage.interfaces.interface.Interface`

Expect interface object.

**clear\_prompts()**

**detach()**

Forget the running subprocess: keep it running but pretend that it's no longer running.

EXAMPLES:

```
sage: a = maxima('y')
sage: saved_expect = maxima._expect # Save this to close later
sage: maxima.detach()
sage: a._check_valid()
Traceback (most recent call last):
...
ValueError: The maxima session in which this object was defined is no longer running.
sage: saved_expect.close() # Close child process
```

Calling `detach()` a second time does nothing:

```
sage: maxima.detach()
```

**eval** (*code*, *strip*=*True*, *synchronize*=*False*, *locals*=*None*, *allow\_use\_file*=*True*, *split\_lines*=*'nofile'*,  
\*\**kws*)

INPUT:

- **code** – text to evaluate
- **strip** – **bool**; whether to strip output prompts, etc. (ignored in the base class).
- **locals** – **None** (ignored); this is used for compatibility with the Sage notebook’s generic system interface.
- **allow\_use\_file** – **bool** (default: **True**); if **True** and **code** exceeds an interface-specific threshold then **code** will be communicated via a temporary file rather than the character-based interface. If **False** then the code will be communicated via the character interface.
- **split\_lines** – **Tri-state** (default: **“nofile”**); if **“nofile”** then **code** is sent line by line unless it gets communicated via a temporary file. If **True** then **code** is sent line by line, but some lines individually might be sent via temporary file. Depending on the interface, this may transform grammatical **code** into ungrammatical input. If **False**, then the whole block of code is evaluated all at once.
- **\*\*kws** – All other arguments are passed onto the **\_eval\_line** method. An often useful example is **reformat=False**.

**expect** ()

**interrupt** (*tries*=5, *timeout*=2.0, *quit\_on\_fail*=*True*)

**is\_local** ()

**is\_remote** ()

**is\_running** ()

Return True if self is currently running.

**path** ()

**pid** ()

Return the PID of the underlying sub-process.

REMARK:

If the interface terminates unexpectedly, the original PID will still be used. But if it was terminated using **quit** (), a new sub-process with a new PID is automatically started.

EXAMPLE:

```
sage: pid = gap.pid()
sage: gap.eval('quit;')
''
sage: pid == gap.pid()
True
sage: gap.quit()
sage: pid == gap.pid()
False
```

**quit** (*verbose*=*False*, *timeout*=*None*)

Quit the running subprocess.

INPUT:

- **verbose** – (boolean, default **False**) print a message when quitting this process?

EXAMPLES:



```

sage: a = maxima('y')
sage: maxima.quit(verbose=True)
Exiting Maxima with PID ... running .../local/bin/maxima ...
sage: a._check_valid()
Traceback (most recent call last):
...
ValueError: The maxima session in which this object was defined is no longer running.

```

Calling `quit()` a second time does nothing:

```
sage: maxima.quit(verbose=True)
```

**user\_dir()**

```
class sage.interfaces.expect.ExpectElement (parent, value, is_name=False, name=None)
Bases: sage.interfaces.interface.InterfaceElement
```

Expect element.

```
class sage.interfaces.expect.ExpectFunction (parent, name)
Bases: sage.interfaces.interface.InterfaceFunction
```

Expect function.

```
class sage.interfaces.expect.FunctionElement (obj, name)
Bases: sage.interfaces.interface.InterfaceFunctionElement
```

Expect function element.

```
class sage.interfaces.expect.StdoutContext (interface, silent=False, stdout=None)
A context in which all communication between Sage and a subprocess interfaced via pexpect is printed to stdout.
```

```
sage.interfaces.expect.console (cmd)
```

```
class sage.interfaces.expect.gc_disabled
Bases: object
```

This is a “with” statement context manager. Garbage collection is disabled within its scope. Nested usage is properly handled.

EXAMPLES:

```

sage: import gc
sage: from sage.interfaces.expect import gc_disabled
sage: gc.isenabled()
True
sage: with gc_disabled():
.....:     print gc.isenabled()
.....:     with gc_disabled():
.....:         print gc.isenabled()
.....:     print gc.isenabled()
False
False
False
sage: gc.isenabled()
True

```

```
sage.interfaces.expect.is_ExpectElement (x)
```



## SAGE WRAPPER AROUND PEXPECT'S SPAWN CLASS

AUTHOR:

- Jeroen Demeyer (2015-02-01): initial version, [trac ticket #17686](#)

```
class sage.interfaces.sagespawn.SageSpawn(*args, **kwds)
    Bases: pexpect.spawn
```

Spawn a subprocess in a pseudo-tty.

- `*args, **kwds`: see `pexpect.spawn`.
- `name` – human-readable name for this process, used for display purposes only.
- `quit_string` – (default: `None`) if not `None`, send this string to the child process before killing it.

EXAMPLES:

```
sage: from sage.interfaces.sagespawn import SageSpawn
sage: SageSpawn("sleep 1", name="Sleeping Beauty")
Sleeping Beauty with PID ... running ...
```

```
close()
```

Quit the child process: send the quit string, close the pseudo-tty and kill the process.

This function returns immediately, it doesn't wait for the child process to die.

EXAMPLES:

```
sage: from sage.interfaces.sagespawn import SageSpawn
sage: s = SageSpawn("sleep 1000")
sage: s.close()
sage: while s.isalive(): # long time (5 seconds)
....:     sleep(0.1)
```

```
expect_peek(*args, **kwds)
```

Like `expect()` but restore the read buffer such that it looks like nothing was actually read. The next reading will continue at the current position.

EXAMPLES:

```
sage: from sage.interfaces.sagespawn import SageSpawn
sage: E = SageSpawn("sh", ["-c", "echo hello world"])
sage: _ = E.expect_peek("w")
sage: E.read()
'hello world\r\n'
```

```
expect_upto(*args, **kwds)
```

Like `expect()` but restore the read buffer starting from the matched string. The next reading will continue starting with the matched string.

EXAMPLES:

```
sage: from sage.interfaces.sagespawn import SageSpawn
sage: E = SageSpawn("sh", ["-c", "echo hello world"])
sage: _ = E.expect_upto("w")
sage: E.read()
'world\r\n'
```

**terminate\_async** (*interval=5.0*)

Terminate the child process group asynchronously.

This function returns immediately, while the child is slowly being killed in the background.

INPUT:

- *interval* – (default: 5) how much seconds to wait between sending two signals.

EXAMPLES:

Run an infinite loop in the shell:

```
sage: from sage.interfaces.sagespawn import SageSpawn
sage: s = SageSpawn("sh", ["-c", "while true; do sleep 1; done"])
```

Check that the process eventually dies after calling `terminate_async`:

```
sage: s.terminate_async(interval=0.2)
sage: while True:
....:     try:
....:         os.kill(s.pid, 0)
....:     except OSError:
....:         sleep(0.1)
....:     else:
....:         break # process got killed
```

## INTERFACE TO AXIOM

TODO:

- Evaluation using a file is not done. Any input line with more than a few thousand characters would hang the system, so currently it automatically raises an exception.
- All completions of a given command.
- Interactive help.

Axiom is a free GPL-compatible (modified BSD license) general purpose computer algebra system whose development started in 1973 at IBM. It contains symbolic manipulation algorithms, as well as implementations of special functions, including elliptic functions and generalized hypergeometric functions. Moreover, Axiom has implementations of many functions relating to the invariant theory of the symmetric group  $S_n$ . For many links to Axiom documentation see <http://wiki.axiom-developer.org>.

AUTHORS:

- Bill Page (2006-10): Created this (based on Maxima interface)

---

**Note:** Bill Page put a huge amount of effort into the Sage Axiom interface over several days during the Sage Days 2 coding sprint. This contribution is greatly appreciated.

---

- William Stein (2006-10): misc touchup.
- Bill Page (2007-08): Minor modifications to support axiom4sage-0.3

---

**Note:** The axiom4sage-0.3.spkg is based on an experimental version of the FriCAS fork of the Axiom project by Waldek Hebisch that uses pre-compiled cached Lisp code to build Axiom very quickly with clisp.

---

If the string “error” (case insensitive) occurs in the output of anything from axiom, a `RuntimeError` exception is raised.

EXAMPLES: We evaluate a very simple expression in axiom.

```
sage: axiom('3 * 5')                                #optional - axiom
15
sage: a = axiom(3) * axiom(5); a                      #optional - axiom
15
```

The type of `a` is `AxiomElement`, i.e., an element of the axiom interpreter.

```
sage: type(a)                                         #optional - axiom
<class 'sage.interfaces.axiom.AxiomElement'>
sage: parent(a)                                       #optional - axiom
Axiom
```

The underlying Axiom type of `a` is also available, via the `type` method:

```
sage: a.type()                                     #optional - axiom
PositiveInteger
```

We factor  $x^5 - y^5$  in Axiom in several different ways. The first way yields a Axiom object.

```
sage: F = axiom.factor('x^5 - y^5'); F             #optional - axiom
      4      3      2 2      3      4
- (y - x)(y + x y + x y + x y + x )
sage: type(F)                                     #optional - axiom
<class 'sage.interfaces.axiom.AxiomElement'>
sage: F.type()                                   #optional - axiom
Factored Polynomial Integer
```

Note that Axiom objects are normally displayed using “ASCII art”.

```
sage: a = axiom(2/3); a                           #optional - axiom
  2
  -
  3
sage: a = axiom('x^2 + 3/7'); a                   #optional - axiom
  2      3
x  + -
      7
```

The `axiom.eval` command evaluates an expression in axiom and returns the result as a string. This is exact as if we typed in the given line of code to axiom; the return value is what Axiom would print out.

```
sage: print axiom.eval('factor(x^5 - y^5)')        #optional - axiom
      4      3      2 2      3      4
- (y - x)(y + x y + x y + x y + x )
Type: Factored Polynomial Integer
```

We can create the polynomial  $f$  as a Axiom polynomial, then call the factor method on it. Notice that the notation `f.factor()` is consistent with how the rest of Sage works.

```
sage: f = axiom('x^5 - y^5')                      #optional - axiom
sage: f^2                                           #optional - axiom
      10      5 5      10
y  - 2x y + x
sage: f.factor()                                   #optional - axiom
      4      3      2 2      3      4
- (y - x)(y + x y + x y + x y + x )
```

Control-C interruption works well with the axiom interface, because of the excellent implementation of axiom. For example, try the following sum but with a much bigger range, and hit control-C.

```
sage: f = axiom('(x^5 - y^5)^10000')              # not tested
Interrupting Axiom...
...
<type 'exceptions.TypeError': Ctrl-c pressed while running Axiom>

sage: axiom('1/100 + 1/101')                      #optional - axiom
      201
-----
      10100
sage: a = axiom('(1 + sqrt(2))^5'); a              #optional - axiom
      +-+
      29\|2  + 41
```

TESTS: We check to make sure the subst method works with keyword arguments.

```
sage: a = axiom(x+2); a #optional - axiom
x + 2
sage: a.subst(x=3) #optional - axiom
5
```

We verify that Axiom floating point numbers can be converted to Python floats.

```
sage: float(axiom(2)) #optional - axiom
2.0
```

```
class sage.interfaces.axiom.Axiom(name='axiom', command='axiom -nox -noclef',
                                   script_subdirectory=None, logfile=None, server=None,
                                   server_tmpdir=None, init_code=['lisp (si::readline-off)'])
    Bases: sage.interfaces.axiom.PanAxiom
```

Create an instance of the Axiom interpreter.

TESTS:

```
sage: axiom == loads(dumps(axiom))
True
```

**console()**

Spawn a new Axiom command-line session.

EXAMPLES:

```
sage: axiom.console() #not tested
AXIOM Computer Algebra System
Version: Axiom (January 2009)
Timestamp: Sunday January 25, 2009 at 07:08:54
```

```
-----
Issue )copyright to view copyright notices.
Issue )summary for a summary of useful system commands.
Issue )quit to leave AXIOM and return to shell.
-----
```

```
class sage.interfaces.axiom.AxiomElement(parent, value, is_name=False, name=None)
    Bases: sage.interfaces.axiom.PanAxiomElement
```

```
class sage.interfaces.axiom.AxiomExpectFunction(parent, name)
    Bases: sage.interfaces.axiom.PanAxiomExpectFunction
```

TESTS:

```
sage: axiom.upperCase_q
upperCase?
sage: axiom.upperCase_e
upperCase!
```

```
class sage.interfaces.axiom.AxiomFunctionElement(object, name)
    Bases: sage.interfaces.axiom.PanAxiomFunctionElement
```

TESTS:

```
sage: a = axiom('Hello') #optional - axiom
sage: a.upperCase_q #optional - axiom
upperCase?
```

```
sage: a.upperCase_e      #optional - axiom
upperCase!
sage: a.upperCase_e()    #optional - axiom
"HELLO"
```

```
class sage.interfaces.axiom.PanAxiom(name='axiom', command='axiom -nox -noclef',
                                     script_subdirectory=None, logfile=None, server=None,
                                     server_tmpdir=None, init_code=['lisp (si::readline-
                                     off)'])
```

Bases: `sage.interfaces.expect.Expect`

Interface to a PanAxiom interpreter.

**get** (*var*)

Get the string value of the Axiom variable *var*.

EXAMPLES:

```
sage: axiom.set('xx', '2')      #optional - axiom
sage: axiom.get('xx')           #optional - axiom
'2'
sage: a = axiom('(1 + sqrt(2))^5') #optional - axiom
sage: axiom.get(a.name())       #optional - axiom
'      +-+\r\r\n  29\\|2  + 41'
```

**set** (*var*, *value*)

Set the variable *var* to the given value.

EXAMPLES:

```
sage: axiom.set('xx', '2')      #optional - axiom
sage: axiom.get('xx')           #optional - axiom
'2'

sage: fricas.set('xx', '2')     #optional - fricas
sage: fricas.get('xx')          #optional - fricas
'2'
```

**trait\_names** (*verbose=True*, *use\_disk\_cache=True*)

Returns a list of all the commands defined in Axiom and optionally (per default) store them to disk.

EXAMPLES:

```
sage: c = axiom.trait_names(use_disk_cache=False, verbose=False) #optional - axiom
sage: len(c) > 100 #optional - axiom
True
sage: 'factor' in c #optional - axiom
True
sage: '**' in c     #optional - axiom
False
sage: 'upperCase?' in c #optional - axiom
False
sage: 'upperCase_q' in c #optional - axiom
True
sage: 'upperCase_e' in c #optional - axiom
True
```

```
class sage.interfaces.axiom.PanAxiomElement(parent, value, is_name=False, name=None)
```

Bases: `sage.interfaces.expect.ExpectElement`



**as\_type**(*type*)

Returns self as type.

## EXAMPLES:

```

sage: a = axiom(1.2); a          #optional - axiom
1.2
sage: a.as_type(axiom.DoubleFloat) #optional - axiom
1.2
sage: _.type()                  #optional - axiom
DoubleFloat

sage: a = fricas(1.2); a        #optional - fricas
1.2
sage: a.as_type(fricas.DoubleFloat) #optional - fricas
1.2
sage: _.type()                  #optional - fricas
DoubleFloat

```

**comma**(*\*args*)

Returns a Axiom tuple from self and args.

## EXAMPLES:

```

sage: two = axiom(2)    #optional - axiom
sage: two.comma(3)      #optional - axiom
[2, 3]
sage: two.comma(3, 4)   #optional - axiom
[2, 3, 4]
sage: _.type()          #optional - axiom
Tuple PositiveInteger

sage: two = fricas(2)   #optional - fricas
sage: two.comma(3)      #optional - fricas
[2, 3]
sage: two.comma(3, 4)   #optional - fricas
[2, 3, 4]
sage: _.type()          #optional - fricas
Tuple(PositiveInteger)

```

**type**()

Returns the type of an AxiomElement.

## EXAMPLES:

```

sage: axiom(x+2).type() #optional - axiom
Polynomial Integer

```

**unparsed\_input\_form**()

Get the linear string representation of this object, if possible (often it isn't).

## EXAMPLES:

```

sage: a = axiom(x^2+1); a          #optional - axiom
      2
      x  + 1
sage: a.unparsed_input_form() #optional - axiom
'x*x+1'

sage: a = fricas(x^2+1)            #optional - fricas
sage: a.unparsed_input_form() #optional - fricas
'x^2+1'

```

```
class sage.interfaces.axiom.PanAxiomExpectFunction (parent, name)
    Bases: sage.interfaces.expect.ExpectFunction
```

TESTS:

```
sage: axiom.upperCase_q
upperCase?
sage: axiom.upperCase_e
upperCase!
```

```
class sage.interfaces.axiom.PanAxiomFunctionElement (object, name)
    Bases: sage.interfaces.expect.FunctionElement
```

TESTS:

```
sage: a = axiom(' "Hello"') #optional - axiom
sage: a.upperCase_q         #optional - axiom
upperCase?
sage: a.upperCase_e         #optional - axiom
upperCase!
sage: a.upperCase_e()       #optional - axiom
"HELLO"
```

```
sage.interfaces.axiom.axiom_console()
```

Spawn a new Axiom command-line session.

EXAMPLES:

```
sage: axiom_console() #not tested
      AXIOM Computer Algebra System
      Version: Axiom (January 2009)
      Timestamp: Sunday January 25, 2009 at 07:08:54
-----
      Issue )copyright to view copyright notices.
      Issue )summary for a summary of useful system commands.
      Issue )quit to leave AXIOM and return to shell.
-----
```

```
sage.interfaces.axiom.is_AxiomElement (x)
```

Returns True if x is of type AxiomElement.

EXAMPLES:

```
sage: from sage.interfaces.axiom import is_AxiomElement
sage: is_AxiomElement(axiom(2)) #optional - axiom
True
sage: is_AxiomElement(2)
False
```

```
sage.interfaces.axiom.reduce_load_Axiom()
```

Returns the Axiom interface object defined in sage.interfaces.axiom.

EXAMPLES:

```
sage: from sage.interfaces.axiom import reduce_load_Axiom
sage: reduce_load_Axiom()
Axiom
```

## THE ELLIPTIC CURVE FACTORIZATION METHOD

The elliptic curve factorization method (ECM) is the fastest way to factor a **known composite** integer if one of the factors is relatively small (up to approximately 80 bits / 25 decimal digits). To factor an arbitrary integer it must be combined with a primality test. The `ECM.factor()` method is an example for how to combine ECM with a primality test to compute the prime factorization of integers.

Sage includes GMP-ECM, which is a highly optimized implementation of Lenstra's elliptic curve factorization method. See <http://ecm.gforge.inria.fr> for more about GMP-ECM.

### AUTHORS:

These people wrote GMP-ECM: Pierrick Gaudry, Jim Fougeron, Laurent Fousse, Alexander Kruppa, Dave Newman, Paul Zimmermann

### BUGS:

Output from ecm is non-deterministic. Doctests should set the random seed, but currently there is no facility to do so.

```
class sage.interfaces.ecm.ECM(B1=10, B2=None, **kws)
```

```
    Bases: sage.structure.sage_object.SageObject
```

Create an interface to the GMP-ECM elliptic curve method factorization program.

See <http://ecm.gforge.inria.fr>

### INPUT:

- `B1` – integer. Stage 1 bound
- `B2` – integer. Stage 2 bound (or interval `B2min-B2max`)

In addition the following keyword arguments can be used:

- `x0` – integer  $x$ . use  $x$  as initial point
- `sigma` – integer  $s$ . Use  $s$  as curve generator [ecm]
- `A` – integer  $a$ . Use  $a$  as curve parameter [ecm]
- `k` – integer  $n$ . Perform  $\geq n$  steps in stage 2
- `power` – integer  $n$ . Use  $x^n$  for Brent-Suyama's extension
- `dickson` – integer  $n$ . Use  $n$ -th Dickson's polynomial for Brent-Suyama's extension
- `c` – integer  $n$ . Perform  $n$  runs for each input
- `pm1` – boolean. perform P-1 instead of ECM
- `pp1` – boolean. perform P+1 instead of ECM
- `q` – boolean. quiet mode

- `v` – boolean. verbose mode
- `timestamp` – boolean. print a time stamp with each number
- `mpzmod` – boolean. use GMP's `mpz_mod` for mod reduction
- `modmuln` – boolean. use Montgomery's `MODMULN` for mod reduction
- `redc` – boolean. use Montgomery's `REDC` for mod reduction
- `nobase2` – boolean. Disable special base-2 code
- `base2` – integer  $n$ . Force base 2 mode with  $2^{n+1}$  ( $n > 0$ ) or  $2^{n-1}$  ( $n < 0$ )
- `save` – string filename. Save residues at end of stage 1 to file
- `savea` – string filename. Like `-save`, appends to existing files
- `resume` – string filename. Resume residues from file, reads from stdin if file is “-“
- `primetest` – boolean. Perform a primality test on input
- `treefile` – string. Store product tree of  $F$  in files `f.0 f.1 ...`
- `i` – integer. increment  $B1$  by this constant on each run
- `I` – integer  $f$ . auto-calculated increment for  $B1$  multiplied by  $f$  scale factor.
- `inp` – string. Use file as input (instead of redirecting stdin)
- `b` – boolean. Use breadth-first mode of file processing
- `d` – boolean. Use depth-first mode of file processing (default)
- `one` – boolean. Stop processing a candidate if a factor is found (looping mode )
- `n` – boolean. Run ecm in ‘nice’ mode (below normal priority)
- `nn` – boolean. Run ecm in ‘very nice’ mode (idle priority)
- `t` – integer  $n$ . Trial divide candidates before  $P-1$ ,  $P+1$  or ECM up to  $n$ .
- `ve` – integer  $n$ . Verbosely show short ( $< n$  character) expressions on each loop
- `cofdec` – boolean. Force cofactor output in decimal (even if expressions are used )
- `B2scale` – integer. Multiplies the default  $B2$  value
- `go` – integer. Preload with group order `val`, which can be a simple expression, or can use  $N$  as a placeholder for the number being factored.
- `prp` – string. use shell command `cmd` to do large primality tests
- `prplen` – integer. only candidates longer than this number of digits are ‘large’
- `prpval` – integer. `value`  $\geq 0$  which indicates the `prp` command found number to be PRP.
- `prptmp` – file. outputs  $n$  value to temp file prior to running (NB. gets deleted)
- `prplog` – file. otherwise get PRP results from this file (NB. gets deleted)
- `prpyes` – string. Literal string found in `prplog` file when number is PRP
- `prpno` – string. Literal string found in `prplog` file when number is composite

**factor** ( $n$ , *factor\_digits=None*, *B1=2000*, *proof=False*, *\*\*kws*)

Return a probable prime factorization of  $n$ .

Combines GMP-ECM with a primality test, see `is_prime()`. The primality test is provable or probabilistic depending on the *proof* flag.

Moreover, for small  $n$  PARI is used directly.

**Warning:** There is no mathematical guarantee that the factors returned are actually prime if `proof=False` (default). It is extremely likely, though. Currently, there are no known examples where this fails.

INPUT:

- `n` – a positive integer
- `factor_digits` – integer or `None` (default). Optional guess at how many digits are in the smallest factor.
- `B1` – initial lower bound, defaults to 2000 (15 digit factors). Used if `factor_digits` is not specified.
- `proof` – boolean (default: `False`). Whether to prove that the factors are prime.
- `kws` – keyword arguments to pass to `ecm-gmp`. See help for [ECM](#) for more details.

OUTPUT:

A list of integers whose product is  $n$ .

---

**Note:** Trial division should typically be performed, but this is not implemented (yet) in this method.

If you suspect that  $n$  is the product of two similarly-sized primes, other methods (such as a quadratic sieve – use the `qsieve` command) will usually be faster.

The best known algorithm for factoring in the case where all factors are large is the general number field sieve. This is not implemented in Sage; You probably want to use a cluster for problems of this size.

---

EXAMPLES:

```
sage: ecm.factor(602400691612422154516282778947806249229526581)
[45949729863572179, 13109994191499930367061460439]
sage: ecm.factor((2^197 + 1)/3) # long time
[197002597249, 1348959352853811313, 251951573867253012259144010843]
sage: ecm.factor(179427217^13) == [179427217] * 13
True
```

**find\_factor** ( $n$ , `factor_digits=None`, `B1=2000`, `**kws`)

Return a factor of  $n$ .

See also `factor()` if you want a prime factorization of  $n$ .

INPUT:

- `n` – a positive integer,
- `factor_digits` – integer or `None` (default). Decimal digits estimate of the wanted factor.
- `B1` – integer. Stage 1 bound (default 2000). This is used as bound if `factor_digits` is not specified.
- `kws` – optional keyword parameters.

OUTPUT:

List of integers whose product is  $n$ . For certain lengths of the factor, this is the best algorithm to find a factor.

EXAMPLES:

```
sage: f = ECM()
sage: n = 508021860739623467191080372196682785441177798407961
sage: f.find_factor(n)
[79792266297612017, 6366805760909027985741435139224233]
```

Note that the input number can't have more than 4095 digits:

```
sage: f=2^2^14+1
sage: ecm.find_factor(f)
Traceback (most recent call last):
...
ValueError: n must have at most 4095 digits
```

### `get_last_params()`

Return the parameters (including the curve) of the last ecm run.

In the case that the number was factored successfully, this will return the parameters that yielded the factorization.

OUTPUT:

A dictionary containing the parameters for the most recent factorization.

EXAMPLES:

```
sage: ecm.factor((2^197 + 1)/3) # long time
[197002597249, 1348959352853811313, 251951573867253012259144010843]
sage: ecm.get_last_params() # random output
{'poly': 'x^1', 'sigma': '1785694449', 'B1': '8885', 'B2': '1002846'}
```

### `interact()`

Interactively interact with the ECM program.

EXAMPLES:

```
sage: ecm.interact() # not tested
```

### `one_curve(n, factor_digits=None, B1=2000, algorithm='ECM', **kws)`

Run one single ECM (or P-1/P+1) curve on input  $n$ .

Note that trying a single curve is not particularly useful by itself. One typically needs to run over thousands of trial curves to factor  $n$ .

INPUT:

- $n$  – a positive integer
- `factor_digits` – integer. Decimal digits estimate of the wanted factor.
- `B1` – integer. Stage 1 bound (default 2000)
- `algorithm` – either “ECM” (default), “P-1” or “P+1”

OUTPUT:

a list  $[p, q]$  where  $p$  and  $q$  are integers and  $n = p * q$ . If no factor was found, then  $p = 1$  and  $q = n$ .

**Warning:** Neither  $p$  nor  $q$  in the output is guaranteed to be prime.

EXAMPLES:

```

sage: f = ECM()
sage: n = 508021860739623467191080372196682785441177798407961
sage: f.one_curve(n, B1=10000, sigma=11)
[1, 508021860739623467191080372196682785441177798407961]
sage: f.one_curve(n, B1=10000, sigma=1022170541)
[79792266297612017, 6366805760909027985741435139224233]
sage: n = 432132887883903108009802143314445113500016816977037257
sage: f.one_curve(n, B1=500000, algorithm="P-1")
[67872792749091946529, 6366805760909027985741435139224233]
sage: n = 2088352670731726262548647919416588631875815083
sage: f.one_curve(n, B1=2000, algorithm="P+1", x0=5)
[328006342451, 6366805760909027985741435139224233]

```

**recommended\_B1** (*factor\_digits*)

Return recommended B1 setting.

INPUT:

- *factor\_digits* – integer. Number of digits.

OUTPUT:

Integer. Recommended settings from [http://www.mersennewiki.org/index.php/Elliptic\\_Curve\\_Method](http://www.mersennewiki.org/index.php/Elliptic_Curve_Method)

EXAMPLES:

```

sage: ecm.recommended_B1(33)
1000000

```

**time** (*n*, *factor\_digits*, *verbose=False*)

Print a runtime estimate.

BUGS:

This method should really return something and not just print stuff on the screen.

INPUT:

- *n* – a positive integer
- *factor\_digits* – the (estimated) number of digits of the smallest factor

OUTPUT:

An approximation for the amount of time it will take to find a factor of size *factor\_digits* in a single process on the current computer. This estimate is provided by GMP-ECM's verbose option on a single run of a curve.

EXAMPLES:

```

sage: n = next_prime(11^23)*next_prime(11^37)
sage: ecm.time(n, 35)                                     # random output
Expected curves: 910, Expected time: 23.95m

sage: ecm.time(n, 30, verbose=True)                       # random output
GMP-ECM 6.4.4 [configured with MPFR 2.6.0, --enable-asm-redc] [ECM]
Running on localhost.localdomain
Input number is 304481639541418099574459496544854621998616257489887231115912293 (63 digits)
Using MODMULN [mulredc:0, sqrredc:0]
Using B1=250000, B2=128992510, polynomial Dickson(3), sigma=3244548117
dF=2048, k=3, d=19110, d2=11, i0=3
Expected number of curves to find a factor of n digits:
35 40 45 50 55 60 65 70 75 80

```

```
4911 70940 1226976 2.5e+07 5.8e+08 1.6e+10 2.7e+13 4e+18 5.4e+23 Inf
Step 1 took 230ms
Using 10 small primes for NTT
Estimated memory usage: 4040K
Initializing tables of differences for F took 0ms
Computing roots of F took 9ms
Building F from its roots took 16ms
Computing 1/F took 9ms
Initializing table of differences for G took 0ms
Computing roots of G took 8ms
Building G from its roots took 16ms
Computing roots of G took 7ms
Building G from its roots took 16ms
Computing G * H took 6ms
Reducing G * H mod F took 5ms
Computing roots of G took 7ms
Building G from its roots took 17ms
Computing G * H took 5ms
Reducing G * H mod F took 5ms
Computing polyeval(F,G) took 34ms
Computing product of all F(g_i) took 0ms
Step 2 took 164ms
Expected time to find a factor of n digits:
35 40 45 50 55 60 65 70 75 80
32.25m 7.76h 5.60d 114.21d 7.27y 196.42y 337811y 5e+10y 7e+15y Inf

Expected curves: 4911, Expected time: 32.25m
```



## INTERFACE TO 4TI2

<http://www.4ti2.de>

You must have the 4ti2 Sage package installed on your computer for this interface to work.

Use `sage -i 4ti2` to install the package.

AUTHORS:

- Mike Hansen (2009): Initial version.
- Bjarke Hammersholt Rouné (2009-06-26): Added Groebner, made code usable as part of the Sage library and added documentation and some doctests.
- Marshall Hampton (2011): Minor fixes to documentation.

**class** `sage.interfaces.four_ti_2.FourTi2` (*directory=None*)  
Bases: `object`

This object defines an interface to the program 4ti2. Each command 4ti2 has is exposed as one method.

**call** (*command, project, verbose=True*)  
Run the 4ti2 program `command` on the project named `project` in the directory `directory()`.

INPUT:

- `command` – The 4ti2 program to run.
- `project` – The file name of the project to run on.
- `verbose` – Display the output of 4ti2 if `True`.

EXAMPLES:

```
sage: from sage.interfaces.four_ti_2 import four_ti_2
sage: four_ti_2.write_matrix([[6,10,15]], "test_file")
sage: four_ti_2.call("groebner", "test_file", False) # optional - 4ti2
sage: four_ti_2.read_matrix("test_file.gro") # optional - 4ti2
[-5  0  2]
[-5  3  0]
```

**circuits** (*mat=None, project=None*)

Run the 4ti2 program `circuits` on the parameters. See <http://www.4ti2.de/> for details.

EXAMPLES:

```
sage: from sage.interfaces.four_ti_2 import four_ti_2
sage: four_ti_2.circuits([1,2,3]) # optional - 4ti2
[ 0  3 -2]
[ 2 -1  0]
[ 3  0 -1]
```

**directory()**

Return the directory where the input files for 4ti2 are written by Sage and where 4ti2 is run.

EXAMPLES:

```
sage: from sage.interfaces.four_ti_2 import FourTi2
sage: f = FourTi2("/tmp/")
sage: f.directory()
'/tmp/'
```

**graver** (*mat=None, lat=None, project=None*)

Run the 4ti2 program graver on the parameters. See <http://www.4ti2.de/> for details.

EXAMPLES:

```
sage: from sage.interfaces.four_ti_2 import four_ti_2
sage: four_ti_2.graver([1,2,3]) # optional - 4ti2
[ 2 -1  0]
[ 3  0 -1]
[ 1  1 -1]
[ 1 -2  1]
[ 0  3 -2]
sage: four_ti_2.graver(lat=[[1,2,3],[1,1,1]]) # optional - 4ti2
[ 1  0 -1]
[ 0  1  2]
[ 1  1  1]
[ 2  1  0]
```

**groebner** (*mat=None, lat=None, project=None*)

Run the 4ti2 program groebner on the parameters. This computes a Toric Groebner basis of a matrix. See <http://www.4ti2.de/> for details.

EXAMPLES:

```
sage: from sage.interfaces.four_ti_2 import four_ti_2
sage: A = [6,10,15]
sage: four_ti_2.groebner(A) # optional - 4ti2
[-5  0  2]
[-5  3  0]
sage: four_ti_2.groebner(lat=[[1,2,3],[1,1,1]]) # optional - 4ti2
[-1  0  1]
[ 2  1  0]
```

**hilbert** (*mat=None, lat=None, project=None*)

Run the 4ti2 program hilbert on the parameters. See <http://www.4ti2.de/> for details.

EXAMPLES:

```
sage: from sage.interfaces.four_ti_2 import four_ti_2
sage: four_ti_2.hilbert(four_ti_2._magic3x3()) # optional - 4ti2
[2 0 1 0 1 2 1 2 0]
[1 0 2 2 1 0 0 2 1]
[0 2 1 2 1 0 1 0 2]
[1 2 0 0 1 2 2 0 1]
[1 1 1 1 1 1 1 1 1]
sage: four_ti_2.hilbert(lat=[[1,2,3],[1,1,1]]) # optional - 4ti2
[2 1 0]
[0 1 2]
[1 1 1]
```

**minimize** (*mat=None, lat=None*)

Run the 4ti2 program minimize on the parameters. See <http://www.4ti2.de/> for details.

EXAMPLES:

```
sage: from sage.interfaces.four_ti_2 import four_ti_2
sage: four_ti_2.minimize() # optional - 4ti2
Traceback (most recent call last):
...
NotImplementedError: 4ti2 command 'minimize' not implemented in Sage.
```

**ppi**(*n*)

Run the 4ti2 program ppi on the parameters. See <http://www.4ti2.de/> for details.

EXAMPLES:

```
sage: from sage.interfaces.four_ti_2 import four_ti_2
sage: four_ti_2.ppi(3) # optional - 4ti2
[-2  1  0]
[ 0 -3  2]
[-1 -1  1]
[-3  0  1]
[ 1 -2  1]
```

**qsolve**(*mat=None, rel=None, sign=None, project=None*)

Run the 4ti2 program qsolve on the parameters. See <http://www.4ti2.de/> for details.

EXAMPLES:

```
sage: from sage.interfaces.four_ti_2 import four_ti_2
sage: A = [[1,1,1],[1,2,3]]
sage: four_ti_2.qsolve(A) # optional - 4ti2
[[], [ 1 -2  1]]
```

**rays**(*mat=None, project=None*)

Run the 4ti2 program rays on the parameters. See <http://www.4ti2.de/> for details.

EXAMPLES:

```
sage: from sage.interfaces.four_ti_2 import four_ti_2
sage: four_ti_2.rays(four_ti_2._magic3x3()) # optional - 4ti2
[0 2 1 2 1 0 1 0 2]
[1 0 2 2 1 0 0 2 1]
[1 2 0 0 1 2 2 0 1]
[2 0 1 0 1 2 1 2 0]
```

**read\_matrix**(*filename*)

Read a matrix in 4ti2 format from the file *filename* in directory *directory*() .

INPUT:

- *filename* – The name of the file to read from.

OUTPUT:

The data from the file as a matrix over  $\mathbb{Z}$ .

EXAMPLES:

```
sage: from sage.interfaces.four_ti_2 import four_ti_2
sage: four_ti_2.write_matrix([[1,2,3],[3,4,6]], "test_file")
sage: four_ti_2.read_matrix("test_file")
[1 2 3]
[3 4 6]
```

**temp\_project()**

Return an input project file name that has not been used yet.

EXAMPLES:

```
sage: from sage.interfaces.four_ti_2 import four_ti_2
sage: four_ti_2.temp_project()
'project_...'
```

**write\_array(array, nrows, ncols, filename)**

Write the matrix array of integers (can be represented as a list of lists) to the file filename in directory directory() in 4ti2 format. The matrix must have nrows rows and ncols columns.

INPUT:

- array – A matrix of integers. Can be represented as a list of lists.
- nrows – The number of rows in array.
- ncols – The number of columns in array.
- file – A file name not including a path.

EXAMPLES:

```
sage: from sage.interfaces.four_ti_2 import four_ti_2
sage: four_ti_2.write_array([[1,2,3],[3,4,5]], 2, 3, "test_file")
```

**write\_matrix(mat, filename)**

Write the matrix mat to the file filename in 4ti2 format.

INPUT:

- mat – A matrix of integers or something that can be converted to that.
- filename – A file name not including a path.

EXAMPLES:

```
sage: from sage.interfaces.four_ti_2 import four_ti_2
sage: four_ti_2.write_matrix([[1,2],[3,4]], "test_file")
```

**write\_single\_row(row, filename)**

Write the list row to the file filename in 4ti2 format as a matrix with one row.

INPUT:

- row – A list of integers.
- filename – A file name not including a path.

EXAMPLES:

```
sage: from sage.interfaces.four_ti_2 import four_ti_2
sage: four_ti_2.write_single_row([1,2,3,4], "test_file")
```

**zsolve(mat=None, rel=None, rhs=None, sign=None, lat=None, project=None)**

Run the 4ti2 program zsolve on the parameters. See <http://www.4ti2.de/> for details.

EXAMPLES:

```
sage: from sage.interfaces.four_ti_2 import four_ti_2
sage: A = [[1,1,1],[1,2,3]]
sage: rel = ['<', '<']
sage: rhs = [2, 3]
sage: sign = [1,0,1]
```

```
sage: four_ti_2.zsolve(A, rel, rhs, sign) # optional - 4ti2
[
      [ 1 -1  0]
      [ 0 -1  0]
[0 0 1] [ 0 -3  2]
[1 1 0] [ 1 -2  1]
[0 1 0], [ 0 -2  1], []
]
sage: four_ti_2.zsolve(lat=[[1,2,3],[1,1,1]]) # optional - 4ti2
[
      [1 2 3]
[0 0 0], [], [1 1 1]
]
```



## INTERFACE TO FRICAS

---

### Todo

- Evaluation using a file is not done. Any input line with more than a few thousand characters would hang the system, so currently it automatically raises an exception.
  - All completions of a given command.
  - Interactive help.
- 

FriCAS is a free GPL-compatible (modified BSD license) general purpose computer algebra system based on Axiom. The FriCAS website can be found at <http://fricas.sourceforge.net/>.

### AUTHORS:

- Mike Hansen (2009-02): Split off the FriCAS interface from the Axiom interface.

If the string “error” (case insensitive) occurs in the output of anything from FriCAS, a `RuntimeError` exception is raised.

EXAMPLES: We evaluate a very simple expression in FriCAS.

```
sage: fricas('3 * 5')                # optional - fricas
15
sage: a = fricas(3) * fricas(5); a    # optional - fricas
15
```

The type of `a` is `FriCASElement`, i.e., an element of the FriCAS interpreter.

```
sage: type(a)                        # optional - fricas
<class 'sage.interfaces.fricas.FriCASElement'>
sage: a.parent()                     # optional - fricas
FriCAS
```

The underlying FriCAS type of `a` is also available, via the `type` method:

```
sage: a.type()                       # optional - fricas
PositiveInteger
```

We factor  $x^5 - y^5$  in FriCAS in several different ways. The first way yields a FriCAS object.

```
sage: F = fricas.factor('x^5 - y^5'); F    # optional - fricas
      4      3      2 2      3      4
- (y - x) (y  + x y  + x y  + x y + x )
sage: type(F)                            # optional - fricas
<class 'sage.interfaces.fricas.FriCASElement'>
```

```
sage: F.type()                                # optional - fricas
Factored(Polynomial(Integer))
```

Note that FriCAS objects are normally displayed using “ASCII art”.

```
sage: a = fricas(2/3); a                      # optional - fricas
  2
  -
  3
sage: a = fricas('x^2 + 3/7'); a              # optional - fricas
  2   3
x  + -
  7
```

The `fricas.eval` command evaluates an expression in FriCAS and returns the result as a string. This is exact as if we typed in the given line of code to FriCAS; the return value is what FriCAS would print out.

```
sage: print fricas.eval('factor(x^5 - y^5)')  # optional - fricas
      4      3      2 2      3      4
- (y - x)(y  + x y  + x y  + x y + x )      Type: Factored(Polynomial(Integer))
```

We can create the polynomial  $f$  as a FriCAS polynomial, then call the `factor` method on it. Notice that the notation `f.factor()` is consistent with how the rest of Sage works.

```
sage: f = fricas('x^5 - y^5')                # optional - fricas
sage: f^2                                     # optional - fricas
      10      5 5      10
y  - 2x y  + x
sage: f.factor()                             # optional - fricas
      4      3      2 2      3      4
- (y - x)(y  + x y  + x y  + x y + x )
```

Control-C interruption works well with the FriCAS interface. For example, try the following sum but with a much bigger range, and hit control-C.

```
sage: f = fricas('(x^5 - y^5)^10000')        # not tested - fricas
Interrupting FriCAS...
...
<type 'exceptions.TypeError': Ctrl-c pressed while running FriCAS>

sage: fricas('1/100 + 1/101')                # optional - fricas
      201
-----
      10100
sage: a = fricas('(1 + sqrt(2))^5'); a        # optional - fricas
      +-+
      29\|2  + 41
```

TESTS: We check to make sure the `subst` method works with keyword arguments.

```
sage: a = fricas(x+2); a                     #optional - fricas
x + 2
sage: a.subst(x=3)                           #optional - fricas
5
```

We verify that FriCAS floating point numbers can be converted to Python floats.



```
sage: float(fricas(2))      #optional - fricas
2.0
```

```
class sage.interfaces.fricas.FriCAS(name='fricas', command='fricas -nox -noclef',
                                     script_subdirectory=None, logfile=None, server=None,
                                     server_tmpdir=None, init_code=['(si::readline-off)'])
```

Bases: `sage.interfaces.axiom.PanAxiom`

Interface to a FriCAS interpreter.

**console()**

Spawn a new FriCAS command-line session.

EXAMPLES:

```
sage: fricas.console() #not tested
FriCAS (AXIOM fork) Computer Algebra System
Version: FriCAS 1.0.5
Timestamp: Thursday February 19, 2009 at 06:57:33
```

```
-----
Issue )copyright to view copyright notices.
Issue )summary for a summary of useful system commands.
Issue )quit to leave AXIOM and return to shell.
-----
```

```
class sage.interfaces.fricas.FriCASElement(parent, value, is_name=False, name=None)
```

Bases: `sage.interfaces.axiom.PanAxiomElement`

```
class sage.interfaces.fricas.FriCASExpectFunction(parent, name)
```

Bases: `sage.interfaces.axiom.PanAxiomExpectFunction`

TESTS:

```
sage: axiom.upperCase_q
upperCase?
sage: axiom.upperCase_e
upperCase!
```

```
class sage.interfaces.fricas.FriCASFunctionElement(object, name)
```

Bases: `sage.interfaces.axiom.PanAxiomFunctionElement`

TESTS:

```
sage: a = axiom("Hello") #optional - axiom
sage: a.upperCase_q      #optional - axiom
upperCase?
sage: a.upperCase_e      #optional - axiom
upperCase!
sage: a.upperCase_e()    #optional - axiom
"HELLO"
```

```
sage.interfaces.fricas.fricas_console()
```

Spawn a new FriCAS command-line session.

EXAMPLES:

```
sage: fricas_console() #not tested
FriCAS (AXIOM fork) Computer Algebra System
Version: FriCAS 1.0.5
Timestamp: Thursday February 19, 2009 at 06:57:33
```

```
-----
Issue )copyright to view copyright notices.
```

```
Issue )summary for a summary of useful system commands.  
Issue )quit to leave AXIOM and return to shell.
```

-----

`sage.interfaces.fricas.is_FriCASElement(x)`

Return True if x is of type FriCASElement.

EXAMPLES:

```
sage: from sage.interfaces.fricas import is_FriCASElement #optional - fricas  
sage: is_FriCASElement(fricas(2)) #optional - fricas  
True  
sage: is_FriCASElement(2) #optional - fricas  
False
```

`sage.interfaces.fricas.reduce_load_fricas()`

Returns the FriCAS interface object defined in `sage.interfaces.fricas`.

EXAMPLES:

```
sage: from sage.interfaces.fricas import reduce_load_fricas  
sage: reduce_load_fricas()  
FriCAS
```

## INTERFACE TO FROBBY FOR FAST COMPUTATIONS ON MONOMIAL IDEALS.

The software package Frobbly provides a number of computations on monomial ideals. The current main feature is the socle of a monomial ideal, which is largely equivalent to computing the maximal standard monomials, the Alexander dual or the irreducible decomposition.

Operations on monomial ideals are much faster than algorithms designed for ideals in general, which is what makes a specialized library for these operations on monomial ideals useful.

AUTHORS:

- Bjarke Hammersholt Roune (2008-04-25): Wrote the Frobbly C++ program and the initial version of the Python interface.

NOTES:

The official source for Frobbly is <<http://www.broune.com/frobbly>>, which also has documentation and papers describing the algorithms used.

**class** sage.interfaces.frobbly.Frobbly

**alexander\_dual** (*monomial\_ideal*)

This function computes the Alexander dual of the passed-in monomial ideal. This ideal is the one corresponding to the simplicial complex whose faces are the complements of the nonfaces of the simplicial complex corresponding to the input ideal.

INPUT:

- *monomial\_ideal* – The monomial ideal to decompose.

OUTPUT:

The monomial corresponding to the Alexander dual.

EXAMPLES:

This is a simple example of computing irreducible decomposition.

```
sage: (a, b, c, d) = QQ['a,b,c,d'].gens() # optional - frobbly
sage: id = ideal(a * b, b * c, c * d, d * a) # optional - frobbly
sage: alexander_dual = frobbly.alexander_dual(id) # optional - frobbly
sage: true_alexander_dual = ideal(b * d, a * c) # optional - frobbly
sage: alexander_dual == true_alexander_dual # use sets to ignore order # optional - frobbly
True
```

We see how it is much faster to compute this with frobbly than the built-in procedure for simplicial complexes.

```
sage: t=simplicial_complexes.PoincareHomologyThreeSphere() # optional - frobby sage:
R=PolynomialRing(QQ,16,'x') # optional - frobby sage: I=R.ideal([prod([R.gen(i-1) for i in a]
for a in t.facets())]) # optional - frobby sage: len(frobby.alexander_dual(I).gens()) # optional -
frobby 643
```

**associated\_primes** (*monomial\_ideal*)

This function computes the associated primes of the passed-in monomial ideal.

INPUT:

- `monomial_ideal` – The monomial ideal to decompose.

OUTPUT:

A list of the associated primes of the monomial ideal. These ideals are constructed in the same ring as `monomial_ideal` is.

EXAMPLES:

```
sage: R.<d,b,c>=QQ[] # optional - frobby
sage: I=[d*b*c,b^2*c,b^10,d^10]*R # optional - frobby
sage: frobby.associated_primes(I) # optional - frobby
[Ideal (d, b) of Multivariate Polynomial Ring in d, b, c over Rational Field,
Ideal (d, b, c) of Multivariate Polynomial Ring in d, b, c over Rational Field]
```

**dimension** (*monomial\_ideal*)

This function computes the dimension of the passed-in monomial ideal.

INPUT:

- `monomial_ideal` – The monomial ideal to decompose.

OUTPUT:

The dimension of the zero set of the ideal.

EXAMPLES:

```
sage: R.<d,b,c>=QQ[] # optional - frobby
sage: I=[d*b*c,b^2*c,b^10,d^10]*R # optional - frobby
sage: frobby.dimension(I) # optional - frobby
1
```

**hilbert** (*monomial\_ideal*)

Computes the multigraded Hilbert-Poincare series of the input ideal. Use the `-univariate` option to get the univariate series.

The Hilbert-Poincare series of a monomial ideal is the sum of all monomials not in the ideal. This sum can be written as a (finite) rational function with  $(x_1 - 1)(x_2 - 1)\dots(x_n - 1)$  in the denominator, assuming the variables of the ring are  $x_1, x_2, \dots, x_n$ . This action computes the polynomial in the numerator of this fraction.

INPUT:

- `monomial_ideal` – A monomial ideal.

OUTPUT:

- A polynomial in the same ring as the ideal.

EXAMPLES:

```
sage: R.<d,b,c>=QQ[] # optional - frobby
sage: I=[d*b*c,b^2*c,b^10,d^10]*R # optional - frobby
```

```
sage: frobby.hilbert(I) # optional - frobby
d^10*b^10*c + d^10*b^10 + d^10*b*c + b^10*c + d^10 + b^10 + d*b^2*c + d*b*c + b^2*c + 1
```

### **irreducible\_decomposition** (*monomial\_ideal*)

This function computes the irreducible decomposition of the passed-in monomial ideal. I.e. it computes the unique minimal list of irreducible monomial ideals whose intersection equals `monomial_ideal`.

INPUT:

- `monomial_ideal` – The monomial ideal to decompose.

OUTPUT:

A list of the unique irredundant irreducible components of `monomial_ideal`. These ideals are constructed in the same ring as `monomial_ideal` is.

EXAMPLES:

This is a simple example of computing irreducible decomposition.

```
sage: (x, y, z) = QQ['x,y,z'].gens() # optional - frobby
sage: id = ideal(x ** 2, y ** 2, x * z, y * z) # optional - frobby
sage: decom = frobby.irreducible_decomposition(id) # optional - frobby
sage: true_decom = [ideal(x, y), ideal(x ** 2, y ** 2, z)] # optional - frobby
sage: set(decom) == set(true_decom) # use sets to ignore order # optional - frobby
True
```

We now try the special case of the zero ideal in different rings.

We should also try `PolynomialRing(QQ, names=[])`, but it has a bug which makes that impossible (see trac ticket 3028).

```
sage: rings = [ZZ['x'], CC['x,y']] # optional - frobby
sage: allOK = True # optional - frobby
sage: for ring in rings: # optional - frobby
...     id0 = ring.ideal(0) # optional - frobby
...     decom0 = frobby.irreducible_decomposition(id0) # optional - frobby
...     allOK = allOK and decom0 == [id0] # optional - frobby
sage: allOK # optional - frobby
True
```

Finally, we try the ideal that is all of the ring in different rings.

```
sage: rings = [ZZ['x'], CC['x,y']] # optional - frobby
sage: allOK = True # optional - frobby
sage: for ring in rings: # optional - frobby
...     id1 = ring.ideal(1) # optional - frobby
...     decom1 = frobby.irreducible_decomposition(id1) # optional - frobby
...     allOK = allOK and decom1 == [id1] # optional - frobby
sage: allOK # optional - frobby
True
```



## INTERFACE TO GAP

Sage provides an interface to the GAP system. This system provides extensive group theory, combinatorics, etc.

The GAP interface will only work if GAP is installed on your computer; this should be the case, since GAP is included with Sage. The interface offers three pieces of functionality:

1. `gap_console()` - A function that dumps you into an interactive command-line GAP session.
2. `gap(expr)` - Evaluation of arbitrary GAP expressions, with the result returned as a string.
3. `gap.new(expr)` - Creation of a Sage object that wraps a GAP object. This provides a Pythonic interface to GAP. For example, if `f=gap.new(10)`, then `f.Factors()` returns the prime factorization of 10 computed using GAP.

### 9.1 First Examples

We factor an integer using GAP:

```
sage: n = gap(20062006); n
20062006
sage: n.parent()
Gap
sage: fac = n.Factors(); fac
[ 2, 17, 59, 73, 137 ]
sage: fac.parent()
Gap
sage: fac[1]
2
```

### 9.2 GAP and Singular

This example illustrates conversion between Singular and GAP via Sage as an intermediate step. First we create and factor a Singular polynomial.

```
sage: singular(389)
389
sage: R1 = singular.ring(0, '(x,y)', 'dp')
sage: f = singular('9*x^16-18*x^13*y^2-9*x^12*y^3+9*x^10*y^4-18*x^11*y^2+36*x^8*y^4+18*x^7*y^5-18*x^5*y^3+9*x^4*y^2-9*x^3*y+9*x^2*y-9*x*y+9')
sage: F = f.factorize()
sage: print F
[1]:
  _[1]=9
```

```
_[2]=x^6-2*x^3*y^2-x^2*y^3+y^4
_[3]=-x^5+y^2
[2]:
1, 1, 2
```

Next we convert the factor  $-x^5 + y^2$  to a Sage multivariate polynomial. Note that it is important to let  $x$  and  $y$  be the generators of a polynomial ring, so the `eval` command works.

```
sage: R.<x,y> = PolynomialRing(QQ,2)
sage: s = F[1][3].sage_polysttring(); s
'-x**5+y**2'
sage: g = eval(s); g
-x^5 + y^2
```

Next we create a polynomial ring in GAP and obtain its indeterminates:

```
sage: R = gap.PolynomialRing('Rationals', 2); R
PolynomialRing( Rationals, ["x_1", "x_2"] )
sage: I = R.IndeterminatesOfPolynomialRing(); I
[ x_1, x_2 ]
```

In order to eval  $g$  in GAP, we need to tell GAP to view the variables  $x_0$  and  $x_1$  as the two generators of  $R$ . This is the one tricky part. In the GAP interpreter the object `I` has its own name (which isn't `I`). We can access its name using `I.name()`.

```
sage: _ = gap.eval("x := %s[1];; y := %s[2];;"%(I.name(), I.name()))
```

Now  $x_0$  and  $x_1$  are defined, so we can construct the GAP polynomial  $f$  corresponding to  $g$ :

```
sage: R.<x,y> = PolynomialRing(QQ,2)
sage: f = gap(str(g)); f
-x_1^5+x_2^2
```

We can call GAP functions on  $f$ . For example, we evaluate the GAP `Value` function, which evaluates  $f$  at the point  $(1, 2)$ .

```
sage: f.Value(I, [1,2])
3
sage: g(1,2)           # agrees
3
```

## 9.3 Saving and loading objects

Saving and loading GAP objects (using the `dumps` method, etc.) is *not* supported, since the output string representation of Gap objects is sometimes not valid input to GAP. Creating classes that wrap GAP objects *is* supported, via simply defining the `_gap_init_` member function that returns a string that when evaluated in GAP constructs the object. See `groups/perm_gps/permgroup.py` for a nontrivial example of this.

## 9.4 Long Input

The GAP interface reads in even very long input (using files) in a robust manner, as long as you are creating a new object.



---

**Note:** Using `gap.eval` for long input is much less robust, and is not recommended.

---

```
sage: t = '"%s"%10^10000' # ten thousand character string.
sage: a = gap(t)
```

## 9.5 Changing which GAP is used

Use this code to change which GAP interpreter is run. E.g.,

```
import sage.interfaces.gap
sage.interfaces.gap.gap_cmd = "/usr/local/bin/gap"
```

AUTHORS:

- David Joyner and William Stein: initial version(s)
- William Stein (2006-02-01): modified `gap_console` command so it uses exactly the same startup command as `Gap.__init__`.
- William Stein (2006-03-02): added tab completions: `gap.[tab]`, `x = gap(...)`, `x.[tab]`, and docs, e.g., `gap.function?` and `x.function?`

```
class sage.interfaces.gap.Gap(max_workspace_size=None, maxread=100000,
                              script_subdirectory=None, use_workspace_cache=True,
                              server=None, server_tmpdir=None, logfile=None, seed=None)
```

Bases: `sage.interfaces.gap.Gap_generic`

Interface to the GAP interpreter.

AUTHORS:

- William Stein and David Joyner

**console()**

Spawn a new GAP command-line session.

EXAMPLES:

```
sage: gap.console() # not tested
*****      GAP, Version 4.5.7 of 14-Dec-2012 (free software, GPL)
*   GAP   *   http://www.gap-system.org
*****      Architecture: x86_64-unknown-linux-gnu-gcc-default64
Libs used:  gmp, readline
Loading the library and packages ...
Packages:   GAPDoc 1.5.1
Try '?help' for help. See also '?copyright' and '?authors'
gap>
```

**cputime(t=None)**

Returns the amount of CPU time that the GAP session has used. If `t` is not `None`, then it returns the difference between the current CPU time and `t`.

EXAMPLES:

```
sage: t = gap.cputime()
sage: t #random
0.13600000000000001
sage: gap.Order(gap.SymmetricGroup(5))
```

```
120
sage: gap.cputime(t)    #random
0.05999999999999998
```

**get** (*var*, *use\_file=False*)

Get the string representation of the variable *var*.

EXAMPLES:

```
sage: gap.set('x', '2')
sage: gap.get('x')
'2'
```

**help** (*s*, *pager=True*)

Print help on a given topic.

EXAMPLES:

```
sage: print gap.help('SymmetricGroup', pager=False)
```

```
50 Group Libraries
```

```
When you start GAP, it already knows several groups. Currently GAP initially
knows the following groups:
```

```
...
```

**save\_workspace** ()

Save the GAP workspace.

TESTS:

We make sure that #9938 (GAP does not start if the path to the GAP workspace file contains more than 82 characters) is fixed:

```
sage: ORIGINAL_WORKSPACE = sage.interfaces.gap.WORKSPACE
sage: sage.interfaces.gap.WORKSPACE = os.path.join(SAGE_TMP, "gap" + "0"*(80-len(SAGE_TMP)))
sage: gap = Gap()
sage: gap('3+2')    # long time (4s on sage.math, 2013)
5
sage: sage.interfaces.gap.WORKSPACE = ORIGINAL_WORKSPACE
```

**set** (*var*, *value*)

Set the variable *var* to the given value.

EXAMPLES:

```
sage: gap.set('x', '2')
sage: gap.get('x')
'2'
```

**set\_seed** (*seed=None*)

Sets the seed for gap interpreter. The seed should be an integer.

EXAMPLES:

```
sage: g = Gap()
sage: g.set_seed(0)
0
sage: [g.Random(1,10) for i in range(5)]
[2, 3, 3, 4, 2]
```

```

trait_names()
    EXAMPLES:
    sage: c = gap.trait_names()
    sage: len(c) > 100
    True
    sage: 'Order' in c
    True

```

```

class sage.interfaces.gap.GapElement (parent, value, is_name=False, name=None)
    Bases: sage.interfaces.gap.GapElement_generic

```

```

str (use_file=False)
    EXAMPLES:
    sage: print gap(2)
    2

```

```

trait_names()
    EXAMPLES:
    sage: s5 = gap.SymmetricGroup(5)
    sage: 'Centralizer' in s5.trait_names()
    True

```

```

class sage.interfaces.gap.GapElement_generic (parent, value, is_name=False, name=None)
    Bases: sage.interfaces.expect.ExpectElement

```

Generic interface to the GAP3/GAP4 interpreters.

AUTHORS:

- William Stein and David Joyner (interface for GAP4)
- Franco Saliola (Feb 2010): refactored to separate out the generic code

```

bool()
    EXAMPLES:
    sage: bool(gap(2))
    True
    sage: gap(0).bool()
    False
    sage: gap('false').bool()
    False

```

```

is_string()
    Tell whether this element is a string.
    EXAMPLES:
    sage: gap('"abc"').is_string()
    True
    sage: gap('[1,2,3]').is_string()
    False

```

```

class sage.interfaces.gap.GapFunction (parent, name)
    Bases: sage.interfaces.expect.ExpectFunction

```

```

class sage.interfaces.gap.GapFunctionElement (obj, name)
    Bases: sage.interfaces.expect.FunctionElement

```

```
class sage.interfaces.gap.Gap_generic(name, prompt, command=None, server=None,
                                     server_tmpdir=None, ulimit=None,
                                     maxread=100000, script_subdirectory=None,
                                     restart_on_ctrlc=False, verbose_start=False,
                                     init_code=[], max_startup_time=None, logfile=None,
                                     eval_using_file_cutoff=0, do_cleaner=True, re-
                                     mote_cleaner=False, path=None, terminal_echo=True)
```

Bases: `sage.interfaces.expect.Expect`

Generic interface to the GAP3/GAP4 interpreters.

AUTHORS:

- William Stein and David Joyner (interface for GAP4)
- Franco Saliola (Feb 2010): refactored to separate out the generic code

**eval** (*x*, *newlines=False*, *strip=True*, *split\_lines=True*, *\*\*kwds*)

Send the code in the string *s* to the GAP interpreter and return the output as a string.

INPUT:

- s* - string containing GAP code.
- newlines* - bool (default: True); if False, remove all backslash-newlines inserted by the GAP output formatter.
- strip* - ignored
- split\_lines* - bool (default: True); if True then each line is evaluated separately. If False, then the whole block of code is evaluated all at once.

EXAMPLES:

```
sage: gap.eval('2+2')
'4'
sage: gap.eval('Print(4); #test\n Print(6);')
'46'
sage: gap.eval('Print("#"); Print(6);')
'#6'
sage: gap.eval('4; \n 6;')
'4\n6'
sage: gap.eval('if 3>2 then\nPrint("hi");\nfi;')
'hi'
sage: gap.eval('## this is a test\nPrint("OK")')
'OK'
sage: gap.eval('Print("This is a test. Oh no, a #");# but this is a comment\nPrint("OK")')
'This is a test. Oh no, a #OK'
sage: gap.eval('if 4>3 then')
''
sage: gap.eval('Print("Hi how are you?")')
'Hi how are you?'
sage: gap.eval('fi')
''
```

**function\_call** (*function*, *args=None*, *kwds=None*)

Calls the GAP function with *args* and *kwds*.

EXAMPLES:

```
sage: gap.function_call('SymmetricGroup', [5])
SymmetricGroup( [ 1 .. 5 ] )
```

If the GAP function does not return a value, but prints something to the screen, then a string of the printed output is returned.

```
sage: s = gap.function_call('Display', [gap.SymmetricGroup(5).CharacterTable()])
sage: type(s)
<class 'sage.interfaces.interface.AsciiArtString'>
sage: s.startswith('CT')
True
```

**get\_record\_element** (*record, name*)

Return the element of a GAP record identified by name.

INPUT:

- record – a GAP record
- name – str

OUTPUT:

- GapElement

EXAMPLES:

```
sage: rec = gap('rec( a := 1, b := "2" )')
sage: gap.get_record_element(rec, 'a')
1
sage: gap.get_record_element(rec, 'b')
2
```

TESTS:

```
sage: rec = gap('rec( a := 1, b := "2" )')
sage: type(gap.get_record_element(rec, 'a'))
<class 'sage.interfaces.gap.GapElement'>
```

**interrupt** (*tries=None, timeout=1, quit\_on\_fail=True*)

Interrupt the GAP process

Gap installs a SIGINT handler, we call it directly instead of trying to sent Ctrl-C. Unlike `interrupt()`, we only try once since we are knowing what we are doing.

Sometimes GAP dies while interrupting.

EXAMPLES:

```
sage: gap._eval_line('while(1=1) do i:=1;; od;', wait_for_prompt=False);
,,
sage: rc = gap.interrupt(timeout=1)
sage: [ gap(i) for i in range(10) ] # check that it is still working
[0, 1, 2, 3, 4, 5, 6, 7, 8, 9]
```

TESTS:

```
sage: gap('"finished computation"); gap.interrupt(); gap('"ok"')
finished computation
True
ok
```

**load\_package** (*pkg, verbose=False*)

Load the Gap package with the given name.

If loading fails, raise a `RuntimeError` exception.

TESTS:

```
sage: gap.load_package("chevie")
Traceback (most recent call last):
...
RuntimeError: Error loading Gap package chevie. You may want to install the gap_packages and
```

**trait\_names()**

EXAMPLES:

```
sage: c = gap.trait_names()
sage: len(c) > 100
True
sage: 'Order' in c
True
```

**unbind(var)**

Clear the variable named var.

EXAMPLES:

```
sage: gap.set('x', '2')
sage: gap.get('x')
'2'
sage: gap.unbind('x')
sage: gap.get('x')
Traceback (most recent call last):
...
RuntimeError: Gap produced error output
Error, Variable: 'x' must have a value
...
```

**version()**

Returns the version of GAP being used.

EXAMPLES:

```
sage: print gap.version()
4.7...
```

`sage.interfaces.gap.gap_command(use_workspace_cache=True, local=True)`

`sage.interfaces.gap.gap_console()`

Spawn a new GAP command-line session.

Note that in gap-4.5.7 you cannot use a workspace cache that had no commandline to restore a gap session with commandline.

EXAMPLES:

```
sage: gap_console() # not tested
*****   GAP, Version 4.5.7 of 14-Dec-2012 (free software, GPL)
*   GAP   *   http://www.gap-system.org
*****   Architecture: x86_64-unknown-linux-gnu-gcc-default64
Libs used: gmp, readline
Loading the library and packages ...
Packages:  GAPDoc 1.5.1
Try '?help' for help. See also  '?copyright' and  '?authors'
gap>
```

TESTS:

```

sage: import subprocess
sage: from sage.interfaces.gap import gap_command
sage: cmd = 'echo "quit;" | ' + gap_command(use_workspace_cache=False)[0]
sage: gap_startup = subprocess.check_output(cmd, shell=True, stderr=subprocess.STDOUT)
sage: 'http://www.gap-system.org' in gap_startup
True
sage: 'Error' not in gap_startup
True
sage: 'sorry' not in gap_startup
True

```

`sage.interfaces.gap.gap_reset_workspace(max_workspace_size=None, verbose=False)`

Call this to completely reset the GAP workspace, which is used by default when Sage first starts GAP.

The first time you start GAP from Sage, it saves the startup state of GAP in a file `$HOME/.sage/gap/workspace-HASH`, where `HASH` is a hash of the directory where Sage is installed.

This is useful, since then subsequent startup of GAP is at least 10 times as fast. Unfortunately, if you install any new code for GAP, it won't be noticed unless you explicitly load it, e.g., with `gap.load_package("my_package")`

The packages `sonata`, `guava`, `factint`, `gapdoc`, `grape`, `design`, `toric`, and `laguna` are loaded in all cases before the workspace is saved, if they are available.

TESTS:

Check that `gap_reset_workspace` still works when `GAP_DIR` doesn't exist, see [trac ticket #14171](#):

```

sage: ORIGINAL_GAP_DIR = sage.interfaces.gap.GAP_DIR
sage: ORIGINAL_WORKSPACE = sage.interfaces.gap.WORKSPACE
sage: sage.interfaces.gap.GAP_DIR = os.path.join(tmp_dir(), "test_gap_dir")
sage: sage.interfaces.gap.WORKSPACE = os.path.join(sage.interfaces.gap.GAP_DIR, "test_workspace")
sage: os.path.isfile(sage.interfaces.gap.WORKSPACE) # long time
False
sage: gap_reset_workspace() # long time
sage: os.path.isfile(sage.interfaces.gap.WORKSPACE) # long time
True
sage: sage.interfaces.gap.GAP_DIR = ORIGINAL_GAP_DIR
sage: sage.interfaces.gap.WORKSPACE = ORIGINAL_WORKSPACE

```

Check that the race condition from [trac ticket #14242](#) has been fixed. We temporarily need to change the worksheet filename.

```

sage: ORIGINAL_WORKSPACE = sage.interfaces.gap.WORKSPACE
sage: sage.interfaces.gap.WORKSPACE = tmp_filename()
sage: from multiprocessing import Process
sage: import time
sage: gap = Gap() # long time (reset GAP session)
sage: P = [Process(target=gap, args=("14242",)) for i in range(4)]
sage: for p in P: # long time, indirect doctest
....:     p.start()
....:     time.sleep(0.2)
sage: for p in P: # long time
....:     p.join()
sage: os.unlink(sage.interfaces.gap.WORKSPACE) # long time
sage: sage.interfaces.gap.WORKSPACE = ORIGINAL_WORKSPACE

```

`sage.interfaces.gap.get_gap_memory_pool_size()`

Get the gap memory pool size for new GAP processes.

EXAMPLES:

```
sage: from sage.interfaces.gap import ... get_gap_memory_pool_size
sage: get_gap_memory_pool_size() # random output
1534059315
```

`sage.interfaces.gap.gfq_gap_to_sage(x, F)`

INPUT:

- `x` – GAP finite field element
- `F` – Sage finite field

OUTPUT: element of `F`

EXAMPLES:

```
sage: x = gap('Z(13)')
sage: F = GF(13, 'a')
sage: F(x)
2
sage: F(gap('0*Z(13)'))
0
sage: F = GF(13^2, 'a')
sage: x = gap('Z(13)')
sage: F(x)
2
sage: x = gap('Z(13^2)^3')
sage: F(x)
12*a + 11
sage: F.multiplicative_generator()^3
12*a + 11
```

TESTS:

Check that [trac ticket #18048](#) is fixed:

```
sage: K.<a> = GF(16)
sage: b = a^2 + a
sage: K(b._gap_())
a^2 + a
```

AUTHOR:

- David Joyner and William Stein

`sage.interfaces.gap.intmod_gap_to_sage(x)`

INPUT:

- `x` – Gap integer mod ring element

EXAMPLES:

```
sage: a = gap(Mod(3, 18)); a
ZmodnZObj( 3, 18 )
sage: b = sage.interfaces.gap.intmod_gap_to_sage(a); b
3
sage: b.parent()
Ring of integers modulo 18

sage: a = gap(Mod(3, 17)); a
Z(17)
sage: b = sage.interfaces.gap.intmod_gap_to_sage(a); b
3
```



```

sage: b.parent()
Finite Field of size 17

sage: a = gap(Mod(0, 17)); a
0*Z(17)
sage: b = sage.interfaces.gap.intmod_gap_to_sage(a); b
0
sage: b.parent()
Finite Field of size 17

sage: a = gap(Mod(3, 65537)); a
ZmodpZObj( 3, 65537 )
sage: b = sage.interfaces.gap.intmod_gap_to_sage(a); b
3
sage: b.parent()
Ring of integers modulo 65537

```

`sage.interfaces.gap.is_GapElement(x)`  
Returns True if x is a GapElement.

EXAMPLES:

```

sage: from sage.interfaces.gap import is_GapElement
sage: is_GapElement(gap(2))
True
sage: is_GapElement(2)
False

```

`sage.interfaces.gap.reduce_load()`  
This is for backwards compatibility only.

To be precise, it only serves at unpickling the invalid gap elements that are stored in the pickle jar.

EXAMPLES:

```

sage: from sage.interfaces.gap import reduce_load
sage: reduce_load()
doctest:...: DeprecationWarning: This function is only used to unpickle invalid objects
See http://trac.sagemath.org/18848 for details.
<repr(<sage.interfaces.gap.GapElement at ...>) failed:
ValueError: The session in which this object was defined is no longer running.>

```

By [trac ticket #18848](#), pickling actually often works:

```

sage: loads(dumps(gap([1, 2, 3])))
[ 1, 2, 3 ]

```

`sage.interfaces.gap.reduce_load_GAP()`  
Returns the GAP interface object defined in `sage.interfaces.gap`.

EXAMPLES:

```

sage: from sage.interfaces.gap import reduce_load_GAP
sage: reduce_load_GAP()
Gap

```

`sage.interfaces.gap.set_gap_memory_pool_size(size_in_bytes)`  
Set the desired gap memory pool size.

Subsequently started GAP/libGAP instances will use this as default. Currently running instances are unchanged.

GAP will only reserve `size_in_bytes` address space. Unless you actually start a big GAP computation, the memory will not be used. However, corresponding swap space will be reserved so that GAP will always be able to use the reserved address space if needed. While nothing is actually written to disc as long as you don't run a big GAP computation, the reserved swap space will not be available for other processes.

INPUT:

- `size_in_bytes` – integer. The desired memory pool size.

EXAMPLES:

```
sage: from sage.interfaces.gap import ... get_gap_memory_pool_size, set_gap_memory
sage: n = get_gap_memory_pool_size()
sage: set_gap_memory_pool_size(n)
sage: n == get_gap_memory_pool_size()
True
sage: n # random output
1534059315
```

## INTERFACE TO GAP3

This module implements an interface to GAP3.

AUTHORS:

- Franco Saliola (February 2010)

**Warning:** GAP3 is not distributed with Sage. You need to install it separately; see the section [Obtaining GAP3](#).

### 10.1 Obtaining GAP3

The GAP3 interface will only work if GAP3 is installed on your computer. Here are some ways to obtain GAP3:

- There is an optional Sage package providing GAP3 pre-packaged with several GAP3 packages:  
[http://trac.sagemath.org/sage\\_trac/ticket/8906](http://trac.sagemath.org/sage_trac/ticket/8906)
- Frank Luebeck maintains a GAP3 Linux executable, optimized for i686 and statically linked for jobs of 2 GByte or more:  
<http://www.math.rwth-aachen.de/~Frank.Luebeck/gap/GAP3>
- Jean Michel maintains a version of GAP3 pre-packaged with CHEVIE and VKCURVE. It can be obtained here:  
<http://people.math.jussieu.fr/~jmichel/chevie/chevie.html>
- Finally, you can download GAP3 from the GAP website below. Since GAP3 is no longer supported, it may not be easy to install this version.  
<http://www.gap-system.org/Gap3/Download3/download.html>

### 10.2 Changing which GAP3 is used

**Warning:** There is a bug in the pexpect module (see trac ticket #8471) that prevents the following from working correctly. For now, just make sure that `gap3` is in your `PATH`.

Sage assumes that GAP3 can be launched with the command `gap3`; that is, Sage assumes that the command `gap3` is in your `PATH`. If this is not the case, then you can start GAP3 using the following command:

```
sage: gap3 = Gap3(command='/usr/local/bin/gap3')           #not tested
```

## 10.3 Functionality and Examples

The interface to GAP3 offers the following functionality.

1. `gap3(expr)` - Evaluation of arbitrary GAP3 expressions, with the result returned as a Sage object wrapping the corresponding GAP3 element:

```
sage: a = gap3('3+2')           #optional - gap3
sage: a                         #optional - gap3
5
sage: type(a)                   #optional - gap3
<class 'sage.interfaces.gap3.GAP3Element'>

sage: S5 = gap3('SymmetricGroup(5)') #optional - gap3
sage: S5                       #optional - gap3
Group( (1,5), (2,5), (3,5), (4,5) )
sage: type(S5)                 #optional - gap3
<class 'sage.interfaces.gap3.GAP3Record'>
```

This provides a Pythonic interface to GAP3. If `gap_function` is the name of a GAP3 function, then the syntax `gap_element.gap_function()` returns the `gap_element` obtained by evaluating the command `gap_function(gap_element)` in GAP3:

```
sage: S5.Size()                #optional - gap3
120
sage: S5.CharTable()           #optional - gap3
CharTable( Group( (1,5), (2,5), (3,5), (4,5) ) )
```

Alternatively, you can instead use the syntax `gap3.gap_function(gap_element)`:

```
sage: gap3.DerivedSeries(S5)   #optional - gap3
[ Group( (1,5), (2,5), (3,5), (4,5) ),
  Subgroup( Group( (1,5), (2,5), (3,5), (4,5) ),
    [ (1,2,5), (1,3,5), (1,4,5) ] ) ]
```

If `gap_element` corresponds to a GAP3 record, then `gap_element.recfield` provides a means to access the record element corresponding to the field `recfield`:

```
sage: S5.IsRec()               #optional - gap3
true
sage: S5.recfields()           #optional - gap3
['isDomain', 'isGroup', 'identity', 'generators', 'operations',
'isPermGroup', 'isFinite', '1', '2', '3', '4', 'degree']
sage: S5.identity              #optional - gap3
()
sage: S5.degree                #optional - gap3
5
sage: S5.1                     #optional - gap3
(1,5)
sage: S5.2                     #optional - gap3
(2,5)
```

2. By typing `%gap3` or `gap3.interact()` at the command-line, you can interact directly with the underlying GAP3 session.

```
sage: gap3.interact()          #not tested

--> Switching to Gap3 <--
```

3. You can start a new GAP3 session as follows:

The interface also has access to the GAP3 help system:

## 10.4 Common Pitfalls

1. If you want to pass a string to GAP3, then you need to wrap it in single quotes as follows:

This is particularly important when a GAP3 package is loaded via the `RequirePackage` method (note that one can instead use the `load_package` method):

## 10.5 Examples

## 10.4. Common Pitfalls

```
sage: gap3.load_package("chevie")           #optional - gap3 chevie
sage: gap3.version() # random               #optional - gap3
'lib: v3r4p4 1997/04/18, src: v3r4p0 1994/07/10, sys: usg gcc ansi'
```

Working with GAP3 lists. Note that GAP3 lists are 1-indexed:

```
sage: L = gap3([1,2,3])                   #optional - gap3
sage: L[1]                                #optional - gap3
1
sage: L[2]                                #optional - gap3
2
sage: 3 in L                               #optional - gap3
True
sage: 4 in L                               #optional - gap3
False
sage: m = gap3([[1,2],[3,4]])             #optional - gap3
sage: m[2,1]                              #optional - gap3
3
sage: [1,2] in m                           #optional - gap3
True
sage: [3,2] in m                           #optional - gap3
False
sage: gap3([1,2]) in m                     #optional - gap3
True
```

Controlling variable names used by GAP3:

```
sage: gap3('2', name='x')                 #optional - gap3
2
sage: gap3('x')                           #optional - gap3
2
sage: gap3.unbind('x')                     #optional - gap3
sage: gap3('x')                           #optional - gap3
Traceback (most recent call last):
...
TypeError: Gap3 produced error output
Error, Variable: 'x' must have a value
...
```

```
class sage.interfaces.gap3.GAP3Element (parent, value, is_name=False, name=None)
```

Bases: `sage.interfaces.gap3.GapElement_generic`

A GAP3 element

---

**Note:** If the corresponding GAP3 element is a GAP3 record, then the class is changed to a GAP3Record.

---

INPUT:

- parent – the GAP3 session
- value – the GAP3 command as a string
- is\_name – bool (default: False); if True, then value is the variable name for the object
- name – str (default: None); the variable name to use for the object. If None, then a variable name is generated.

---

**Note:** If you pass E, X or Z for name, then an error is raised because these are sacred variable names in GAP3 that should never be redefined. Sage raises an error because GAP3 does not!

## EXAMPLES:

```

sage: from sage.interfaces.gap3 import GAP3Element      #optional - gap3
sage: gap3 = Gap3()                                    #optional - gap3
sage: GAP3Element(gap3, value='3+2')                    #optional - gap3
5
sage: GAP3Element(gap3, value='sage0', is_name=True)    #optional - gap3
5

```

## TESTS:

```

sage: GAP3Element(gap3, value='3+2', is_name=False, name='X') #optional - gap3
Traceback (most recent call last):
...

```

**ValueError:** you are attempting to redefine X; but you should never redefine E, X or Z in gap3 (b

## AUTHORS:

- Franco Saliola (Feb 2010)

**class** `sage.interfaces.gap3.GAP3Record` (*parent, value, is\_name=False, name=None*)

Bases: `sage.interfaces.gap3.GAP3Element`

A GAP3 record

---

**Note:** This class should not be called directly, use `GAP3Element` instead. If the corresponding GAP3 element is a GAP3 record, then the class is changed to a `GAP3Record`.

---

## AUTHORS:

- Franco Saliola (Feb 2010)

**operations()**

Return a list of the GAP3 operations for the record.

## OUTPUT:

- list of strings - operations of the record

## EXAMPLES:

```

sage: S5 = gap3.SymmetricGroup(5)                        #optional - gap3
sage: S5.operations()                                    #optional - gap3
[ ..., 'NormalClosure', 'NormalIntersection', 'Normalizer',
'NumberConjugacyClasses', 'PCore', 'Radical', 'SylowSubgroup',
'TrivialSubgroup', 'FusionConjugacyClasses', 'DerivedSeries', ...]
sage: S5.DerivedSeries()                                #optional - gap3
[ Group( (1,5), (2,5), (3,5), (4,5) ),
  Subgroup( Group( (1,5), (2,5), (3,5), (4,5) ),
    [ (1,2,5), (1,3,5), (1,4,5) ] ) ]

```

**recfields()**

Return a list of the fields for the record. (Record fields are akin to object attributes in Sage.)

## OUTPUT:

- list of strings - the field records

## EXAMPLES:

```

sage: S5 = gap3.SymmetricGroup(5)                        #optional - gap3
sage: S5.recfields()                                     #optional - gap3

```

```

['isDomain', 'isGroup', 'identity', 'generators',
 'operations', 'isPermGroup', 'isFinite', '1', '2',
 '3', '4', 'degree']
sage: S5.degree                                     #optional - gap3
5

```

**trait\_names()**

Defines the list of methods and attributes that will appear for tab completion.

OUTPUT:

- list of strings – the available fields and operations of the record

EXAMPLES:

```

sage: S5 = gap3.SymmetricGroup(5)                 #optional - gap3
sage: S5.trait_names()                             #optional - gap3
[... 'ConjugacyClassesTry', 'ConjugateSubgroup', 'ConjugateSubgroups',
 'Core', 'DegreeOperation', 'DerivedSeries', 'DerivedSubgroup',
 'Difference', 'DimensionsLoewyFactors', 'DirectProduct', ...]

```

**class** sage.interfaces.gap3.**Gap3**(command='gap3')

Bases: sage.interfaces.gap.Gap\_generic

A simple Expect interface to GAP3.

EXAMPLES:

```

sage: from sage.interfaces.gap3 import Gap3
sage: gap3 = Gap3(command='gap3')

```

TESTS:

```

sage: gap3(2) == gap3(3)                         #optional - gap3
False
sage: gap3(2) == gap3(2)                         #optional - gap3
True
sage: gap3.trait_names()                         #optional - gap3
[]

```

We test the interface behaves correctly after a keyboard interrupt:

```

sage: gap3(2)                                     #optional - gap3
2
sage: try:
.....:     gap3._keyboard_interrupt()
.....: except KeyboardInterrupt:
.....:     pass
sage: gap3(2)                                     #optional - gap3
2

```

We test that the interface busts out of GAP3's break loop correctly:

```

sage: f = gap3('function(L) return L[0]; end;;') #optional - gap3
sage: f([1,2,3])                                #optional - gap3
Traceback (most recent call last):
...
RuntimeError: Gap3 produced error output
Error, List Element: <position> must be a positive integer at
return L[0] ...

```

AUTHORS:



- Franco Saliola (Feb 2010)

```
console ()
```

Spawn a new GAP3 command-line session.

EXAMPLES:

```
sage: gap3.console()                                #not tested
```

```
#####  
###   ####  
##      ##  
##         #  
##          #  
####        ##  
#####       ##  
#####      ##  
#####     ##  
#####    ##  
#####   ##  
#####  ##  
#####  
  
Lehrstuhl D fuer Mathematik  
RWTH Aachen  
  
#####  
#           ##  
#           ##  
##          #  
##          #  
#####  
  
#####  
##  #          ##  
#            ##  
#            ##  
##           ##  
##           ##  
#####  
#  
#  
#  
#  
#  
  
Version 3  
Release 4.4  
18 Apr 97  
  
##  # Alice Niemeyer, Werner Nickel, Martin Schoenert  
##  # Johannes Meier, Alex Wegner, Thomas Bishops  
##  # Frank Celler, Juergen Mnich, Udo Polis  
### ## Thomas Breuer, Goetz Pfeiffer, Hans U. Besche  
##### Volkmar Felsch, Heiko Theissen, Alexander Hulpke  
Ansgar Kaup, Akos Seress, Erzsebet Horvath  
Bettina Eick  
For help enter: ?<return>
```

```
gap>
```

```
cputime ( $t=None$ )
```

Returns the amount of CPU time that the GAP session has used in seconds. If `τ` is not `None`, then it returns the difference between the current CPU time and `τ`.

EXAMPLES:

```
sage: t = gap3.cputime() #optional - gap3
sage: t #random #optional - gap3
0.02
sage: gap3.SymmetricGroup(5).Size() #optional - gap3
120
sage: gap3.cputime() #random #optional - gap3
0.14999999999999999
sage: gap3.cputime(t) #random #optional - gap3
0.13
```

### `help (topic, pager=True)`

Print help on the given topic.

INPUT:

- topic – string

EXAMPLES:

```
sage: gap3.help('help', pager=False) #optional - gap3
Help _____...
```

This section describes together with the following section the help system. The help system lets you read the manual interactively.

```
sage: gap3.help('SymmetricGroup', pager=False) #optional - gap3
no section with this name was found
```

#### TESTS:

```
sage: m = gap3([[1,2,3],[4,5,6]]); m #optional - gap3
[ [ 1, 2, 3 ], [ 4, 5, 6 ] ]
sage: gap3.help('help', pager=False) #optional - gap3
Help _____...
sage: m #optional - gap3
[ [ 1, 2, 3 ], [ 4, 5, 6 ] ]
sage: m.Print() #optional - gap3
[ [ 1, 2, 3 ], [ 4, 5, 6 ] ]
sage: gap3.help('Group', pager=False) #optional - gap3
Group _____...
sage: m #optional - gap3
[ [ 1, 2, 3 ], [ 4, 5, 6 ] ]
sage: m.Print() #optional - gap3
[ [ 1, 2, 3 ], [ 4, 5, 6 ] ]
```

```
sage.interfaces.gap3.gap3_console()
Spawn a new GAP3 command-line session.
```

#### EXAMPLES:

```
sage: gap3.console() #not tested
```

```
#####
###      ###
##        ##
##        #
##        #
##        #
####      #
#####    ##
#####    #
#
##
###
## #
## #
# # Alice Niemeyer, Werner Nickel, Martin Schoenert
# # Johannes Meier, Alex Wegner, Thomas Bischops
## # Frank Celler, Juergen Mnich, Udo Polis
### ## Thomas Breuer, Goetz Pfeiffer, Hans U. Besche
##### Volkmar Felsch, Heiko Theissen, Alexander Hulpke
Ansgar Kaup, Akos Seress, Erzsebet Horvath
Bettina Eick
For help enter: ?<return>

gap>
```

```
sage.interfaces.gap3.gap3_version()
Return the version of GAP3 that you have in your PATH on your computer.
```

#### EXAMPLES:

```
sage: gap3_version()                                     # random, optional - gap3
'lib: v3r4p4 1997/04/18, src: v3r4p0 1994/07/10, sys: usg gcc ansi'
```



## INTERFACE TO GROEBNER FAN

AUTHOR:

- Anders Nedergaard Jensen: Write gfan C++ program, which implements algorithms many of which were invented by Jensen, Komei Fukuda, and Rekha Thomas.
- William Stein (2006-03-18): wrote gfan interface (first version)
- Marshall Hampton (2008-03-17): modified to use gfan-0.3, subprocess instead of os.popen2

TODO – much functionality of gfan-0.3 is still not exposed:

```
* at most 52 variables:

    - use gfan_substitute to make easier (?)
    MH: I think this is now irrelevant since gfan can accept the original ring variables

* --symmetry is really useful
    - permutations are 0-based *not* cycle notation; a <---> 0
    output is broken up much more nicely.

* -- can work in  $\mathbb{Z}/p\mathbb{Z}$  for  $p \leq 32749$ 

* -- can compute individual GB's for lex and revlex (via buchberger)
```

**class** sage.interfaces.gfan.**Gfan**

Interface to Anders Jensen's Groebner Fan program.



## INTERFACE TO GIAC

(adapted by F. Han from William Stein and Gregg Musiker maple’s interface)

You must have the optional Giac interpreter installed and available as the command `giac` in your `PATH` in order to use this interface. You need a `giac` version supporting “`giac –sage`” (roughly after 0.9.1). In this case you do not have to install any optional Sage packages. If `giac` is not already installed, you can download binaries or sources or spkg (follow the sources link) from the homepage:

Homepage <<http://www-fourier.ujf-grenoble.fr/~parisse/giac.html>>

Type `giac.[tab]` for a list of all the functions available from your Giac install. Type `giac.[tab]?` for Giac’s help about a given function. Type `giac(...)` to create a new Giac object, and `giac.eval(...)` to run a string using Giac (and get the result back as a string).

If the `giac` spkg is installed, you should find the full html documentation there:

`$SAGE_LOCAL/share/giac/doc/en/cascmd_local/index.html`

### EXAMPLES:

```
sage: giac('3 * 5')                                # optional - giac
15
sage: giac.eval('ifactor(2005)')                    # optional - giac
'5*401'
sage: giac.ifactor(2005)                             # optional - giac
2005
sage: l=giac.ifactors(2005) ; l; l[2]                # optional - giac
[5,1,401,1]
401
sage: giac.fsolve('x^2=cos(x)+4', 'x', '0..5')      # optional - giac
[1.9140206190...]
sage: giac.factor('x^5 - y^5')                      # optional - giac
(x-y)*(x^4+x^3*y+x^2*y^2+x*y^3+y^4)
sage: R.<x,y>=QQ[]; f=(x+y)^5; f2=giac(f); (f-f2).normal() #optional - giac
0
sage: x,y=giac('x,y'); giac.int(y/(cos(2*x)+cos(x)),x) #random; optional - giac
y*2*((-tan(x/2))/6+(-2*1/6/sqrt(3))*ln(abs(6*tan(x/2)-2*sqrt(3))/abs(6*tan(x/2)+2*sqrt(3))))
```

If the string “error” (case insensitive) occurs in the output of anything from Giac, a `RuntimeError` exception is raised.

## 12.1 Tutorial

### AUTHORS:

- Gregg Musiker (2006-02-02): initial version.

(adapted to giac by F.Han)

This tutorial is based on the Maple Tutorial for number theory from <http://www.math.mun.ca/~drideout/m3370/numtheory.html>.

There are several ways to use the Giac Interface in Sage. We will discuss two of those ways in this tutorial.

1. If you have a giac expression such as

```
factor( (x^5-1));
```

We can write that in sage as

```
sage: giac('factor(x^5-1)') # optional - giac
(x-1)*(x^4+x^3+x^2+x+1)
```

Notice, there is no need to use a semicolon.

2. Since Sage is written in Python, we can also import giac commands and write our scripts in a pythonic way. For example, `factor()` is a giac command, so we can also factor in Sage using

```
sage: giac(' (x^5-1) ').factor() # optional - giac
(x-1)*(x^4+x^3+x^2+x+1)
```

where `expression.command()` means the same thing as `command(expression)` in Giac. We will use this second type of syntax whenever possible, resorting to the first when needed.

```
sage: giac(' (x^12-1)/(x-1) ').normal() # optional - giac
x^11+x^10+x^9+x^8+x^7+x^6+x^5+x^4+x^3+x^2+x+1
```

The normal command will reduce a rational function to the lowest terms. In giac, simplify is slower than normal because it tries more sophisticated simplifications (ex algebraic extensions) The factor command will factor a polynomial with rational coefficients into irreducible factors over the ring of integers (if your default configuration of giac (cf .xcasrc) has not allowed square roots). So for example,

```
sage: giac(' (x^12-1) ').factor() # optional - giac
(x-1)*(x+1)*(x^2+1)*(x^2-x+1)*(x^2+x+1)*(x^4-x^2+1)
```

```
sage: giac(' (x^28-1) ').factor() # optional - giac
(x-1)*(x+1)*(x^2+1)*(x^6-x^5+x^4-x^3+x^2-x+1)*(x^6+x^5+x^4+x^3+x^2+x+1)*(x^12-x^10+x^8-x^6+x^4-x^2+1)
```

Another important feature of giac is its online help. We can access this through sage as well. After reading the description of the command, you can press q to immediately get back to your original prompt.

Incidentally you can always get into a giac console by the command.

```
sage: giac.console() # not tested
sage: !giac # not tested
```

Note that the above two commands are slightly different, and the first is preferred.

For example, for help on the giac command factors, we type

```
sage: giac.help('factors') # not tested
```

```
sage: alpha = giac((1+sqrt(5))/2) # optional - giac
sage: beta = giac(1-sqrt(5))/2 # optional - giac
sage: f19 = alpha^19 - beta^19/sqrt(5) # optional - giac
sage: f19 # optional - giac
(sqrt(5)/2+1/2)^19-((-sqrt(5)+1)/2)^19/sqrt(5)
```



```
sage: (f19-(5778*sqrt(5)+33825)/5).normal() # optional - giac
0
```

Let's say we want to write a giac program now that squares a number if it is positive and cubes it if it is negative. In giac, that would look like

```
mysqcu := proc(x)
if x > 0 then x^2;
else x^3; fi;
end;
```

In Sage, we write

```
sage: mysqcu = giac('proc(x) if x > 0 then x^2 else x^3 fi end') # optional - giac
sage: mysqcu(5) # optional - giac
25
sage: mysqcu(-5) # optional - giac
-125
```

More complicated programs should be put in a separate file and loaded.

```
class sage.interfaces.giac.Giac(maxread=10000, script_subdirectory=None, server=None,
server_tmpdir=None, logfile=None)
Bases: sage.interfaces.expect.Expect
```

Interface to the Giac interpreter.

You must have the optional Giac interpreter installed and available as the command `giac` in your `PATH` in order to use this interface. Try the command: `print giac._install_hints()` for more informations on giac installation.

Type `giac.[tab]` for a list of all the functions available from your Giac install. Type `giac.[tab]?` for Giac's help about a given function. Type `giac(...)` to create a new Giac object.

Full html documentation for giac is available from your giac installation at `$PREFIX/share/giac/doc/en/cascmd_en/index.html`

EXAMPLES:

Any Giac instruction can be evaluated as a string by the `giac` command. You can access the `giac` functions by adding the `giac.` prefix to the usual Giac name.

```
sage: l=giac('normal((y+sqrt(2))^4)'); l # optional - giac
y^4+4*sqrt(2)*y^3+12*y^2+8*sqrt(2)*y+4
sage: f=giac('(u,v)->{ if (u<v){ [u,v] } else { [v,u] }}'); f(1,2), f(3,1) # optional - giac
([1,2], [1,3])
```

The output of the `giac` command is a Giac object, and it can be used for another `giac` command.

```
sage: l.factors() #optional - giac
[y+sqrt(2), 4]
sage: giac('(x^12-1)').factor() # optional - giac
(x-1)*(x+1)*(x^2+1)*(x^2-x+1)*(x^2+x+1)*(x^4-x^2+1)
sage: giac('assume(y>0)'); giac('y^2=3').solve('y') #optional - giac
y
...[sqrt(3)]
```

You can create some Giac elements and avoid many quotes like this:

```
sage: x,y,z=giac('x,y,z'); type(y) # optional - giac
<class 'sage.interfaces.giac.GiacElement'>
sage: I1=(1/(cos(2*y)+cos(y))).integral(y,0,pi/4).simplify() #optional - giac
sage: (I1-((-2*ln((sqrt(3)-3*tan(1/8*pi))/(sqrt(3)+3*tan(1/8*pi)))*sqrt(3)-3*tan(1/8*pi))/9)).no
```

```
0
sage: ((y+z*sqrt(5))*(y-sqrt(5)*z)).normal() # optional - giac
y^2-5*z^2
```

Polynomials or elements of SR can be evaluated directly by the giac interface.

```
sage: R.<a,b>=QQ[];f=(2+a+b);p=giac.gcd(f^3+5*f^5,f^2+f^5);p;R(p); #optional - giac
a^2+2*a*b+4*a+b^2+4*b+4
a^2 + 2*a*b + b^2 + 4*a + 4*b + 4
```

Variable names in python and giac are independant.

```
sage: a=sqrt(2);giac('Digits=30;a:=5');a,giac('a'),giac(a),giac(a).evalf() # optional - giac
30
(sqrt(2), 5, sqrt(2), 1.41421356237309504880168872421)
```

**clear**(var)

Clear the variable named var.

EXAMPLES:

```
sage: giac.set('xx', '2') # optional - giac
sage: giac.get('xx') # optional - giac
'2'
sage: giac.clear('xx') # optional - giac
sage: giac.get('xx') # optional - giac
'xx'
```

**completions**(s)

Return all commands that complete the command starting with the string s.

EXAMPLES:

```
sage: c = giac.completions('cas') # optional - giac
sage: 'cas_setup' in c # optional - giac
True
```

**console**()

Spawn a new Giac command-line session.

EXAMPLES:

```
sage: giac_console() # not tested - giac
...
Homepage http://www-fourier.ujf-grenoble.fr/~parisse/giac.html
Released under the GPL license 3.0 or above
See http://www.gnu.org for license details
-----
Press CTRL and D simultaneously to finish session
Type ?commandname for help
0>>
```

**cputime**(t=None)

Returns the amount of CPU time that the Giac session has used. If t is not None, then it returns the difference between the current CPU time and t.

EXAMPLES:

```
sage: t = giac.cputime() # optional - giac
sage: t # random; optional - giac
0.02
sage: x = giac('x') # optional - giac
```

```

sage: giac.diff(x^2, x) # optional - giac
2*x
sage: giac.cputime(t)    # random; optional - giac
0.0

```

**eval** (code, strip=True, \*\*kwds)

Send the code x to the Giac interpreter. Remark: To enable multi-lines codes in the notebook magic mode: %giac, the \n are removed before sending the code to giac.

INPUT:

- code – str
- strip – Default is True and removes \n

EXAMPLES:

```

sage: giac.eval("2+2;\n3") #optional - giac
'4,3'
sage: giac.eval("2+2;\n3",False) # optional - giac
'4\n3'
sage: s='g(x) :={\nx+1;\nx+2;\n}' # optional - giac
sage: giac(s)                    # optional - giac
(x)->{
x+1;
x+2;
}
sage: giac.g(5)                  # optional - giac
7

```

**expect** ()

Returns the pexpect object for this Giac session.

EXAMPLES:

```

sage: m = Giac()                # optional - giac
sage: m.expect() is None        # optional - giac
True
sage: m._start()                # optional - giac
sage: m.expect()                # optional - giac
Giac with PID ... running .../giac --sage
sage: m.quit()                  # optional - giac

```

**get** (var)

Get the value of the variable var.

EXAMPLES:

```

sage: giac.set('xx', '2') # optional - giac
sage: giac.get('xx')      # optional - giac
'2'

```

**help** (str)

Display Giac help about str. This is the same as typing "?str" in the Giac console.

INPUT:

- str - a string to search for in the giac help system

EXAMPLES:

```
sage: giac.help('Psi')           # not tested - depends of giac and $LANG
Psi(a,n)=nth-derivative of the function DiGamma (=ln@Gamma) at point a (Psi(a,0)=Psi(a))...
```

**set** (*var, value*)

Set the variable *var* to the given value.

EXAMPLES:

```
sage: giac.set('xx', '2') # optional - giac
sage: giac.get('xx')      # optional - giac
'2'
```

**trait\_names** (*verbose=True, use\_disk\_cache=True*)

Returns a list of all the commands defined in Giac and optionally (per default) store them to disk.

EXAMPLES:

```
sage: c = giac.trait_names(use_disk_cache=False, verbose=False) # optional - giac
sage: len(c) > 100 # optional - giac
True
sage: 'factors' in c # optional - giac
True
```

**version** ()

Wrapper for *giac*'s *version()*.

EXAMPLES:

```
sage: giac.version() # optional - giac
"giac..."
```

**class** `sage.interfaces.giac.GiacElement` (*parent, value, is\_name=False, name=None*)

Bases: `sage.interfaces.expect.ExpectElement`

**integral** (*var='x', min=None, max=None*)

Return the integral of self with respect to the variable *x*.

INPUT:

- *var* - variable
- *min* - default: None
- *max* - default: None

Returns the definite integral if *xmin* is not None, otherwise returns an indefinite integral.

EXAMPLES:

```
sage: y=giac('y'); f=(sin(2*y)/y).integral(y).simplify(); f # optional - giac
Si(2*y)
sage: f.diff(y).simplify() # optional - giac
sin(2*y)/y
```

```
sage: f = giac('exp(x^2)').integral('x',0,1) ; f # optional - giac
integra...
```

```
sage: f.evalf(10) # optional - giac
1.462651746
```

```
sage: x,y=giac('x'),giac('y'); integrate(cos(x+y),'x=0..pi').simplify() # optional - giac
-2*sin(y)
```

**integrate** (*var='x', min=None, max=None*)

Return the integral of self with respect to the variable *x*.

INPUT:

- var - variable
- min - default: None
- max - default: None

Returns the definite integral if xmin is not None, otherwise returns an indefinite integral.

EXAMPLES:

```
sage: y=giac('y'); f=(sin(2*y)/y).integral(y).simplify(); f      # optional - giac
Si(2*y)
sage: f.diff(y).simplify()                                       # optional - giac
sin(2*y)/y
```

```
sage: f = giac('exp(x^2)').integral('x',0,1) ; f                # optional - giac
integra...
```

```
sage: f.evalf(10)                                                # optional - giac
1.462651746
```

```
sage: x,y=giac('x'),giac('y'); integrate(cos(x+y),'x=0..pi').simplify() # optional - giac
-2*sin(y)
```

**sum**(var, min=None, max=None)

Return the sum of self with respect to the variable x.

INPUT:

- var - variable
- min - default: None
- max - default: None

Returns the definite integral if xmin is not None, otherwise returns an indefinite integral.

EXAMPLES:

```
sage: giac('1/(1+k^2)').sum('k',-oo,+infinity).simplify()      # optional - giac
(pi*exp(pi)^2+pi)/(exp(pi)^2-1)
```

**trait\_names**()

EXAMPLES:

```
sage: a = giac(2) # optional - giac
sage: 'sin' in a.trait_names() # optional - giac
True
```

**unapply**(var)

Creates a Giac function in the given arguments from a Giac symbol.

EXAMPLES:

```
sage: f=giac('y^3+1+t')      # optional - giac
sage: g=(f.unapply('y,t'))   # optional - giac
sage: g                       # optional - giac
(y,t)->y^3+1+t
sage: g(1,2)                  # optional - giac
4
```

**class** sage.interfaces.giac.**GiacFunction**(parent, name)

Bases: sage.interfaces.expect.ExpectFunction

**class** `sage.interfaces.giac.GiacFunctionElement` (*obj, name*)  
Bases: `sage.interfaces.expect.FunctionElement`

`sage.interfaces.giac.giac_console()`  
Spawn a new Giac command-line session.

EXAMPLES:

```
sage: giac.console()  # not tested - giac
...
Homepage http://www-fourier.ujf-grenoble.fr/~parisse/giac.html
Released under the GPL license 3.0 or above
See http://www.gnu.org for license details
-----
Press CTRL and D simultaneously to finish session
Type ?commandname for help
```

`sage.interfaces.giac.reduce_load_Giac()`  
Returns the `giac` object created in `sage.interfaces.giac`.

EXAMPLES:

```
sage: from sage.interfaces.giac import reduce_load_Giac
sage: reduce_load_Giac()
Giac
```

## INTERFACE TO THE GNUPLOT INTERPRETER

**class** `sage.interfaces.gnuplot.Gnuplot`

Bases: `sage.structure.sage_object.SageObject`

Interface to the Gnuplot interpreter.

**console** ()

**gnuplot** ()

**interact** (*cmd*)

**plot** (*cmd*, *file=None*, *verbose=True*, *reset=True*)

Draw the plot described by *cmd*, and possibly also save to an eps or png file.

INPUT:

- *cmd* - string
- *file* - string (default: None), if specified save plot to given file, which may be either an eps (default) or png file.
- *verbose* - print some info
- *reset* - True: reset gnuplot before making graph

OUTPUT: displays graph

---

**Note:** Note that  $\wedge$  s are replaced by  $**$  s before being passed to gnuplot.

---

**plot3d** (*f*, *xmin=-1*, *xmax=1*, *ymin=-1*, *ymax=1*, *zmin=-1*, *zmax=1*, *title=None*, *samples=25*, *isosamples=20*, *xlabel='x'*, *ylabel='y'*, *interact=True*)

**plot3d\_parametric** (*f*=*'cos(u)\*(3 + v\*cos(u/2)), sin(u)\*(3 + v\*cos(u/2)), v\*sin(u/2)'*, *range1='[u=-pi:pi]'*, *range2='[v=-0.2:0.2]'*, *samples=50*, *title=None*, *interact=True*)

Draw a parametric 3d surface and rotate it interactively.

INPUT:

- *f* - (string) a function of two variables, e.g., *'cos(u)\*(3 + v\*cos(u/2)), sin(u)\*(3 + v\*cos(u/2)), v\*sin(u/2)'*
- *range1* - (string) range of values for one variable, e.g., *'[u=-pi:pi]'*
- *range2* - (string) range of values for another variable, e.g., *'[v=-0.2:0.2]'*
- *samples* - (int) number of sample points to use
- *title* - (string) title of the graph.

EXAMPLES:

```
sage: gnuplot.plot3d_parametric('v^2*sin(u), v*cos(u), v*(1-v)') # optional - gnuplot (no
sage.interfaces.gnuplot.gnuplot_console()
```



## INTERFACE TO THE GP CALCULATOR OF PARI/GP

Type `gp . [tab]` for a list of all the functions available from your Gp install. Type `gp . [tab] ?` for Gp's help about a given function. Type `gp ( . . . )` to create a new Gp object, and `gp.eval ( . . . )` to evaluate a string using Gp (and get the result back as a string).

EXAMPLES: We illustrate objects that wrap GP objects (gp is the PARI interpreter):

```
sage: M = gp('[1,2;3,4]')
sage: M
[1, 2; 3, 4]
sage: M * M
[7, 10; 15, 22]
sage: M + M
[2, 4; 6, 8]
sage: M.matdet()
-2

sage: E = gp.ellinit([1,2,3,4,5])
sage: E.ellglobalred()
[10351, [1, -1, 0, -1], 1, [11, 1; 941, 1], [[1, 5, 0, 1], [1, 5, 0, 1]]]
sage: E.ellan(20)
[1, 1, 0, -1, -3, 0, -1, -3, -3, -3, -1, 0, 1, -1, 0, -1, 5, -3, 4, 3]
```

```
sage: primitive_root(7)
3
sage: x = gp("znlog( Mod(2,7), Mod(3,7) )")
sage: 3^x % 7
2
```

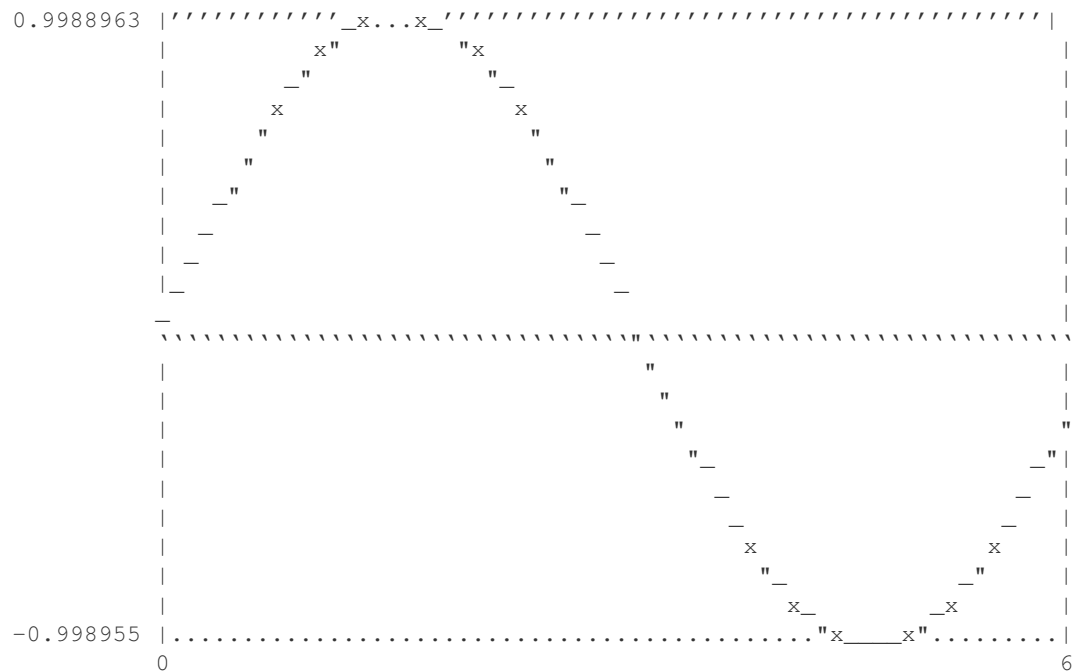
```
sage: print gp("taylor(sin(x),x)")
x - 1/6*x^3 + 1/120*x^5 - 1/5040*x^7 + 1/362880*x^9 - 1/39916800*x^11 + 1/6227020800*x^13 - 1/130767
```

GP has a powerful very efficient algorithm for numerical computation of integrals.

```
sage: gp("a = intnum(x=0,6,sin(x))")
0.03982971334963397945434770208 # 32-bit
0.039829713349633979454347702077075594548 # 64-bit
sage: gp("a")
0.03982971334963397945434770208 # 32-bit
0.039829713349633979454347702077075594548 # 64-bit
sage: gp.kill("a")
sage: gp("a")
a
```

Note that gp ASCII plots *do* work in Sage, as follows:

```
sage: print gp.eval("plot(x=0,6,sin(x))")
```



The GP interface reads in even very long input (using files) in a robust manner, as long as you are creating a new object.

```
sage: t = '%"s"%10^10000 # ten thousand character string.
sage: a = gp.eval(t)
sage: a = gp(t)
```

In Sage, the PARI large Galois groups datafiles should be installed by default:

```
sage: f = gp('x^9 - x - 2')
sage: f.polgalois()
[362880, -1, 34, "S9"]
```

## TESTS:

Test error recovery:

```
sage: x = gp('1/0')
Traceback (most recent call last):
...
TypeError: Error executing code in GP:
CODE:
    sage[...]=1/0;
PARI/GP ERROR:
*** at top-level: sage[...]=1/0
***                               ^__
*** _/_: impossible inverse in gdiv: 0.
```

## AUTHORS:

- William Stein
- David Joyner: some examples

- William Stein (2006-03-01): added tab completion for methods: `gp.[tab]` and `x = gp(blah); x.[tab]`
- William Stein (2006-03-01): updated to work with PARI 2.2.12-beta
- William Stein (2006-05-17): updated to work with PARI 2.2.13-beta

**class** `sage.interfaces.gp.Gp` (*stacksize=10000000, maxread=100000, script\_subdirectory=None, logfile=None, server=None, server\_tmpdir=None, init\_list\_length=1024, seed=None*)

Bases: `sage.interfaces.expect.Expect`

Interface to the PARI gp interpreter.

Type `gp.[tab]` for a list of all the functions available from your Gp install. Type `gp.[tab]?` for Gp's help about a given function. Type `gp(...)` to create a new Gp object, and `gp.eval(...)` to evaluate a string using Gp (and get the result back as a string).

INPUT:

- `stacksize` (int, default 10000000) – the initial PARI stacksize in bytes (default 10MB)
- `maxread` (int, default 100000) – ??
- `script_subdirectory` (string, default None) – name of the subdirectory of `SAGE_EXTCODE/pari` from which to read scripts
- `logfile` (string, default None) – log file for the pexpect interface
- `server` – name of remote server
- `server_tmpdir` – name of temporary directory on remote server
- `init_list_length` (int, default 1024) – length of initial list of local variables.
- `seed` (int, default random) – random number generator seed for pari

EXAMPLES:

```
sage: Gp()
PARI/GP interpreter
```

**console()**

Spawn a new GP command-line session.

EXAMPLES:

```
sage: gp.console() # not tested
GP/PARI CALCULATOR Version 2.4.3 (development svn-12577)
amd64 running linux (x86-64/GMP-4.2.1 kernel) 64-bit version
compiled: Jul 21 2010, gcc-4.6.0 20100705 (experimental) (GCC)
(readline v6.0 enabled, extended help enabled)
```

**cputime** (*t=None*)

cputime for pari - cputime since the pari process was started.

INPUT:

- `t` - (default: None); if not None, then returns time since `t`

**Warning:** If you call `gettime` explicitly, e.g., `gp.eval('gettime')`, you will throw off this clock.

EXAMPLES:

```
sage: gp.cputime()           # random output
0.00800000000000000002
sage: gp.factor('2^157-1')
[852133201, 1; 60726444167, 1; 1654058017289, 1; 2134387368610417, 1]
sage: gp.cputime()           # random output
0.2690000000000000002
```

**get (var)**

Get the value of the GP variable var.

INPUT:

- var (string) – a valid GP variable identifier

EXAMPLES:

```
sage: gp.set('x', '2')
sage: gp.get('x')
'2'
```

**get\_default (var)**

Return the current value of a PARI gp configuration variable.

INPUT:

- var (string) – the name of a PARI gp configuration variable. (See gp.default () for a list.)

OUTPUT:

(string) the value of the variable.

EXAMPLES:

```
sage: gp.get_default('log')
0
sage: gp.get_default('datadir')
'.../local/share/pari'
sage: gp.get_default('seriesprecision')
16
sage: gp.get_default('realprecision')
28          # 32-bit
38          # 64-bit
```

**get\_precision ()**

Return the current PARI precision for real number computations.

EXAMPLES:

```
sage: gp.get_precision()
28          # 32-bit
38          # 64-bit
```

**get\_real\_precision ()**

Return the current PARI precision for real number computations.

EXAMPLES:

```
sage: gp.get_precision()
28          # 32-bit
38          # 64-bit
```

**get\_series\_precision ()**

Return the current PARI power series precision.

EXAMPLES:

```
sage: gp.get_series_precision()
16
```

**help** (*command*)

Returns GP's help for *command*.

EXAMPLES:

```
sage: gp.help('gcd')
'gcd(x,{y}): greatest common divisor of x and y.'
```

**kill** (*var*)

Kill the value of the GP variable *var*.

INPUT:

- *var* (string) – a valid GP variable identifier

EXAMPLES:

```
sage: gp.set('xx', '22')
sage: gp.get('xx')
'22'
sage: gp.kill('xx')
sage: gp.get('xx')
'xx'
```

**new\_with\_bits\_prec** (*s*, *precision=0*)

Creates a GP object from *s* with *precision* bits of precision. GP actually automatically increases this precision to the nearest word (i.e. the next multiple of 32 on a 32-bit machine, or the next multiple of 64 on a 64-bit machine).

EXAMPLES:

```
sage: pi_def = gp(pi); pi_def
3.141592653589793238462643383          # 32-bit
3.1415926535897932384626433832795028842 # 64-bit
sage: pi_def.precision()
28          # 32-bit
38          # 64-bit
sage: pi_150 = gp.new_with_bits_prec(pi, 150)
sage: new_prec = pi_150.precision(); new_prec
48          # 32-bit
57          # 64-bit
sage: old_prec = gp.set_precision(new_prec); old_prec
28          # 32-bit
38          # 64-bit
sage: pi_150
3.14159265358979323846264338327950288419716939938 # 32-bit
3.14159265358979323846264338327950288419716939937510582098 # 64-bit
sage: gp.set_precision(old_prec)
48          # 32-bit
57          # 64-bit
sage: gp.get_precision()
28          # 32-bit
38          # 64-bit
```

**set** (*var*, *value*)

Set the GP variable *var* to the given value.

INPUT:

- var (string) – a valid GP variable identifier
- value – a value for the variable

EXAMPLES:

```
sage: gp.set('x', '2')
sage: gp.get('x')
'2'
```

**set\_default** (var, value)

Set a PARI gp configuration variable, and return the old value.

INPUT:

- var (string) – the name of a PARI gp configuration variable. (See `gp.default()` for a list.)
- value – the value to set the variable to.

EXAMPLES:

```
sage: old_prec = gp.set_default('realprecision', 110)
sage: gp.get_default('realprecision')
115
sage: gp.set_default('realprecision', old_prec)
115
sage: gp.get_default('realprecision')
28          # 32-bit
38          # 64-bit
```

**set\_precision** (prec)

Sets the PARI precision (in decimal digits) for real computations, and returns the old value.

---

**Note:** PARI/GP rounds up precisions to the nearest machine word, so the result of `get_precision()` is not always the same as the last value inputted to `set_precision()`.

---

EXAMPLES:

```
sage: old_prec = gp.set_precision(53); old_prec
28          # 32-bit
38          # 64-bit
sage: gp.get_precision()
57
sage: gp.set_precision(old_prec)
57
sage: gp.get_precision()
28          # 32-bit
38          # 64-bit
```

**set\_real\_precision** (prec)

Sets the PARI precision (in decimal digits) for real computations, and returns the old value.

---

**Note:** PARI/GP rounds up precisions to the nearest machine word, so the result of `get_precision()` is not always the same as the last value inputted to `set_precision()`.

---

EXAMPLES:

```
sage: old_prec = gp.set_precision(53); old_prec
28          # 32-bit
38          # 64-bit
```

```

sage: gp.get_precision()
57
sage: gp.set_precision(old_prec)
57
sage: gp.get_precision()
28          # 32-bit
38          # 64-bit

```

**set\_seed**(*seed=None*)

Sets the seed for gp interpreter. The seed should be an integer.

EXAMPLES:

```

sage: g = Gp()
sage: g.set_seed(1)
1
sage: [g.random() for i in range(5)]
[1546275796, 879788114, 1745191708, 771966234, 1247963869]

```

**set\_series\_precision**(*prec=None*)

Sets the PARI power series precision, and returns the old precision.

EXAMPLES:

```

sage: old_prec = gp.set_series_precision(50); old_prec
16
sage: gp.get_series_precision()
50
sage: gp.set_series_precision(old_prec)
50
sage: gp.get_series_precision()
16

```

**trait\_names**()

EXAMPLES:

```

sage: c = gp.trait_names()
sage: len(c) > 100
True
sage: 'gcd' in c
True

```

**version**()

Returns the version of GP being used.

EXAMPLES:

```

sage: gp.version() # not tested
((2, 4, 3), 'GP/PARI CALCULATOR Version 2.4.3 (development svn-12577)')

```

**class** sage.interfaces.gp.**GpElement**(*parent, value, is\_name=False, name=None*)

Bases: sage.interfaces.expect.ExpectElement

EXAMPLES: This example illustrates dumping and loading GP elements to compressed strings.

```

sage: a = gp(39393)
sage: loads(a.dumps()) == a
True

```

Since dumping and loading uses the string representation of the object, it need not result in an identical object from the point of view of PARI:

```
sage: E = gp('ellinit([1,2,3,4,5])')
sage: loads(dumps(E)) == E
True
sage: x = gp.Pi()/3
sage: loads(dumps(x)) == x
False
sage: x
1.047197551196597746154214461          # 32-bit
1.0471975511965977461542144610931676281 # 64-bit
sage: loads(dumps(x))
1.047197551196597746154214461          # 32-bit
1.0471975511965977461542144610931676281 # 64-bit
```

The two elliptic curves look the same, but internally the floating point numbers are slightly different.

**bool()**

EXAMPLES:

```
sage: gp(2).bool()
True
sage: bool(gp(2))
True
sage: bool(gp(0))
False
```

**is\_string()**

Tell whether this element is a string.

EXAMPLES:

```
sage: gp('"abc"').is_string()
True
sage: gp('[1,2,3]').is_string()
False
```

**trait\_names()**

EXAMPLES:

```
sage: 'gcd' in gp(2).trait_names()
True
```

**class** sage.interfaces.gp.**GpFunction**(parent, name)

Bases: sage.interfaces.expect.ExpectFunction

**class** sage.interfaces.gp.**GpFunctionElement**(obj, name)

Bases: sage.interfaces.expect.FunctionElement

sage.interfaces.gp.**gp\_console()**

Spawn a new GP command-line session.

EXAMPLES:

```
sage: gp.console() # not tested
GP/PARI CALCULATOR Version 2.4.3 (development svn-12577)
amd64 running linux (x86-64/GMP-4.2.1 kernel) 64-bit version
compiled: Jul 21 2010, gcc-4.6.0 20100705 (experimental) (GCC)
(readline v6.0 enabled, extended help enabled)
```

sage.interfaces.gp.**gp\_version()**

EXAMPLES:



```
sage: gp.version() # not tested
((2, 4, 3), 'GP/PARI CALCULATOR Version 2.4.3 (development svn-12577)')
```

`sage.interfaces.gp.is_GpElement(x)`

Returns True if `x` is a `GpElement`.

EXAMPLES:

```
sage: from sage.interfaces.gp import is_GpElement
sage: is_GpElement(gp(2))
True
sage: is_GpElement(2)
False
```

`sage.interfaces.gp.reduce_load_GP()`

Returns the GP interface object defined in `sage.interfaces.gp`.

EXAMPLES:

```
sage: from sage.interfaces.gp import reduce_load_GP
sage: reduce_load_GP()
PARI/GP interpreter
```



## INTERFACE FOR EXTRACTING DATA AND GENERATING IMAGES FROM JMWOL READABLE FILES.

JmolData is a no GUI version of Jmol useful for extracting data from files Jmol reads and for generating image files.

AUTHORS:

- Jonathan Gutow (2012-06-14): complete doctest coverage
- Jonathan Gutow (2012-03-21): initial version

**class** `sage.interfaces.jmoldata.JmolData`  
Bases: `sage.structure.sage_object.SageObject`

---

### Todo

Create an animated image file (GIF) if spin is on and put data extracted from a file into a variable/string/structure to return

---

**export\_image** (*targetfile, datafile, datafile\_cmd='script', image\_type='PNG', figsize=5, \*\*kws*)  
This executes JmolData.jar to make an image file.

INPUT:

- *targetfile* – the full path to the file where the image should be written.
- *datafile* – full path to the data file Jmol can read or text of a script telling Jmol what to read or load. If it is a script and the platform is cygwin, the filenames in the script should be in native windows format.
- *datafile\_cmd* – (default 'script') 'load' or 'script' should be "load" for a data file.
- *image\_type* – (default "PNG") 'PNG' 'JPG' or 'GIF'
- *figsize* – number (default 5) equal to (pixels/side)/100

OUTPUT:

Image file, .png, .gif or .jpg (default .png)

---

**Note:** Examples will generate an error message if a functional Java Virtual Machine (JVM) is not installed on the machine the Sage instance is running on.

**Warning:** Programmers using this module should check that the JVM is available before making calls to avoid the user getting error messages. Check for the JVM using the function `is_jvm_available()`, which returns True if a JVM is available.

EXAMPLES:

Use Jmol to load a pdb file containing some DNA from a web data base and make an image of the DNA.

If you execute this in the notebook, the image will appear in the output cell:

```
sage: from sage.interfaces.jmoldata import JmolData
sage: JData = JmolData()
sage: script = "load =llcd;display DNA;moveto 0.0 { -473 -713 -518 59.94} 100.0 0.0 0.0 {21.
sage: testfile = tmp_filename(ext='DNA.png')
sage: JData.export_image(targetfile=testfile,datafile=script,image_type='PNG') # optional -
sage: print os.path.exists(testfile) # optional -- java internet
True
```

Use Jmol to save an image of a 3-D object created in Sage. This method is used internally by plot3d to generate static images. This example doesn't have correct scaling:

```
sage: from sage.interfaces.jmoldata import JmolData
sage: JData = JmolData()
sage: D=dodecahedron()
sage: from sage.misc.misc import SAGE_TMP
sage: archive_name=os.path.join(SAGE_TMP, "archive.jmol.zip")
sage: D.export_jmol(archive_name) #not scaled properly...need some more steps.
sage: archive_native = archive_name
sage: import sys
sage: if sys.platform == 'cygwin':
....:     from subprocess import check_output, STDOUT
....:     archive_native = check_output(['cygpath', '-w', archive_native],
....:                                   stderr=STDOUT).rstrip()
sage: script = 'set defaultdirectory "{0}"\n script SCRIPT\n'.format(archive_native)
sage: testfile = os.path.join(SAGE_TMP, "testimage.png")
sage: JData.export_image(targetfile=testfile, datafile=script, image_type="PNG") # optional
sage: print os.path.exists(testfile) # optional -- java
True
```

**is\_jvm\_available()**

Returns True if the Java Virtual Machine is available and False if not.

EXAMPLES:

Check that it returns a boolean:

```
sage: from sage.interfaces.jmoldata import JmolData
sage: JData = JmolData()
sage: type(JData.is_jvm_available())
<type 'bool'>
```

## INTERFACE TO KASH

Sage provides an interface to the KASH computer algebra system, which is a *free* (as in beer!) but *closed source* program for algebraic number theory that shares much common code with Magma. To use KASH, you must install the appropriate optional Sage package by typing something like “sage -i kash3-linux-2005.11.22” or “sage -i kash3\_osx-2005.11.22”. For a list of optional packages type “sage -optional”. If you type one of the above commands, the (about 16MB) package will be downloaded automatically (you don’t have to do that).

It is not enough to just have KASH installed on your computer. Note that the KASH Sage package is currently only available for Linux and OSX. If you need Windows, support contact me ([wstein@gmail.com](mailto:wstein@gmail.com)).

The KASH interface offers three pieces of functionality:

1. `kash_console()` - A function that dumps you into an interactive command-line KASH session. Alternatively, type `!kash` from the Sage prompt.
2. `kash(expr)` - Creation of a Sage object that wraps a KASH object. This provides a Pythonic interface to KASH. For example, if `f=kash.new(10)`, then `f.Factors()` returns the prime factorization of 10 computed using KASH.
3. `kash.function_name(args ...)` - Call the indicated KASH function with the given arguments and return the result as a KASH object.
4. `kash.eval(expr)` - Evaluation of arbitrary KASH expressions, with the result returned as a string.

### 16.1 Issues

For some reason hitting Control-C to interrupt a calculation doesn’t work correctly. (TODO)

### 16.2 Tutorial

The examples in this tutorial require that the optional kash package be installed.

#### 16.2.1 Basics

Basic arithmetic is straightforward. First, we obtain the result as a string.

```
sage: kash.eval('(9 - 7) * (5 + 6)')           # optional -- kash
'22'
```

Next we obtain the result as a new KASH object.

```
sage: a = kash('(9 - 7) * (5 + 6)'); a          # optional -- kash
22
sage: a.parent()                             # optional -- kash
Kash
```

We can do arithmetic and call functions on KASH objects:

```
sage: a*a                                     # optional -- kash
484
sage: a.Factorial()                         # optional -- kash
112400072777607680000
```

## 16.2.2 Integrated Help

Use the `kash.help(name)` command to get help about a given command. This returns a list of help for each of the definitions of `name`. Use `print kash.help(name)` to nicely print out all signatures.

## 16.2.3 Arithmetic

Using the `kash.new` command we create Kash objects on which one can do arithmetic.

```
sage: a = kash(12345)                       # optional -- kash
sage: b = kash(25)                          # optional -- kash
sage: a/b                                   # optional -- kash
2469/5
sage: a*b                                   # optional -- kash
1937659030411463935651167391656422626577614411586152317674869233464019922771432158872187137603759765
```

## 16.2.4 Variable assignment

Variable assignment using `kash` takes place in Sage.

```
sage: a = kash('32233')                    # optional -- kash
sage: a                                     # optional -- kash
32233
```

In particular, `a` is not defined as part of the KASH session itself.

```
sage: kash.eval('a')                       # optional -- kash
"Error, the variable 'a' must have a value"
```

Use `a.name()` to get the name of the KASH variable:

```
sage: a.name()                             # somewhat random; optional - kash
'sage0'
sage: kash(a.name())                       # optional -- kash
32233
```

## 16.2.5 Integers and Rationals

We illustrate arithmetic with integers and rationals in KASH.

```
sage: kash.GCD(15,25)
5
sage: kash.LCM(15,25)
75
sage: kash.Div(25,15)
1
sage: kash(17) % kash(5)
2
sage: kash.IsPrime(10007)
TRUE
sage: kash.IsPrime(2005)
FALSE
sage: kash.NextPrime(10007)
10009
```

[illegible]

```
sage: kash.Log('3+4*I') # optional -- kash
1.60943791243410037460075933323 + 0.927295218001612232428512462922*I
sage: kash.Log('I') # optional -- kash
1.57079632679489661923132169164*I

sage: kash.Sqrt(4) # optional -- kash
2.00000000000000000000000000000000
sage: kash.Sqrt(2) # optional -- kash
1.41421356237309504880168872421

sage: kash.Floor('9/5') # optional -- kash
1
sage: kash.Floor('3/5') # optional -- kash
0

sage: x_c = kash('3+I') # optional -- kash
sage: x_c.Argument() # optional -- kash
0.321750554396642193401404614359
sage: x_c.Imaginary() # optional -- kash
1.00000000000000000000000000000000
```

### 16.2.7 Lists

Note that list appends are completely different in KASH than in Python. Use underscore after the function name for the mutation version.

```
sage: v = kash([1,2,3]); v
[ 1, 2, 3 ]
sage: v[1]
1
sage: v[3]
3
sage: v.Append([5])
[ 1, 2, 3, 5 ]
sage: v
[ 1, 2, 3 ]
sage: v.Append_([5, 6])
SUCCESS
sage: v
[ 1, 2, 3, 5, 6 ]
sage: v.Add(5)
[ 1, 2, 3, 5, 6, 5 ]
sage: v
[ 1, 2, 3, 5, 6 ]
sage: v.Add_(5)
SUCCESS
sage: v
[ 1, 2, 3, 5, 6, 5 ]
```

The `Apply` command applies a function to each element of a list.

```
:: sage: L = kash([1,2,3,4]) # optional - kash sage: L.Apply('i -> 3*i') # optional - kash [ 3, 6, 9, 12 ] sage: L #
optional - kash [ 1, 2, 3, 4 ] sage: L.Apply('IsEven') # optional - kash [ FALSE, TRUE, FALSE, TRUE ] sage:
L # optional - kash [ 1, 2, 3, 4 ]
```



## 16.2.8 Ranges

the following are examples of ranges.

```
sage: L = kash(' [1..10]')          # optional -- kash
sage: L                             # optional -- kash
[ 1 .. 10 ]
sage: L = kash(' [2,4..100]')      # optional -- kash
sage: L                             # optional -- kash
[ 2, 4 .. 100 ]
```

## 16.2.9 Sequences

## 16.2.10 Tuples

## 16.2.11 Polynomials

```
sage: f = kash('X^3 + X + 1')      # optional -- kash
sage: f + f                       # optional -- kash
2*X^3 + 2*X + 2
sage: f * f                       # optional -- kash
X^6 + 2*X^4 + 2*X^3 + X^2 + 2*X + 1
sage: f.Evaluate(10)              # optional -- kash
1011
sage: Qx = kash.PolynomialAlgebra('Q') # optional -- kash
sage: Qx.gen(1)**5 + kash('7/3')    # sage1 below somewhat random; optional -- kash
sage1.1^5 + 7/3
```

## 16.2.12 Number Fields

We create an equation order.

```
sage: f = kash('X^5 + 4*X^4 - 56*X^2 -16*X + 192') # optional -- kash
sage: OK = f.EquationOrder()                       # optional -- kash
sage: OK                                            # optional -- kash
Equation Order with defining polynomial X^5 + 4*X^4 - 56*X^2 - 16*X + 192 over Z

sage: f = kash('X^5 + 4*X^4 - 56*X^2 -16*X + 192') # optional -- kash
sage: O = f.EquationOrder()                       # optional -- kash
sage: a = O.gen(2)                                # optional -- kash
sage: a                                            # optional -- kash
[0, 1, 0, 0, 0]
sage: O.Basis()                                   # output somewhat random; optional -- kash
[
  _NG.1,
  _NG.2,
  _NG.3,
  _NG.4,
  _NG.5
]
sage: O.Discriminant()                           # optional -- kash
1364202618880
sage: O.MaximalOrder()                           # name sage2 below somewhat random; optional -- kash
Maximal Order of sage2
```

```
sage: O = kash.MaximalOrder('X^3 - 77')           # optional -- kash
sage: I = O.Ideal(5, [2, 1, 0])                   # optional -- kash
sage: I                                           # name sage14 below random; optional -- kash
Ideal of sage14
Two element generators:
[5, 0, 0]
[2, 1, 0]

sage: F = I.Factorisation()                       # optional -- kash
sage: F                                           # name sage14 random; optional -- kash
[
<Prime Ideal of sage14
Two element generators:
[5, 0, 0]
[2, 1, 0], 1>
]
```

Determining whether an ideal is principal.

```
sage: I.IsPrincipal()                             # optional -- kash
FALSE, extended by:
ext1 := Unassign
```

Computation of class groups and unit groups:

```
sage: f = kash('X^5 + 4*X^4 - 56*X^2 -16*X + 192') # optional -- kash
sage: O = kash.EquationOrder(f)                   # optional -- kash
sage: OK = O.MaximalOrder()                       # optional -- kash
sage: OK.ClassGroup()                             # name sage32 below random; optional -- kash
Abelian Group isomorphic to Z/6
  Defined on 1 generator
  Relations:
  6*sage32.1 = 0, extended by:
  ext1 := Mapping from: grp^abl: sage32 to ids/ord^num: _AA

sage: U = OK.UnitGroup()                           # optional -- kash
sage: U                                             # name sage34 below random; optional -- kash
Abelian Group isomorphic to Z/2 + Z + Z
  Defined on 3 generators
  Relations:
  2*sage34.1 = 0, extended by:
  ext1 := Mapping from: grp^abl: sage34 to ord^num: sage30

sage: kash.Apply('x->%s.ext1(x)'%U.name(), U.Generators().List()) # optional -- kash
[ [1, -1, 0, 0, 0], [1, 1, 0, 0, 0], [-1, 0, 0, 0, 0] ]
```

### 16.2.13 Function Fields

```
sage: k = kash.FiniteField(25)                    # optional -- kash
sage: kT = k.RationalFunctionField()               # optional -- kash
sage: kTy = kT.PolynomialAlgebra()                # optional -- kash
sage: T = kT.gen(1)                               # optional -- kash
sage: y = kTy.gen(1)                              # optional -- kash
sage: f = y**3 + T**4 + 1                         # optional -- kash
```

## 16.3 Long Input

The KASH interface reads in even very long input (using files) in a robust manner, as long as you are creating a new object.

---

**Note:** Using `kash.eval` for long input is much less robust, and is not recommended.

---

```
sage: a = kash(range(10000))                                # optional -- kash
```

Note that KASH seems to not support string or integer literals with more than 1024 digits, which is why the above example uses a list unlike for the other interfaces.

```
class sage.interfaces.kash.Kash(max_workspace_size=None,          maxread=100000,
                                script_subdirectory=None, restart_on_ctrlc=True, logfile=None,
                                server=None, server_tmpdir=None)
```

Bases: `sage.interfaces.expect.Expect`

Interface to the Kash interpreter.

AUTHORS:

- William Stein and David Joyner

**console**()

**eval**(*x*, *newlines=False*, *strip=True*, *\*\*kws*)

Send the code in the string *s* to the Kash interpreter and return the output as a string.

INPUT:

- s* - string containing Kash code.
- newlines* - bool (default: True); if False, remove all backslash-newlines inserted by the Kash output formatter.
- strip* - ignored

**get**(*var*)

Get the value of the variable *var*.

**help**(*name=None*)

Return help on KASH commands.

Returns help on all commands with a given name. If *name* is None, return the location of the installed Kash HTML documentation.

EXAMPLES:

```
sage: X = kash.help('IntegerRing')    # optional -- kash
```

There is one entry in *X* for each item found in the documentation for this function: If you type `print X[0]` you will get help on about the first one, printed nicely to the screen.

AUTHORS:

- Sebastion Pauli (2006-02-04): during Sage coding sprint

**help\_search**(*name*)

**set**(*var*, *value*)

Set the variable *var* to the given value.

**version**()

```
class sage.interfaces.kash.KashDocumentation
    Bases: list

class sage.interfaces.kash.KashElement (parent, value, is_name=False, name=None)
    Bases: sage.interfaces.expect.ExpectElement

sage.interfaces.kash.is_KashElement(x)

sage.interfaces.kash.kash_console()

sage.interfaces.kash.kash_version()

sage.interfaces.kash.reduce_load_Kash()
```

## INTERFACE TO LIE

LiE is a software package under development at CWI since January 1988. Its purpose is to enable mathematicians and physicists to obtain on-line information as well as to interactively perform computations of a Lie group theoretic nature. It focuses on the representation theory of complex semisimple (reductive) Lie groups and algebras, and on the structure of their Weyl groups and root systems.

Type `lie.[tab]` for a list of all the functions available from your LiE install. Type `lie.[tab]?` for LiE's help about a given function. Type `lie(...)` to create a new LiE object, and `lie.eval(...)` to run a string using LiE (and get the result back as a string).

To access the LiE interpreter directly, run `lie_console()`.

EXAMPLES:

```
sage: a4 = lie('A4') # optional - lie
sage: lie.diagram('A4') # optional - lie
O---O---O---O
1   2   3   4
A4

sage: lie.diagram(a4) # optional - lie
O---O---O---O
1   2   3   4
A4

sage: a4.diagram() # optional - lie
O---O---O---O
1   2   3   4
A4

sage: a4.Cartan() # optional - lie
[[ 2,-1, 0, 0]
, [-1, 2,-1, 0]
, [ 0,-1, 2,-1]
, [ 0, 0,-1, 2]
]
sage: lie.LR_tensor([3,1],[2,2]) # optional - lie
1X[5,3]
```

### 17.1 Tutorial

The following examples are taken from Section 2.1 of the LiE manual.

You can perform basic arithmetic operations in LiE.

```
sage: lie.eval('19+68') # optional - lie
'87'
sage: a = lie('111111111*111111111') # optional - lie
sage: a # optional - lie
1234567900987654321
sage: a/111111111 # optional - lie
1111111111
sage: a = lie('345') # optional - lie
sage: a^2+3*a-5 # optional - lie
120055
sage: _ / 7*a # optional - lie
5916750
```

Vectors in LiE are created using square brackets. Notice that the indexing in LiE is 1-based, unlike Python/Sage which is 0-based.

```
sage: v = lie('[3,2,6873,-38]') # optional - lie
sage: v # optional - lie
[3,2,6873,-38]
sage: v[3] # optional - lie
6873
sage: v+v # optional - lie
[6,4,13746,-76]
sage: v*v # optional - lie
47239586
sage: v+234786 # optional - lie
[3,2,6873,-38,234786]
sage: v-3 # optional - lie
[3,2,-38]
sage: v^v # optional - lie
[3,2,6873,-38,3,2,6873,-38]
```

You can also work with matrices in LiE.

```
sage: m = lie('[ [1,0,3,3], [12,4,-4,7], [-1,9,8,0], [3,-5,-2,9] ]') # optional - lie
sage: m # optional - lie
[[ 1, 0, 3, 3]
 , [12, 4, -4, 7]
 , [-1, 9, 8, 0]
 , [ 3, -5, -2, 9]
 ]
sage: print lie.eval('*'+m.__name) # optional - lie
[[1,12,-1, 3]
 , [0, 4, 9,-5]
 , [3,-4, 8,-2]
 , [3, 7, 0, 9]
 ]

sage: m^3 # optional - lie
[[ 220, 87, 81, 375]
 , [-168,-1089, 13,1013]
 , [1550, 357,-55,1593]
 , [-854, -652, 98,-170]
 ]
sage: v*m # optional - lie
[-6960,62055,55061,-319]
sage: m*v # optional - lie
[20508,-27714,54999,-14089]
```

```
sage: v*m*v # optional - lie
378549605
```

```
sage: m+v # optional - lie
[[ 1, 0, 3, 3]
, [12, 4, -4, 7]
, [-1, 9, 8, 0]
, [ 3, -5, -2, 9]
, [ 3, 2, 6873, -38]
]
```

```
sage: m-2 # optional - lie
[[ 1, 0, 3, 3]
, [-1, 9, 8, 0]
, [ 3, -5, -2, 9]
]
```

LiE handles multivariate (Laurent) polynomials.

```
sage: lie('X[1,2]') # optional - lie
1X[1,2]
sage: -3*_ # optional - lie
-3X[1,2]
sage: _ + lie('4X[-1,4]') # optional - lie
4X[-1,4] - 3X[ 1,2]
sage: _^2 # optional - lie
16X[-2,8] - 24X[ 0,6] + 9X[ 2,4]
sage: lie('(4X[-1,4]-3X[1,2])*(X[2,0]-X[0,-4])') # optional - lie
-4X[-1, 0] + 3X[ 1,-2] + 4X[ 1, 4] - 3X[ 3, 2]
sage: _ - _ # optional - lie
0X[0,0]
```

You can call LiE's built-in functions using `lie.functionname`.

```
sage: lie.partitions(6) # optional - lie
[[6,0,0,0,0,0]
, [5,1,0,0,0,0]
, [4,2,0,0,0,0]
, [4,1,1,0,0,0]
, [3,3,0,0,0,0]
, [3,2,1,0,0,0]
, [3,1,1,1,0,0]
, [2,2,2,0,0,0]
, [2,2,1,1,0,0]
, [2,1,1,1,1,0]
, [1,1,1,1,1,1]
]
sage: lie.diagram('E8') # optional - lie
      O 2
      |
      |
O---O---O---O---O---O---O
1   3   4   5   6   7   8
E8
```

You can define your own functions in LiE using `lie.eval`. Once you've defined a function (say `f`), you can call it using `lie.f`; however, user-defined functions do not show up when using tab-completion.

```
sage: lie.eval('f(int x) = 2*x') # optional - lie
''
sage: lie.f(984) # optional - lie
1968
sage: lie.eval('f(int n) = a=3*n-7; if a < 0 then a = -a fi; 7^a+a^3-4*a-57') # optional - lie
''
sage: lie.f(2) # optional - lie
-53
sage: lie.f(5) # optional - lie
5765224
```

LiE's help can be accessed through `lie.help('functionname')` where `functionname` is the function you want to receive help for.

```
sage: print lie.help('diagram') # optional - lie
diagram(g). Prints the Dynkin diagram of g, also indicating
the type of each simple component printed, and labeling the nodes as
done by Bourbaki (for the second and further simple components the
labels are given an offset so as to make them disjoint from earlier
labels). The labeling of the vertices of the Dynkin diagram prescribes
the order of the coordinates of root- and weight vectors used in LiE.
```

This can also be accessed with `lie.functionname?` .

With the exception of groups, all LiE data types can be converted into native Sage data types by calling the `.sage()` method.

Integers:

```
sage: a = lie('1234') # optional - lie
sage: b = a.sage(); b # optional - lie
1234
sage: type(b) # optional - lie
<type 'sage.rings.integer.Integer'>
```

Vectors:

```
sage: a = lie('[1,2,3]') # optional - lie
sage: b = a.sage(); b # optional - lie
[1, 2, 3]
sage: type(b) # optional - lie
<type 'list'>
```

Matrices:

```
sage: a = lie('[ [1,2], [3,4] ]') # optional - lie
sage: b = a.sage(); b # optional - lie
[1 2]
[3 4]
sage: type(b) # optional - lie
<type 'sage.matrix.matrix_integer_dense.Matrix_integer_dense'>
```

Polynomials:

```
sage: a = lie('X[1,2] - 2*X[2,1]') # optional - lie
sage: b = a.sage(); b # optional - lie
-2*x0^2*x1 + x0*x1^2
sage: type(b) # optional - lie
<type 'sage.rings.polynomial.multi_polynomial_libsingular.MPolynomial_libsingular'>
```



Text:

```
sage: a = lie('"text"') # optional - lie
sage: b = a.sage(); b # optional - lie
'text'
sage: type(b) # optional - lie
<type 'str'>
```

LiE can be programmed using the Sage interface as well. Section 5.1.5 of the manual gives an example of a function written in LiE's language which evaluates a polynomial at a point. Below is a (roughly) direct translation of that program into Python / Sage.

```
sage: def eval_pol(p, pt): # optional - lie
...     s = 0
...     for i in range(1,p.length().sage()+1):
...         m = 1
...         for j in range(1,pt.size().sage()+1):
...             m *= pt[j]^p.expon(i)[j]
...         s += p.coef(i)*m
...     return s
sage: a = lie('X[1,2]') # optional - lie
sage: b1 = lie('[1,2]') # optional - lie
sage: b2 = lie('[2,3]') # optional - lie
sage: eval_pol(a, b1) # optional - lie
4
sage: eval_pol(a, b2) # optional - lie
18
```

AUTHORS:

- Mike Hansen 2007-08-27
- William Stein (template)

```
class sage.interfaces.lie.LiE(maxread=100000, script_subdirectory=None, logfile=None,
                             server=None)
Bases: sage.interfaces.expect.Expect
```

Interface to the LiE interpreter.

Type `lie.[tab]` for a list of all the functions available from your LiE install. Type `lie.[tab]?` for LiE's help about a given function. Type `lie(...)` to create a new LiE object, and `lie.eval(...)` to run a string using LiE (and get the result back as a string).

**console()**

Spawn a new LiE command-line session.

EXAMPLES:

```
sage: lie.console() # not tested
LiE version 2.2.2 created on Sep 26 2007 at 18:13:19
Authors: Arjeh M. Cohen, Marc van Leeuwen, Bert Lisser.
Free source code distribution
...
```

**eval (code, strip=True, \*\*kws)**

EXAMPLES:

```
sage: lie.eval('2+2') # optional - lie
'4'
```

**function\_call** (*function*, *args=None*, *kws=None*)

EXAMPLES:

```
sage: lie.function_call("diagram", args=['A4']) # optional - lie
O---O---O---O
1   2   3   4
A4
```

**get** (*var*)

Get the value of the variable *var*.

EXAMPLES:

```
sage: lie.set('x', '2') # optional - lie
sage: lie.get('x')      # optional - lie
'2'
```

**get\_using\_file** (*var*)

EXAMPLES:

```
sage: lie.get_using_file('x')
Traceback (most recent call last):
...
NotImplementedError
```

**help** (*command*)

Returns a string of the LiE help for *command*.

EXAMPLES:

```
sage: lie.help('diagram') # optional - lie
'diagram(g) ...'
```

**read** (*filename*)

EXAMPLES:

```
sage: filename = tmp_filename()
sage: f = open(filename, 'w')
sage: f.write('x = 2\n')
sage: f.close()
sage: lie.read(filename) # optional - lie
sage: lie.get('x')       # optional - lie
'2'
sage: import os
sage: os.unlink(filename)
```

**set** (*var*, *value*)

Set the variable *var* to the given value.

EXAMPLES:

```
sage: lie.set('x', '2') # optional - lie
sage: lie.get('x')      # optional - lie
'2'
```

**trait\_names** (*type=None*, *verbose=False*, *use\_disk\_cache=True*)

EXAMPLES:

```
sage: lie.trait_names() # optional - lie
['Adams',
...
'Cartan_type',
```

```

...
'cent_roots',
...
'n_comp',
...
'write']

```

**version()**

EXAMPLES:

```

sage: lie.version() # optional - lie
'2.1'

```

**class** sage.interfaces.lie.**LiEElement** (parent, value, is\_name=False, name=None)

Bases: sage.interfaces.expect.ExpectElement

**trait\_names()**

Returns the possible tab completions for self.

EXAMPLES:

```

sage: a4 = lie('A4') # optional - lie
sage: a4.trait_names() # optional - lie
['Cartan',
...
'center',
'det_Cartan',
'diagram',
...
'n_comp',
...
'res_mat']

```

**type()**

EXAMPLES:

```

sage: m = lie('[[1,0,3,3],[12,4,-4,7],[-1,9,8,0],[3,-5,-2,9]]') # optional - lie
sage: m.type() # optional - lie
'mat'

```

**class** sage.interfaces.lie.**LiEFunction** (parent, name)

Bases: sage.interfaces.expect.ExpectFunction

**class** sage.interfaces.lie.**LiEFunctionElement** (obj, name)

Bases: sage.interfaces.expect.FunctionElement

sage.interfaces.lie.**is\_LiEElement** (x)

EXAMPLES:

```

sage: from sage.interfaces.lie import is_LiEElement
sage: l = lie(2) # optional - lie
sage: is_LiEElement(l) # optional - lie
True
sage: is_LiEElement(2)
False

```

sage.interfaces.lie.**lie\_console** ()

Spawn a new LiE command-line session.

EXAMPLES:

```
sage: from sage.interfaces.lie import lie_console
sage: lie_console() # not tested
LiE version 2.2.2 created on Sep 26 2007 at 18:13:19
Authors: Arjeh M. Cohen, Marc van Leeuwen, Bert Lisser.
Free source code distribution
...
```

sage.interfaces.lie.lie\_version()

EXAMPLES:

```
sage: from sage.interfaces.lie import lie_version
sage: lie_version() # optional - lie
'2.1'
```

sage.interfaces.lie.reduce\_load\_lie()

EXAMPLES:

```
sage: from sage.interfaces.lie import reduce_load_lie
sage: reduce_load_lie()
LiE Interpreter
```

## LISP INTERFACE

### EXAMPLES:

```
sage: lisp.eval('(* 4 5)')
'20'
sage: a = lisp(3); b = lisp(5)
sage: a + b
8
sage: a * b
15
sage: a / b
3/5
sage: a - b
-2
sage: a.sin()
0.14112
sage: b.cos()
0.2836622
sage: a.exp()
20.085537
sage: lisp.eval('( + %s %s)'%(a.name(), b.name()))
'8'
```

One can define functions and the interface supports object-oriented notation for calling them:

```
sage: lisp.eval('(defun factorial (n) (if (= n 1) 1 (* n (factorial (- n 1)))))')
'FACTORIAL'
sage: lisp('(factorial 10)')
3628800
sage: lisp(10).factorial()
3628800
sage: a = lisp(17)
sage: a.factorial()
355687428096000
```

**AUTHORS:** – William Stein (first version) – William Stein (2007-06-20): significant improvements.

```
class sage.interfaces.lisp.Lisp(maxread=100000, script_subdirectory=None, logfile=None,
                                server=None, server_tmpdir=None)
    Bases: sage.interfaces.expect.Expect
```

### EXAMPLES:

```
sage: lisp == loads(dumps(lisp))
True
```

**console()**

Spawn a new Lisp command-line session.

## EXAMPLES:

```
sage: lisp.console() #not tested
ECL (Embeddable Common-Lisp) ...
Copyright (C) 1984 Taiichi Yuasa and Masami Hagiya
Copyright (C) 1993 Giuseppe Attardi
Copyright (C) 2000 Juan J. Garcia-Ripoll
ECL is free software, and you are welcome to redistribute it
under certain conditions; see file 'Copyright' for details.
Type :h for Help.  Top level.
...
```

**eval** (*code*, *strip=True*, *\*\*kwds*)

## EXAMPLES:

```
sage: lisp.eval('( + 2 2)')
'4'
```

## TEST:

Verify that it works when input == output:

```
sage: lisp.eval('2')
'2'
```

**function\_call** (*function*, *args=None*, *kwds=None*)

Calls the Lisp function with given args and kwds. For Lisp functions, the kwds are ignored.

## EXAMPLES:

```
sage: lisp.function_call('sin', ['2'])
0.9092974
sage: lisp.sin(2)
0.9092974
```

**get** (*var*)

## EXAMPLES:

```
sage: lisp.set('x', '2')
sage: lisp.get('x')
'2'
```

**help** (*command*)

## EXAMPLES:

```
sage: lisp.help('setq')
Traceback (most recent call last):
...
NotImplementedError
```

**kill** (*var*)

## EXAMPLES:

```
sage: lisp.kill('x')
Traceback (most recent call last):
...
NotImplementedError
```

**set** (*var*, *value*)

Set the variable `var` to the given value.

EXAMPLES:

```
sage: lisp.set('x', '2')
sage: lisp.get('x')
'2'
```

TEST:

It must also be possible to eval the variable by name:

```
sage: lisp.eval('x')
'2'
```

**trait\_names()**

EXAMPLES:

```
sage: lisp.trait_names()
Traceback (most recent call last):
...
NotImplementedError
```

**version()**

Returns the version of Lisp being used.

EXAMPLES:

```
sage: lisp.version()
'Version information is given by lisp.console().'
```

**class** `sage.interfaces.lisp.LispElement` (*parent, value, is\_name=False, name=None*)  
 Bases: `sage.interfaces.expect.ExpectElement`

**bool()**

EXAMPLES:

```
sage: lisp(2).bool()
True
sage: lisp(0).bool()
False
sage: bool(lisp(2))
True
```

**class** `sage.interfaces.lisp.LispFunction` (*parent, name*)  
 Bases: `sage.interfaces.expect.ExpectFunction`

**class** `sage.interfaces.lisp.LispFunctionElement` (*obj, name*)  
 Bases: `sage.interfaces.expect.FunctionElement`

`sage.interfaces.lisp.is_LispElement` (*x*)

EXAMPLES:

```
sage: from sage.interfaces.lisp import is_LispElement
sage: is_LispElement(lisp(2))
True
sage: is_LispElement(2)
False
```

`sage.interfaces.lisp.lisp_console()`  
 Spawn a new Lisp command-line session.

EXAMPLES:

```
sage: lisp.console() #not tested
ECL (Embeddable Common-Lisp) ...
Copyright (C) 1984 Taiichi Yuasa and Masami Hagiya
Copyright (C) 1993 Giuseppe Attardi
Copyright (C) 2000 Juan J. Garcia-Ripoll
ECL is free software, and you are welcome to redistribute it
under certain conditions; see file 'Copyright' for details.
Type :h for Help.  Top level.
...
```

```
sage.interfaces.lisp.reduce_load_Lisp()
```

EXAMPLES:

```
sage: from sage.interfaces.lisp import reduce_load_Lisp
sage: reduce_load_Lisp()
Lisp Interpreter
```



## INTERFACE TO MACAULAY2

---

**Note:** You must have Macaulay2 installed on your computer for this interface to work. Macaulay2 is not included with Sage, but you can obtain it from <http://www.math.uiuc.edu/Macaulay2/>. Note additional optional Sage packages are required.

---

Sage provides an interface to the Macaulay2 computational algebra system. This system provides extensive functionality for commutative algebra. You do not have to install any optional packages.

The Macaulay2 interface offers three pieces of functionality:

- `Macaulay2_console()` – A function that dumps you into an interactive command-line Macaulay2 session.
- `Macaulay2(expr)` – Evaluation of arbitrary Macaulay2 expressions, with the result returned as a string.
- `Macaulay2.new(expr)` – Creation of a Sage object that wraps a Macaulay2 object. This provides a Pythonic interface to Macaulay2. For example, if `f=Macaulay2.new(10)`, then `f.gcd(25)` returns the GCD of 10 and 25 computed using Macaulay2.

EXAMPLES:

```
sage: print macaulay2('3/5 + 7/11') # optional - macaulay2
68
--
55
sage: f = macaulay2('f = i -> i^3') # optional - macaulay2
sage: f                               # optional - macaulay2
f
sage: f(5)                             # optional - macaulay2
125

sage: R = macaulay2('ZZ/5[x,y,z]') # optional - macaulay2
sage: print R                       # optional - macaulay2
ZZ
--[x..z, Degrees => {3:1}, Heft => {1}, MonomialOrder => {MonomialSize => 32}, DegreeRank => 1]
5
                                     {GRevLex => {3:1} }
                                     {Position => Up    }

sage: x = macaulay2('x')             # optional - macaulay2
sage: y = macaulay2('y')             # optional - macaulay2
sage: print (x+y)^5                  # optional - macaulay2
5      5
x  + y
sage: parent((x+y)^5)                 # optional - macaulay2
Macaulay2

sage: R = macaulay2('QQ[x,y,z,w]') # optional - macaulay2
sage: f = macaulay2('x^4 + 2*x*y^3 + x*y^2*w + x*y*z*w + x*y*w^2 + 2*x*z*w^2 + y^4 + y^3*w + 2*y^2*z*w + y^2*w^2 + y*w^3 + w^4')
```

```

sage: print f
      4      3      4      4      2      3      2      2      2      4
x  + 2x*y  + y  + z  + x*y w + y w + x*y*z*w + 2y z*w + x*y*w  + 2x*z*w  + w
sage: g = f * macaulay2('x+y^5')      # optional - macaulay2
sage: print g.factor()                # optional - macaulay2
      4      3      4      4      2      3      2      2      2      4      5
(x  + 2x*y  + y  + z  + x*y w + y w + x*y*z*w + 2y z*w + x*y*w  + 2x*z*w  + w ) (y  + x)

```

## AUTHORS:

- Kiran Kedlaya and David Roe (2006-02-05, during Sage coding sprint)
- William Stein (2006-02-09): inclusion in Sage; prompt uses regexp, calling of Macaulay2 functions via `__call__`.
- William Stein (2006-02-09): fixed bug in reading from file and improved output cleaning.
- Kiran Kedlaya (2006-02-12): added ring and ideal constructors, list delimiters, `is_Macaulay2Element`, `sage_polystring`, `__floordiv__`, `__mod__`, `__iter__`, `__len__`; stripped extra leading space and trailing newline from output.

## TODO:

- get rid of all numbers in output, e.g., in ideal function below.

```

class sage.interfaces.macaulay2.Macaulay2(maxread=10000, script_subdirectory=None, log-
                                         file=None, server=None, server_tmpdir=None)

```

Bases: `sage.interfaces.expect.Expect`

Interface to the Macaulay2 interpreter.

**console()**

Spawn a new M2 command-line session.

EXAMPLES:

```

sage: macaulay2.console()      # not tested
Macaulay 2, version 1.1
with packages: Classic, Core, Elimination, IntegralClosure, LLLBases, Parsing, PrimaryDecomp
...

```

**cputime(t=None)**

EXAMPLES:

```

sage: R = macaulay2("QQ[x,y]")      # optional - macaulay2
sage: x,y = R.gens()                # optional - macaulay2
sage: a = (x+y+1)^20                # optional - macaulay2
sage: macaulay2.cputime()            # optional - macaulay2; random
0.483937000000000001

```

**eval(code, strip=True, \*\*kws)**

Send the code `x` to the Macaulay2 interpreter and return the output as a string suitable for input back into Macaulay2, if possible.

INPUT:

- `code` – str
- `strip` – ignored

EXAMPLES:

```

sage: macaulay2.eval("2+2")      # optional - macaulay2
4

```

**get** (*var*)Get the value of the variable *var*.

EXAMPLES:

```
sage: macaulay2.set("a", "2") # optional - macaulay2
sage: macaulay2.get("a")      # optional - macaulay2
2
```

**help** (*s*)

EXAMPLES:

```
sage: macaulay2.help("load") # optional - macaulay2
load -- read Macaulay2 commands
*****
...
* "input" -- read Macaulay2 commands and echo
* "notify" -- whether to notify the user when a file is loaded
```

**ideal** (*\*gens*)Return the ideal generated by *gens*.

INPUT:

- *gens* – list or tuple of Macaulay2 objects (or objects that can be made into Macaulay2 objects via evaluation)

OUTPUT:

the Macaulay2 ideal generated by the given list of *gens*

EXAMPLES:

```
sage: R2 = macaulay2.ring('QQ', '[x, y]'); R2 # optional - macaulay2
QQ[x..y, Degrees => {2:1}, Heft => {1}, MonomialOrder => {MonomialSize => 16}, DegreeRank =>
{Lex => 2}
{Position => Up}

sage: I = macaulay2.ideal( ('y^2 - x^3', 'x - y') ); I # optional - macaulay2
      3      2
ideal (- x  + y , x - y)
sage: J = I^3; J.gb().gens().transpose() # optional - macaulay2
{-9} | y9-3y8+3y7-y6 |
{-7} | xy6-2xy5+xy4-y7+2y6-y5 |
{-5} | x2y3-x2y2-2xy4+2xy3+y5-y4 |
{-3} | x3-3x2y+3xy2-y3 |
```

**new\_from** (*type, value*)Returns a new Macaulay2Element of type *type* constructed from *value*.

EXAMPLES:

```
sage: l = macaulay2.new_from("MutableList", [1,2,3]) # optional - macaulay2
sage: l # optional - macaulay2
MutableList{...3...}
sage: list(l) # optional - macaulay2
[1, 2, 3]
```

**restart** ()

Restart Macaulay2 interpreter.

TEST:

```
sage: macaulay2.restart() # optional - macaulay2
```

**ring** (*base\_ring*='ZZ', *vars*='[x]', *order*='Lex')

Create a Macaulay2 ring.

INPUT:

- *base\_ring* – base ring (see examples below)
- *vars* – a tuple or string that defines the variable names
- *order* – string – the monomial order (default: 'Lex')

OUTPUT:

- a Macaulay2 ring (with base ring ZZ)

EXAMPLES:

This is a ring in variables named a through d over the finite field of order 7, with graded reverse lex ordering:

```
sage: R1 = macaulay2.ring('ZZ/7', '[a..d]', 'GRevLex'); R1 # optional - macaulay2
ZZ
--[a..d, Degrees => {4:1}, Heft => {1}, MonomialOrder => {MonomialSize => 16}, DegreeRank =>
7                                     {GRevLex => {4:1} }
                                     {Position => Up   }
sage: R1.char() # optional - macaulay2
7
```

This is a polynomial ring over the rational numbers:

```
sage: R2 = macaulay2.ring('QQ', '[x, y]'); R2 # optional - macaulay2
QQ[x..y, Degrees => {2:1}, Heft => {1}, MonomialOrder => {MonomialSize => 16}, DegreeRank =>
                                     {Lex => 2 }
                                     {Position => Up   }
```

**set** (*var*, *value*)

Set the variable *var* to the given value.

EXAMPLES:

```
sage: macaulay2.set("a", "2") # optional - macaulay2
sage: macaulay2.get("a") # optional - macaulay2
2
```

**trait\_names** ()

Return a list of tab completions for Macaulay2.

Returns dynamically built sorted list of commands obtained using Macaulay2 “apropos” command. Return type: list of strings

TESTS:

```
sage: names = macaulay2.trait_names() # optional - macaulay2
sage: 'ring' in names # optional - macaulay2
True
sage: macaulay2.eval("abccabc = 4") # optional - macaulay2
4
sage: names = macaulay2.trait_names() # optional - macaulay2
sage: "abccabc" in names # optional - macaulay2
True
```

**use(R)**

Use the Macaulay2 ring R.

## EXAMPLES:

```

sage: R = macaulay2("QQ[x,y]") # optional - macaulay2
sage: P = macaulay2("ZZ/7[symbol x, symbol y]") # optional - macaulay2
sage: macaulay2("x").cls() # optional - macaulay2
ZZ
--[x..y, Degrees => {2:1}, Heft => {1}, MonomialOrder => {MonomialSize => 32}, DegreeRank =>
7 {GRevLex => {2:1} }
{Position => Up }
sage: macaulay2.use(R) # optional - macaulay2
sage: macaulay2("x").cls() # optional - macaulay2
QQ[x..y, Degrees => {2:1}, Heft => {1}, MonomialOrder => {MonomialSize => 32}, DegreeRank =>
{GRevLex => {2:1} }
{Position => Up }

```

**version()**

Returns the version of Macaulay2.

## EXAMPLES:

```

sage: macaulay2.version() # optional - macaulay2
(1, 3, 1)

```

```

class sage.interfaces.macaulay2.Macaulay2Element (parent, value, is_name=False,
name=None)

```

Bases: `sage.interfaces.expect.ExpectElement`

**cls()**

Since class is a keyword in Python, we have to use `cls` to call Macaulay2's class. In Macaulay2, class corresponds to Sage's notion of parent.

## EXAMPLES:

```

sage: macaulay2(ZZ).cls() # optional - macaulay2
Ring

```

**dot(x)**

## EXAMPLES:

```

sage: d = macaulay2.new("MutableHashTable") # optional - macaulay2
sage: d["k"] = 4 # optional - macaulay2
sage: d.dot("k") # optional - macaulay2
4

```

**external\_string()**

## EXAMPLES:

```

sage: R = macaulay2("QQ[symbol x, symbol y]") # optional - macaulay2
sage: R.external_string() # optional - macaulay2
'QQ[x..y, Degrees => {2:1}, Heft => {1}, MonomialOrder => VerticalList{MonomialSize => 32, G

```

**repr()**

## EXAMPLES:

```

sage: R = macaulay2("QQ[x,y,z]/(x^3-y^3-z^3)") # optional - macaulay2
sage: x = macaulay2('x') # optional - macaulay2
sage: y = macaulay2('y') # optional - macaulay2
sage: print x+y # optional - macaulay2
x + y

```

```

sage: print macaulay2("QQ[x,y,z]") # optional - macaulay2
QQ[x..z, Degrees => {3:1}, Heft => {1}, MonomialOrder => {MonomialSize => 32}, DegreeRank =>
{GRevLex => {3:1} }
{Position => Up }

sage: print macaulay2("QQ[x,y,z]/(x+y+z)") # optional - macaulay2
QQ[x, y, z]
-----
x + y + z

```

**sage\_polystring()**

If this Macaulay2 element is a polynomial, return a string representation of this polynomial that is suitable for evaluation in Python. Thus `*` is used for multiplication and `**` for exponentiation. This function is primarily used internally.

EXAMPLES:

```

sage: R = macaulay2.ring('QQ', '(x,y)') # optional - macaulay2
sage: f = macaulay2('x^3 + 3*y^11 + 5') # optional - macaulay2
sage: print f # optional - macaulay2
      3      11
x  + 3y  + 5
sage: f.sage_polystring() # optional - macaulay2
'x**3+3*y**11+5'

```

**sharp(x)**

EXAMPLES:

```

sage: a = macaulay2([1,2,3]) # optional - macaulay2
sage: a.sharp(0) # optional - macaulay2
1

```

**starstar(x)**

The binary operator `**` in Macaulay2 is usually used for tensor or Cartesian power.

EXAMPLES:

```

sage: a = macaulay2([1,2]).set() # optional - macaulay2
sage: a.starstar(a) # optional - macaulay2
set {(1, 1), (1, 2), (2, 1), (2, 2)}

```

**structure\_sheaf()**

EXAMPLES:

```

sage: S = macaulay2('QQ[a..d]') # optional - macaulay2
sage: R = S/macaulay2('a^3+b^3+c^3+d^3') # optional - macaulay2
sage: X = R.Proj() # optional - macaulay2
sage: print X.structure_sheaf() # optional - macaulay2
OO
sage...

```

**subs(\*args, \*\*kws)**

Note that we have to override the substitute method so that we get the default one from Macaulay2 instead of the one provided by Element.

EXAMPLES:

```

sage: R = macaulay2("QQ[x]") # optional - macaulay2
sage: P = macaulay2("ZZ/7[symbol x]") # optional - macaulay2
sage: x, = R.gens() # optional - macaulay2
sage: a = x^2 + 1 # optional - macaulay2

```

```

sage: a = a.substitute(P)                # optional - macaulay2
sage: a.to_sage().parent()              # optional - macaulay2
Univariate Polynomial Ring in x over Finite Field of size 7

```

### **substitute** (\*args, \*\*kws)

Note that we have to override the substitute method so that we get the default one from Macaulay2 instead of the one provided by Element.

#### EXAMPLES:

```

sage: R = macaulay2("QQ[x]")            # optional - macaulay2
sage: P = macaulay2("ZZ/7[symbol x]")   # optional - macaulay2
sage: x, = R.gens()                     # optional - macaulay2
sage: a = x^2 + 1                       # optional - macaulay2
sage: a = a.substitute(P)               # optional - macaulay2
sage: a.to_sage().parent()              # optional - macaulay2
Univariate Polynomial Ring in x over Finite Field of size 7

```

### **to\_sage** ()

#### EXAMPLES:

```

sage: macaulay2(ZZ).to_sage()           # optional - macaulay2
Integer Ring
sage: macaulay2(QQ).to_sage()           # optional - macaulay2
Rational Field

sage: macaulay2(2).to_sage()             # optional - macaulay2
2
sage: macaulay2(1/2).to_sage()           # optional - macaulay2
1/2
sage: macaulay2(2/1).to_sage()           # optional - macaulay2
2
sage: _.parent()                         # optional - macaulay2
Rational Field
sage: macaulay2([1,2,3]).to_sage()       # optional - macaulay2
[1, 2, 3]

sage: m = matrix([[1,2],[3,4]])
sage: macaulay2(m).to_sage()             # optional - macaulay2
[1 2]
[3 4]

sage: macaulay2(QQ['x,y']).to_sage()     # optional - macaulay2
Multivariate Polynomial Ring in x, y over Rational Field
sage: macaulay2(QQ['x']).to_sage()        # optional - macaulay2
Univariate Polynomial Ring in x over Rational Field
sage: macaulay2(GF(7)['x,y']).to_sage()  # optional - macaulay2
Multivariate Polynomial Ring in x, y over Finite Field of size 7

sage: macaulay2(GF(7)).to_sage()          # optional - macaulay2
Finite Field of size 7
sage: macaulay2(GF(49, 'a')).to_sage()    # optional - macaulay2
Finite Field in a of size 7^2

sage: R.<x,y> = QQ[]
sage: macaulay2(x^2+y^2+1).to_sage()      # optional - macaulay2
x^2 + y^2 + 1

sage: R = macaulay2("QQ[x,y]")           # optional - macaulay2

```

```

sage: I = macaulay2("ideal (x,y)")      # optional - macaulay2
sage: I.to_sage()                       # optional - macaulay2
Ideal (x, y) of Multivariate Polynomial Ring in x, y over Rational Field

sage: X = R/I                          # optional - macaulay2
sage: X.to_sage()                      # optional - macaulay2
Quotient of Multivariate Polynomial Ring in x, y over Rational Field by the ideal (x, y)

sage: R = macaulay2("QQ^2")            # optional - macaulay2
sage: R.to_sage()                     # optional - macaulay2
Vector space of dimension 2 over Rational Field

sage: m = macaulay2("'hello'")         # optional - macaulay2
sage: m.to_sage()                     # optional - macaulay2
'hello'

```

**trait\_names()**

Return a list of tab completions for self.

Returns dynamically built sorted list of commands obtained using Macaulay2 “methods” command. All returned functions can take self as their first argument

Return type: list of strings

TEST:

```

sage: a = macaulay2("QQ[x,y]")          # optional - macaulay2
sage: traits = a.trait_names()          # optional - macaulay2
sage: "generators" in traits            # optional - macaulay2
True

```

**underscore(x)**

EXAMPLES:

```

sage: a = macaulay2([1,2,3])           # optional - macaulay2
sage: a.underscore(0)                  # optional - macaulay2
1

```

**class** sage.interfaces.macaulay2.**Macaulay2Function**(parent, name)

Bases: sage.interfaces.expect.ExpectFunction

sage.interfaces.macaulay2.**is\_Macaulay2Element**(x)

EXAMPLES:

```

sage: from sage.interfaces.macaulay2 import is_Macaulay2Element
sage: is_Macaulay2Element(2)           # optional - macaulay2
False
sage: is_Macaulay2Element(macaulay2(2)) # optional - macaulay2
True

```

sage.interfaces.macaulay2.**macaulay2\_console**()

Spawn a new M2 command-line session.

EXAMPLES:

```

sage: macaulay2_console()              # not tested
Macaulay 2, version 1.1
with packages: Classic, Core, Elimination, IntegralClosure, LLBases, Parsing, PrimaryDecomposit
...

```



`sage.interfaces.macauly2.reduce_load_macauly2()`

Used for reconstructing a copy of the Macaulay2 interpreter from a pickle.

EXAMPLES:

```
sage: from sage.interfaces.macauly2 import reduce_load_macauly2
sage: reduce_load_macauly2()
Macauly2
```

`sage.interfaces.macauly2.remove_output_labels(s)`

Remove output labels of Macaulay2 from a string.

•s: output of Macaulay2

•s: string

Returns: the input string with  $n$  symbols removed from the beginning of each line, where  $n$  is the minimal number of spaces or symbols of Macaulay2 output labels (looking like 'o39 = ') present on every non-empty line.

Return type: string

---

**Note:** If  $s$  consists of several outputs and their labels have different width, it is possible that some strings will have leading spaces (or maybe even pieces of output labels). However, this function will try not cut any messages.

---

EXAMPLES:

```
sage: from sage.interfaces.macauly2 import remove_output_labels
sage: output = 'o1 = QQ [x, y]\n\no1 : PolynomialRing\n'
sage: remove_output_labels(output)
'QQ [x, y]\n\nPolynomialRing\n'
```



## INTERFACE TO MAGMA

Sage provides an interface to the Magma computational algebra system. This system provides extensive functionality for number theory, group theory, combinatorics and algebra.

---

**Note:** You must have Magma installed on your computer for this interface to work. Magma is not free, so it is not included with Sage, but you can obtain it from <http://magma.maths.usyd.edu.au/>.

---

The Magma interface offers three pieces of functionality:

1. `magma_console()` - A function that dumps you into an interactive command-line Magma session.
2. `magma.new(obj)` and alternatively `magma(obj)` - Creation of a Magma object from a Sage object `obj`. This provides a Pythonic interface to Magma. For example, if `f=magma.new(10)`, then `f.Factors()` returns the prime factorization of 10 computed using Magma. If `obj` is a string containing an arbitrary Magma expression, then the expression is evaluated in Magma to create a Magma object. An example is `magma.new('10 div 3')`, which returns Magma integer 3.
3. `magma.eval(expr)` - Evaluation of the Magma expression `expr`, with the result returned as a string.

Type `magma.[tab]` for a list of all functions available from your Magma. Type `magma.Function?` for Magma's help about the Magma Function.

### 20.1 Parameters

Some Magma functions have optional “parameters”, which are arguments that in Magma go after a colon. In Sage, you pass these using named function arguments. For example,

```
sage: E = magma('EllipticCurve([0,1,1,-1,0])')           # optional - magma
sage: E.Rank(Bound = 5)                                # optional - magma
0
```

### 20.2 Multiple Return Values

Some Magma functions return more than one value. You can control how many you get using the `nvals` named parameter to a function call:

```
sage: n = magma(100)                                   # optional - magma
sage: n.IsSquare(nvals = 1)                             # optional - magma
true
sage: n.IsSquare(nvals = 2)                             # optional - magma
(true, 10)
```

```
sage: n = magma(-2006)                                # optional - magma
sage: n.Factorization()                                # optional - magma
[ <2, 1>, <17, 1>, <59, 1> ]
sage: n.Factorization(nvals=2)                          # optional - magma
([ <2, 1>, <17, 1>, <59, 1> ], -1)
```

We verify that an obviously principal ideal is principal:

```
sage: _ = magma.eval('R<x> := PolynomialRing(RationalField())') # optional - magma
sage: O = magma.NumberField('x^2+23').MaximalOrder()           # optional - magma
sage: I = magma('ideal<%s| %s.1>' % (O.name(), O.name()))      # optional - magma
sage: I.IsPrincipal(nvals=2)                                    # optional - magma
(true, [1, 0])
```

## 20.3 Long Input

The Magma interface reads in even very long input (using files) in a robust manner.

```
sage: t = '"%s"' % 10^10000 # ten thousand character string. # optional - magma
sage: a = magma.eval(t)     # optional - magma
sage: a = magma(t)          # optional - magma
```

## 20.4 Garbage Collection

There is a subtle point with the Magma interface, which arises from how garbage collection works. Consider the following session:

First, create a matrix *m* in Sage:

```
sage: m=matrix(ZZ,2,[1,2,3,4]) # optional - magma
```

Then I create a corresponding matrix *A* in Magma:

```
sage: A = magma(m) # optional - magma
```

It is called `_sage_ [...]` in Magma:

```
sage: s = A.name(); s # optional - magma
'_sage_ [...]'
```

It's there:

```
sage: magma.eval(s) # optional - magma
'[1 2]\n[3 4]'
```

Now I delete the reference to that matrix:

```
sage: del A # optional - magma
```

Now `_sage_ [...]` is “zeroed out” in the Magma session:

```
sage: magma.eval(s) # optional - magma
'0'
```

If Sage did not do this garbage collection, then every single time you ever create any magma object from a sage object, e.g., by doing `magma(m)`, you would use up a lot of memory in that Magma session. This would lead to a horrible memory leak situation, which would make the Magma interface nearly useless for serious work.

## 20.5 Other Examples

We compute a space of modular forms with character.

```
sage: N = 20
sage: D = 20
sage: eps_top = fundamental_discriminant(D)
sage: eps = magma.KroneckerCharacter(eps_top, RationalField()) # optional - magma
sage: M2 = magma.ModularForms(eps) # optional - magma
sage: print M2 # optional - magma
Space of modular forms on Gamma_1(5) ...
sage: print M2.Basis() # optional - magma
[
1 + 10*q^2 + 20*q^3 + 20*q^5 + 60*q^7 + ...
q + q^2 + 2*q^3 + 3*q^4 + 5*q^5 + 2*q^6 + ...
]
```

In Sage/Python (and sort of C++) coercion of an element  $x$  into a structure  $S$  is denoted by  $S(x)$ . This also works for the Magma interface:

```
sage: G = magma.DirichletGroup(20) # optional - magma
sage: G.AssignNames(['a', 'b']) # optional - magma
sage: (G.1).Modulus() # optional - magma
20
sage: e = magma.DirichletGroup(40)(G.1) # optional - magma
sage: print e # optional - magma
$.1
sage: print e.Modulus() # optional - magma
40
```

We coerce some polynomial rings into Magma:

```
sage: R.<y> = PolynomialRing(QQ)
sage: S = magma(R) # optional - magma
sage: print S # optional - magma
Univariate Polynomial Ring in y over Rational Field
sage: S.1 # optional - magma
y
```

This example illustrates that Sage doesn't magically extend how Magma implicit coercion (what there is, at least) works. The errors below are the result of Magma having a rather limited automatic coercion system compared to Sage's:

```
sage: R.<x> = ZZ[]
sage: x * 5
5*x
sage: x * 1.0
x
sage: x * (2/3)
2/3*x
sage: y = magma(x) # optional - magma
sage: y * 5 # optional - magma
```

```
5*x
sage: y * 1.0                                     # optional - magma
$.1
sage: y * (2/3)                                    # optional - magma
Traceback (most recent call last):
...
TypeError: Error evaluating Magma code.
...
Argument types given: RngUPolElt[RngInt], FldRatElt
```

**AUTHORS:**

- William Stein (2005): initial version
- William Stein (2006-02-28): added extensive tab completion and interactive IPython documentation support.
- William Stein (2006-03-09): added `nvals` argument for `magma.functions...`

```
class sage.interfaces.magma.Magma (maxread=10000, script_subdirectory=None, logfile=None,
                                     server=None, server_tmpdir=None, user_config=False,
                                     seed=None)
    Bases: sage.interfaces.expect.Expect
```

Interface to the Magma interpreter.

Type `magma.[tab]` for a list of all the functions available from your Magma install. Type `magma.Function?` for Magma's help about a given `Function`. Type `magma(...)` to create a new Magma object, and `magma.eval(...)` to run a string using Magma (and get the result back as a string).

**Note:** If you do not own a local copy of Magma, try using the `magma_free` command instead, which uses the free demo web interface to Magma.

EXAMPLES:

You must use `nvals = 0` to call a function that doesn't return anything, otherwise you'll get an error. (`nvals` is the number of return values.)

[illegible]

**Attach** (*filename*)

Attach the given file to the running instance of Magma.

Attaching a file in Magma makes all intrinsics defined in the file available to the shell. Moreover, if the file doesn't start with the `freeze;` command, then the file is reloaded whenever it is changed. Note that functions and procedures defined in the file are *not* available. For only those, use `magma.load(filename)`.

INPUT:

- filename - a string

EXAMPLES: Attaching a file that exists is fine:

```
sage: SAGE_EXTCODE = SAGE_ENV['SAGE_EXTCODE'] # optional - magma
sage: magma.attach('%s/magma/sage/basic.m'%SAGE_EXTCODE) # optional - magma
```

Attaching a file that doesn't exist raises an exception:

```
sage: SAGE_EXTCODE = SAGE_ENV['SAGE_EXTCODE'] # optional - magma
sage: magma.attach('%s/magma/sage/basic2.m'%SAGE_EXTCODE) # optional - magma
```

```
Traceback (most recent call last):
...
RuntimeError: Error evaluating Magma code...
```

**AttachSpec** (*filename*)

Attach the given spec file to the running instance of Magma.

This can attach numerous other files to the running Magma (see the Magma documentation for more details).

INPUT:

- filename - a string

EXAMPLES:

```
sage: SAGE_EXTCODE = SAGE_ENV['SAGE_EXTCODE']           # optional - magma
sage: magma.attach_spec('%s/magma/spec'%SAGE_EXTCODE)  # optional - magma
sage: magma.attach_spec('%s/magma/spec2'%SAGE_EXTCODE) # optional - magma
Traceback (most recent call last):
...
RuntimeError: Can't open package spec file .../magma/spec2 for reading (No such file or dire
```

**GetVerbose** (*type*)

Get the verbosity level of a given algorithm class etc. in Magma.

INPUT:

- type - string (e.g. 'Groebner'), see Magma documentation

---

**Note:** This method is provided to be consistent with the Magma naming convention.

---

EXAMPLES:

```
sage: magma.SetVerbose("Groebner", 2)           # optional - magma
sage: magma.GetVerbose("Groebner")              # optional - magma
2
```

**SetVerbose** (*type, level*)

Set the verbosity level for a given algorithm class etc. in Magma.

INPUT:

- type - string (e.g. 'Groebner'), see Magma documentation
- level - integer = 0

---

**Note:** This method is provided to be consistent with the Magma naming convention.

---

```
sage: magma.SetVerbose("Groebner", 2)           # optional - magma
sage: magma.GetVerbose("Groebner")              # optional - magma
2
```

**attach** (*filename*)

Attach the given file to the running instance of Magma.

Attaching a file in Magma makes all intrinsics defined in the file available to the shell. Moreover, if the file doesn't start with the `freeze;` command, then the file is reloaded whenever it is changed. Note that functions and procedures defined in the file are *not* available. For only those, use `magma.load(filename)`.

INPUT:

- filename - a string

EXAMPLES: Attaching a file that exists is fine:

```
sage: SAGE_EXTCODE = SAGE_ENV['SAGE_EXTCODE'] # optional - magma
sage: magma.attach('%s/magma/sage/basic.m'%SAGE_EXTCODE) # optional - magma
```

Attaching a file that doesn't exist raises an exception:

```
sage: SAGE_EXTCODE = SAGE_ENV['SAGE_EXTCODE'] # optional - magma
sage: magma.attach('%s/magma/sage/basic2.m'%SAGE_EXTCODE) # optional - magma
Traceback (most recent call last):
...
RuntimeError: Error evaluating Magma code...
```

### **attach\_spec** (filename)

Attach the given spec file to the running instance of Magma.

This can attach numerous other files to the running Magma (see the Magma documentation for more details).

INPUT:

- filename - a string

EXAMPLES:

```
sage: SAGE_EXTCODE = SAGE_ENV['SAGE_EXTCODE'] # optional - magma
sage: magma.attach_spec('%s/magma/spec'%SAGE_EXTCODE) # optional - magma
sage: magma.attach_spec('%s/magma/spec2'%SAGE_EXTCODE) # optional - magma
Traceback (most recent call last):
...
RuntimeError: Can't open package spec file ../magma/spec2 for reading (No such file or dire
```

### **bar\_call** (left, name, gens, nvals=1)

This is a wrapper around the Magma constructor

nameleft gens

returning nvals.

INPUT:

- left - something coerceable to a magma object
- name - name of the constructor, e.g., sub, quo, ideal, etc.
- gens - if a list/tuple, each item is coerced to magma; otherwise gens itself is converted to magma
- nvals - positive integer; number of return values

OUTPUT: a single magma object if nvals == 1; otherwise a tuple of nvals magma objects.

EXAMPLES: The bar\_call function is used by the sub, quo, and ideal methods of Magma elements. Here we illustrate directly using bar\_call to create quotients:

```
sage: V = magma.RModule(ZZ, 3) # optional - magma
sage: V # optional - magma
RModule(IntegerRing(), 3)
sage: magma.bar_call(V, 'quo', [[1,2,3]], nvals=1) # optional - magma
RModule(IntegerRing(), 2)
sage: magma.bar_call(V, 'quo', [[1,2,3]], nvals=2) # optional - magma
(RModule(IntegerRing(), 2),
 Mapping from: RModule(IntegerRing(), 3) to RModule(IntegerRing(), 2))
sage: magma.bar_call(V, 'quo', V, nvals=2) # optional - magma
```



```
(RModule(IntegerRing(), 0),
 Mapping from: RModule(IntegerRing(), 3) to RModule(IntegerRing(), 0))
```

**chdir** (*dir*)

Change the Magma interpreter's current working directory.

INPUT:

- *dir* – a string

EXAMPLES:

```
sage: magma.chdir('/')          # optional - magma
sage: magma.eval('System("pwd")') # optional - magma
'/'
```

**clear** (*var*)

Clear the variable named *var* and make it available to be used again.

INPUT:

- *var* – a string

EXAMPLES:

```
sage: magma = Magma()          # optional - magma
sage: magma.clear('foo')       # sets foo to 0 in magma; optional - magma
sage: magma.eval('foo')        # optional - magma
'0'
```

Because we cleared *foo*, it is set to be used as a variable name in the future:

```
sage: a = magma('10')          # optional - magma
sage: a.name()                 # optional - magma
'foo'
```

The following tests that the whole variable clearing and freeing system is working correctly.

```
sage: magma = Magma()          # optional - magma
sage: a = magma('100')         # optional - magma
sage: a.name()                 # optional - magma
'_sage_[1]'
sage: del a                    # optional - magma
sage: b = magma('257')         # optional - magma
sage: b.name()                 # optional - magma
'_sage_[1]'
sage: del b                    # optional - magma
sage: magma('_sage_[1]')       # optional - magma
0
```

**console** ()

Run a command line Magma session. This session is completely separate from this Magma interface.

EXAMPLES:

```
sage: magma.console()          # not tested
Magma V2.14-9      Sat Oct 11 2008 06:36:41 on one      [Seed = 1157408761]
Type ? for help.  Type <Ctrl>-D to quit.
>
Total time: 2.820 seconds, Total memory usage: 3.95MB
```

**cputime** (*t=None*)

Return the CPU time in seconds that has elapsed since this Magma session started. This is a floating point number, computed by Magma.

If *t* is given, then instead return the floating point time from when *t* seconds had elapsed. This is useful for computing elapsed times between two points in a running program.

INPUT:

- *t* - float (default: None); if not None, return cputime since *t*

OUTPUT:

- float - seconds

EXAMPLES:

```
sage: type(magma.cputime())           # optional - magma
<type 'float'>
sage: magma.cputime()                 # random, optional - magma
1.9399999999999999
sage: t = magma.cputime()             # optional - magma
sage: magma.cputime(t)                 # random, optional - magma
0.02
```

**eval** (*x, strip=True, \*\*kws*)

Evaluate the given block *x* of code in Magma and return the output as a string.

INPUT:

- *x* - string of code
- *strip* - ignored

OUTPUT: string

EXAMPLES:

We evaluate a string that involves assigning to a variable and printing.

```
sage: magma.eval("a := 10; print 2+a;")   # optional - magma
'12'
```

We evaluate a large input line (note that no weird output appears and that this works quickly).

```
sage: magma.eval("a := %s;"%(10^10000))   # optional - magma
''
```

Verify that trac 9705 is fixed:

```
sage: nl=chr(10) # newline character
sage: magma.eval( # optional - magma
...  "_<x>:=PolynomialRing(Rationals());"+nl+
...  "repeat"+nl+
...  "  g:=3*b*x^4+18*c*x^3-6*b^2*x^2-6*b*c*x-b^3-9*c^2 where b:=Random([-10..10]) where c:=F
...  "until g ne 0 and Roots(g) ne [];" +nl+
...  "print \"success\";")
'success'
```

Verify that trac 11401 is fixed:

```
sage: nl=chr(10) # newline character
sage: magma.eval("a:=3;"+nl+"b:=5;") == nl # optional - magma
True
```

```
sage: magma.eval("[a,b];") # optional - magma
['[ 3, 5 ]']
```

**function\_call** (*function*, *args*=[], *params*={}, *nvals*=1)

Return result of evaluating a Magma function with given input, parameters, and asking for nvals as output.

INPUT:

- *function* - string, a Magma function name
- *args* - list of objects coercible into this magma interface
- *params* - Magma parameters, passed in after a colon
- *nvals* - number of return values from the function to ask Magma for

OUTPUT: MagmaElement or tuple of nvals MagmaElement's

EXAMPLES:

```
sage: magma.function_call('Factorization', 100) # optional - magma
[ <2, 2>, <5, 2> ]
sage: magma.function_call('NextPrime', 100, {'Proof':False}) # optional - magma
101
sage: magma.function_call('PolynomialRing', [QQ,2]) # optional - magma
Polynomial ring of rank 2 over Rational Field
Order: Lexicographical
Variables: $.1, $.2
```

Next, we illustrate multiple return values:

```
sage: magma.function_call('IsSquare', 100) # optional - magma
true
sage: magma.function_call('IsSquare', 100, nvals=2) # optional - magma
(true, 10)
sage: magma.function_call('IsSquare', 100, nvals=3) # optional - magma
Traceback (most recent call last):
...
RuntimeError: Error evaluating Magma code...
Runtime error in :=: Expected to assign 3 value(s) but only computed 2 value(s)
```

**get** (*var*)

Get the value of the variable *var*.

INPUT:

- *var* - string; name of a variable defined in the Magma session

OUTPUT:

- *string* - string representation of the value of the variable.

EXAMPLES:

```
sage: magma.set('abc', '2 + 3/5') # optional - magma
sage: magma.get('abc') # optional - magma
'13/5'
```

**get\_verbose** (*type*)

Get the verbosity level of a given algorithm class etc. in Magma.

INPUT:

- *type* - string (e.g. 'Groebner'), see Magma documentation

EXAMPLES:

```
sage: magma.set_verbose("Groebner", 2)      # optional - magma
sage: magma.get_verbose("Groebner")         # optional - magma
2
```

**help**(s)

Return Magma help on string s.

This returns what typing ?s would return in Magma.

INPUT:

- s - string

OUTPUT: string

EXAMPLES:

```
sage: magma.help("NextPrime")              # optional - magma
=====
PATH: /magma/ring-field-algebra/integer/prime/next-previous/NextPrime
KIND: Intrinsic
=====
NextPrime(n) : RngIntElt -> RngIntElt
NextPrime(n: parameter) : RngIntElt -> RngIntElt
...
```

**ideal**(L)

Return the Magma ideal defined by L.

INPUT:

- L - a list of elements of a Sage multivariate polynomial ring.

OUTPUT: The magma ideal generated by the elements of L.

EXAMPLES:

```
sage: R.<x,y> = QQ[]
sage: magma.ideal([x^2, y^3*x])           # optional - magma
Ideal of Polynomial ring of rank 2 over Rational Field
Order: Graded Reverse Lexicographical
Variables: x, y
Homogeneous
Basis:
[
x^2,
x*y^3
]
```

**load**(filename)

Load the file with given filename using the 'load' command in the Magma shell.

Loading a file in Magma makes all the functions and procedures in the file available. The file should not contain any intrinsics (or you'll get errors). It also runs code in the file, which can produce output.

INPUT:

- filename - string

OUTPUT: output printed when loading the file

EXAMPLES:

```

sage: filename = os.path.join(SAGE_TMP, 'a.m')
sage: open(filename, 'w').write('function f(n) return n^2; end function;\nprint "hi";')
sage: print magma.load(filename)          # optional - magma
Loading ".../a.m"
hi
sage: magma('f(12)')                     # optional - magma
144

```

**objgens** (*value, gens*)

Create a new object with given value and gens.

INPUT:

- **value** - something coercible to an element of this Magma interface
- **gens** - string; comma separated list of variable names

OUTPUT: new Magma element that is equal to value with given gens

EXAMPLES:

```

sage: R = magma.objgens('PolynomialRing(Rationals(),2)', 'alpha,beta') # optional - magma
sage: R.gens()                  # optional - magma
[alpha, beta]

```

Because of how Magma works you can use this to change the variable names of the generators of an object:

```

sage: S = magma.objgens(R, 'X,Y') # optional - magma
sage: R # optional - magma
Polynomial ring of rank 2 over Rational Field
Order: Lexicographical
Variables: X, Y
sage: S # optional - magma
Polynomial ring of rank 2 over Rational Field
Order: Lexicographical
Variables: X, Y

```

**set** (*var, value*)

Set the variable var to the given value in the Magma interpreter.

INPUT:

- **var** - string; a variable name
- **value** - string; what to set var equal to

EXAMPLES:

```

sage: magma.set('abc', '2 + 3/5') # optional - magma
sage: magma('abc')                # optional - magma
13/5

```

**set\_seed** (*seed=None*)

Sets the seed for the Magma interpreter. The seed should be an integer.

EXAMPLES:

```

sage: m = Magma() # optional - magma
sage: m.set_seed(1) # optional - magma
1
sage: [m.Random(100) for i in range(5)] # optional - magma
[95, 20, 61, 59, 24]

```

**set\_verbose** (*type, level*)

Set the verbosity level for a given algorithm, class, etc. in Magma.

INPUT:

- *type* - string (e.g. 'Groebner')
- *level* - integer = 0

EXAMPLES:

```
sage: magma.set_verbose("Groebner", 2)      # optional - magma
sage: magma.get_verbose("Groebner")        # optional - magma
2
```

**trait\_names** (*verbose=True, use\_disk\_cache=True*)

Return a list of all Magma commands.

This is used as a hook to enable custom command completion.

Magma doesn't provide any fast way to make a list of all commands, which is why caching is done by default. Note that an adverse impact of caching is that *new* commands are not picked up, e.g., user defined variables or functions.

INPUT:

- *verbose* - bool (default: True); whether to verbosely output status info the first time the command list is built
- *use\_disk\_cache* - bool (default: True); use cached command list, which is saved to disk.

OUTPUT: list of strings

EXAMPLES:

```
sage: len(magma.trait_names(verbose=False))  # random, optional - magma
7261
```

**version** ()

Return the version of Magma that you have in your PATH on your computer.

OUTPUT:

- *numbers* - 3-tuple: major, minor, etc.
- *string* - version as a string

EXAMPLES:

```
sage: magma.version()      # random, optional - magma
((2, 14, 9), 'V2.14-9')
```

**class** `sage.interfaces.magma.MagmaElement` (*parent, value, is\_name=False, name=None*)Bases: `sage.interfaces.expect.ExpectElement`**AssignNames** (*names*)

EXAMPLES:

```
sage: G = magma.DirichletGroup(20)      # optional - magma
sage: G.AssignNames(['a', 'b'])         # optional - magma
sage: G.1                                # optional - magma
a

sage: G.Elements()                      # optional - magma
[
1,
```

```
a,
b,
a*b
]
```

**assign\_names** (*names*)

EXAMPLES:

```
sage: G = magma.DirichletGroup(20)      # optional - magma
sage: G.AssignNames(['a','b'])          # optional - magma
sage: G.1                                # optional - magma
a
```

```
sage: G.Elements()                      # optional - magma
[
1,
a,
b,
a*b
]
```

**eval** (*\*args*)

Evaluate self at the inputs.

INPUT:

•\*args - import arguments

OUTPUT: self(\*args)

EXAMPLES:

```
sage: f = magma('Factorization')        # optional - magma
sage: f.evaluate(15)                     # optional - magma
[ <3, 1>, <5, 1> ]
sage: f(15)                              # optional - magma
[ <3, 1>, <5, 1> ]
sage: f = magma('GCD')                   # optional - magma
sage: f.evaluate(15,20)                  # optional - magma
5
```

**evaluate** (*\*args*)

Evaluate self at the inputs.

INPUT:

•\*args - import arguments

OUTPUT: self(\*args)

EXAMPLES:

```
sage: f = magma('Factorization')        # optional - magma
sage: f.evaluate(15)                     # optional - magma
[ <3, 1>, <5, 1> ]
sage: f(15)                              # optional - magma
[ <3, 1>, <5, 1> ]
sage: f = magma('GCD')                   # optional - magma
sage: f.evaluate(15,20)                  # optional - magma
5
```

**gen**(*n*)

Return the *n*-th generator of this Magma element. Note that generators are 1-based in Magma rather than 0 based!

INPUT:

- *n* - a *positive* integer

OUTPUT: MagmaElement

EXAMPLES:

```
sage: k.<a> = GF(9)
sage: magma(k).gen(1)          # optional -- magma
a
sage: R.<s,t,w> = k[]
sage: m = magma(R)            # optional -- magma
sage: m.gen(1)                 # optional -- magma
s
sage: m.gen(2)                 # optional -- magma
t
sage: m.gen(3)                 # optional -- magma
w
sage: m.gen(0)                 # optional -- magma
Traceback (most recent call last):
...
IndexError: index must be positive since Magma indexes are 1-based
sage: m.gen(4)                 # optional -- magma
Traceback (most recent call last):
...
IndexError: list index out of range
```

**gen\_names**()

Return list of Magma variable names of the generators of self.

---

**Note:** As illustrated below, these are not the print names of the the generators of the Magma object, but special variable names in the Magma session that reference the generators.

---

EXAMPLES:

```
sage: R.<x,zw> = QQ[]
sage: S = magma(R)            # optional - magma
sage: S.gen_names()           # optional - magma
('_sage_...', '_sage_...')
sage: magma(S.gen_names()[1]) # optional - magma
zw
```

**gens**()

Return generators for self.

If self is named *X* is Magma, this function evaluates *X*.1, *X*.2, etc., in Magma until an error occurs. It then returns a Sage list of the resulting *X*.*i*. Note - I don't think there is a Magma command that returns the list of valid *X*.*i*. There are numerous ad hoc functions for various classes but nothing systematic. This function gets around that problem. Again, this is something that should probably be reported to the Magma group and fixed there.

AUTHORS:

- William Stein (2006-07-02)

EXAMPLES:



```

sage: magma("VectorSpace(RationalField(),3)").gens()           # optional - magma
[(1 0 0), (0 1 0), (0 0 1)]
sage: magma("AbelianGroup(EllipticCurve([1..5]))").gens()      # optional - magma
[$.1]

```

**get\_magma\_attribute(attrname)**

Return value of a given Magma attribute. This is like self.attrname in Magma.

OUTPUT: MagmaElement

EXAMPLES:

```

sage: V = magma("VectorSpace(RationalField(),10)")           # optional - magma
sage: V.set_magma_attribute('M', "hello")                   # optional - magma
sage: V.get_magma_attribute('M')                             # optional - magma
hello
sage: V.M                                                      # optional - magma
hello

```

**ideal(gens)**

Return the ideal of self with given list of generators.

INPUT:

- gens - object or list/tuple of generators

OUTPUT:

- magma element - a Magma ideal

EXAMPLES:

```

sage: R = magma('PolynomialRing(RationalField())')           # optional - magma
sage: R.assign_names(['x'])                                   # optional - magma
sage: x = R.1                                                  # optional - magma
sage: R.ideal([x^2 - 1, x^3 - 1])                             # optional - magma
Ideal of Univariate Polynomial Ring in x over Rational Field generated by x - 1

```

**list\_attributes()**

Return the attributes of self, obtained by calling the ListAttributes function in Magma.

OUTPUT: list of strings

EXAMPLES: We observe that vector spaces in Magma have numerous funny and mysterious attributes.

```

sage: V = magma("VectorSpace(RationalField(),2)")           # optional - magma
sage: v = V.list_attributes(); v.sort(); v                   # optional - magma
['Coroots', 'Involution', 'M', 'RootDatum', 'Roots', 'StrLocalData', 'T', 'decomp', 'eisen',

```

**methods(any=False)**

Return signatures of all Magma intrinsics that can take self as the first argument, as strings.

INPUT:

- any - (bool: default is False) if True, also include signatures with Any as first argument.

OUTPUT: list of strings

EXAMPLES:

```

sage: v = magma('2/3').methods()                             # optional - magma
sage: v[0]                                                     # optional - magma
"/ *'..."

```

**quo** (*gens*)

Return the quotient of self by the given object or list of generators.

INPUT:

- gens - object or list/tuple of generators

OUTPUT:

- magma element - the quotient object
- magma element - mapping from self to the quotient object

EXAMPLES:

```
sage: V = magma('VectorSpace(RationalField(),3)') # optional - magma
sage: V.quo([[1,2,3], [1,1,2]]) # optional - magma
(Full Vector space of degree 1 over Rational Field, Mapping from: Full Vector space of degree 1 over Rational Field)
```

We illustrate quotienting out by an object instead of a list of generators:

```
sage: W = V.sub([ [1,2,3], [1,1,2] ]) # optional - magma
sage: V.quo(W) # optional - magma
(Full Vector space of degree 1 over Rational Field, Mapping from: Full Vector space of degree 1 over Rational Field)
```

We quotient a ZZ module out by a submodule.

```
sage: V = magma.RModule(ZZ,3); V # optional - magma
RModule(IntegerRing(), 3)
sage: W, phi = V.quo([[1,2,3]]) # optional - magma
sage: W # optional - magma
RModule(IntegerRing(), 2)
sage: phi # optional - magma
Mapping from: RModule(IntegerRing(), 3) to RModule(IntegerRing(), 2)
```

**set\_magma\_attribute** (*attrname, value*)

INPUTS: attrname - string value - something coercible to a MagmaElement

EXAMPLES:

```
sage: V = magma("VectorSpace(RationalField(),2)") # optional - magma
sage: V.set_magma_attribute('M',10) # optional - magma
sage: V.get_magma_attribute('M') # optional - magma
10
sage: V.M # optional - magma
10
```

**sub** (*gens*)

Return the sub-object of self with given gens.

INPUT:

- gens - object or list/tuple of generators

EXAMPLES:

```
sage: V = magma('VectorSpace(RationalField(),3)') # optional - magma
sage: W = V.sub([ [1,2,3], [1,1,2] ]); W # optional - magma
Vector space of degree 3, dimension 2 over Rational Field
Generators:
(1 2 3)
(1 1 2)
Echelonized basis:
```

```
(1 0 1)
(0 1 1)
```

**trait\_names()**

Return all Magma functions that have this Magma element as first input. This is used for tab completion.

**Note:** This function can unfortunately be slow if there are a very large number of functions, e.g., when self is an integer. (This could be fixed by the addition of an appropriate function to the Magma kernel, which is something that can only be done by the Magma developers.)

OUTPUT:

- list - sorted list of distinct strings

EXAMPLES:

```
sage: R.<x> = ZZ[]                                # optional - magma
sage: v = magma(R).trait_names()                  # optional - magma
sage: v                                            # optional - magma
['*', '+', '.', '/', 'eq', 'in', 'meet', 'subset', ...]
```

**class** sage.interfaces.magma.**MagmaFunction**(parent, name)

Bases: sage.interfaces.expect.ExpectFunction

**class** sage.interfaces.magma.**MagmaFunctionElement**(obj, name)

Bases: sage.interfaces.expect.FunctionElement

**class** sage.interfaces.magma.**MagmaGBDefaultContext**(magma=None)

Context to force preservation of verbosity options for Magma's Groebner basis computation.

**class** sage.interfaces.magma.**MagmaGBLogPrettyPrinter**(verbosity=1, style='magma')

A device which filters Magma Groebner basis computation logs.

**flush()**

EXAMPLE:

```
sage: from sage.interfaces.magma import MagmaGBLogPrettyPrinter
sage: logs = MagmaGBLogPrettyPrinter()
sage: logs.flush()
```

**write(s)**

EXAMPLE:

```
sage: P.<x,y,z> = GF(32003)[]
sage: I = sage.rings.ideal.Katsura(P)
sage: _ = I.groebner_basis('magma', prot=True) # indirect doctest, optional - magma

Homogeneous weights search
...
Total Faugere F4 time: ..., real time: ...
```

sage.interfaces.magma.**extcode\_dir()**

Return directory that contains all the Magma extcode. This is put in a writable directory owned by the user, since when attached, Magma has to write sig and lck files.

EXAMPLES:

```
sage: sage.interfaces.magma.extcode_dir()
'...dir_.../data/'
```

`sage.interfaces.magma.is_MagmaElement(x)`

Return True if x is of type MagmaElement, and False otherwise.

INPUT:

- x - any object

OUTPUT: bool

EXAMPLES:

```
sage: from sage.interfaces.magma import is_MagmaElement
sage: is_MagmaElement(2)
False
sage: is_MagmaElement(magma(2))           # optional - magma
True
```

`sage.interfaces.magma.magma_console()`

Run a command line Magma session.

EXAMPLES:

```
sage: magma_console()           # not tested
Magma V2.14-9      Sat Oct 11 2008 06:36:41 on one      [Seed = 1157408761]
Type ? for help.  Type <Ctrl>-D to quit.
>
Total time: 2.820 seconds, Total memory usage: 3.95MB
```

`sage.interfaces.magma.magma_gb_standard_options(func)`

Decorator to force default options for Magma.

EXAMPLE:

```
sage: P.<a,b,c,d,e> = PolynomialRing(GF(127))
sage: J = sage.rings.ideal.Cyclic(P).homogenize()
sage: from sage.misc.sageinspect import sage_getsource
sage: "mself" in sage_getsource(J._groebner_basis_magma)
True
```

`sage.interfaces.magma.magma_version()`

Return the version of Magma that you have in your PATH on your computer.

OUTPUT:

- numbers - 3-tuple: major, minor, etc.
- string - version as a string

EXAMPLES:

```
sage: magma_version()           # random, optional - magma
((2, 14, 9), 'V2.14-9')
```

`sage.interfaces.magma.reduce_load_Magma()`

Used in unpickling a Magma interface.

This functions just returns the global default Magma interface.

EXAMPLES:

```
sage: sage.interfaces.magma.reduce_load_Magma()
Magma
```

## INTERFACE TO THE FREE ONLINE MAGMA CALCULATOR

**class** `sage.interfaces.magma_free.MagmaExpr`  
Bases: `str`

**class** `sage.interfaces.magma_free.MagmaFree`  
Evaluate MAGMA code without requiring that MAGMA be installed on your computer by using the free online MAGMA calculator.

**EXAMPLES:** `sage: magma_free("Factorization(9290348092384)")` # optional - internet [ <2, 5>  
<290323377887, 1> ]

**eval** (`x`, *\*\*kws*)

`sage.interfaces.magma_free.magma_free_eval` (`code`, *strip=True*, *columns=0*)

Use the free online MAGMA calculator to evaluate the given input code and return the answer as a string.

**LIMITATIONS:** The code must evaluate in at most 20 seconds and there is a limitation on the amount of RAM.

**EXAMPLES:** `sage: magma_free("Factorization(9290348092384)")` # optional - internet [ <2, 5>  
<290323377887, 1> ]



## INTERFACE TO MAPLE

### AUTHORS:

- William Stein (2005): maple interface
- Gregg Musiker (2006-02-02): tutorial
- William Stein (2006-03-05): added tab completion, e.g., `maple.[tab]`, and help, e.g., `maple.sin?`.

You must have the optional commercial Maple interpreter installed and available as the command `maple` in your PATH in order to use this interface. You do not have to install any optional Sage packages.

Type `maple.[tab]` for a list of all the functions available from your Maple install. Type `maple.[tab]?` for Maple's help about a given function. Type `maple(...)` to create a new Maple object, and `maple.eval(...)` to run a string using Maple (and get the result back as a string).

### EXAMPLES:

```
sage: maple('3 * 5')                # optional - maple
15
sage: maple.eval('ifactor(2005)')    # optional - maple
''(5)*''(401)'
sage: maple.ifactor(2005)            # optional - maple
''(5)*''(401)
sage: maple.fsolve('x^2=cos(x)+4', 'x=0..5') # optional - maple
1.914020619
sage: maple.factor('x^5 - y^5')      # optional - maple
(x-y)*(x^4+x^3*y+x^2*y^2+x*y^3+y^4)
```

If the string “error” (case insensitive) occurs in the output of anything from Maple, a `RuntimeError` exception is raised.

## 22.1 Tutorial

### AUTHORS:

- Gregg Musiker (2006-02-02): initial version.

This tutorial is based on the Maple Tutorial for number theory from <http://www.math.mun.ca/~drideout/m3370/numtheory.html>.

There are several ways to use the Maple Interface in Sage. We will discuss two of those ways in this tutorial.

1. If you have a maple expression such as

```
factor( (x^5-1));
```

We can write that in sage as

```
sage: maple('factor(x^5-1)') # optional - maple
(x-1)*(x^4+x^3+x^2+x+1)
```

Notice, there is no need to use a semicolon.

2. Since Sage is written in Python, we can also import maple commands and write our scripts in a Pythonic way. For example, `factor()` is a maple command, so we can also factor in Sage using

```
sage: maple('(x^5-1)').factor() # optional - maple
(x-1)*(x^4+x^3+x^2+x+1)
```

where `expression.command()` means the same thing as `command(expression)` in Maple. We will use this second type of syntax whenever possible, resorting to the first when needed.

```
sage: maple('(x^12-1)/(x-1)').simplify() # optional - maple
x^11+x^10+x^9+x^8+x^7+x^6+x^5+x^4+x^3+x^2+x+1
```

The normal command will always reduce a rational function to the lowest terms. The factor command will factor a polynomial with rational coefficients into irreducible factors over the ring of integers. So for example,

```
sage: maple('(x^12-1)').factor() # optional - maple
(x-1)*(x+1)*(x^2+x+1)*(x^2-x+1)*(x^2+1)*(x^4-x^2+1)
```

```
sage: maple('(x^28-1)').factor() # optional - maple
(x-1)*(x^6+x^5+x^4+x^3+x^2+x+1)*(x+1)*(1-x+x^2-x^3+x^4-x^5+x^6)*(x^2+1)*(x^12-x^10+x^8-x^6+x^4-x^2+1)
```

Another important feature of maple is its online help. We can access this through sage as well. After reading the description of the command, you can press q to immediately get back to your original prompt.

Incidentally you can always get into a maple console by the command

```
sage: maple.console() # not tested
sage: !maple # not tested
```

Note that the above two commands are slightly different, and the first is preferred.

For example, for help on the maple command `fibonacci`, we type

```
sage: maple.help('fibonacci') # not tested, since it uses a pager
```

We see there are two choices. Type

```
sage: maple.help('combinat, fibonacci') # not tested, since it uses a pager
```

We now see how the Maple command `fibonacci` works under the combinatorics package. Try typing in

```
sage: maple.fibonacci(10) # optional - maple
fibonacci(10)
```

You will get `fibonacci(10)` as output since Maple has not loaded the combinatorics package yet. To rectify this type

```
sage: maple('combinat[fibonacci]')(10) # optional - maple
55
```

instead.

If you want to load the combinatorics package for future calculations, in Sage this can be done as



```
sage: maple.with_package('combinat')           # optional - maple
```

or

```
sage: maple.load('combinat')                   # optional - maple
```

Now if we type `maple.fibonacci(10)`, we get the correct output:

```
sage: maple.fibonacci(10)                       # optional - maple
55
```

Some common maple packages include `combinat`, `linalg`, and `numtheory`. To produce the first 19 Fibonacci numbers, use the `sequence` command.

```
sage: maple('seq(fibonacci(i),i=1..19)')        # optional - maple
1, 1, 2, 3, 5, 8, 13, 21, 34, 55, 89, 144, 233, 377, 610, 987, 1597, 2584,
4181
```

Two other useful Maple commands are `ifactor` and `isprime`. For example

```
sage: maple.isprime(maple.fibonacci(27))        # optional - maple
false
sage: maple.ifactor(maple.fibonacci(27))        # optional - maple
``(2)*``(17)*``(53)*``(109)
```

Note that the `isprime` function that is included with Sage (which uses PARI) is better than the Maple one (it is faster and gives a provably correct answer, whereas Maple is sometimes wrong).

```
sage: alpha = maple('(1+sqrt(5))/2')            # optional - maple
sage: beta = maple('(1-sqrt(5))/2')            # optional - maple
sage: f19 = alpha^19 - beta^19/maple('sqrt(5)') # optional - maple
sage: f19                                       # optional - maple
(1/2+1/2*5^(1/2))^19-1/5*(1/2-1/2*5^(1/2))^19*5^(1/2)
sage: f19.simplify()                          # somewhat randomly ordered output; optional - maple
6765+5778/5*5^(1/2)
```

Let's say we want to write a maple program now that squares a number if it is positive and cubes it if it is negative. In maple, that would look like

```
mysqcu := proc(x)
if x > 0 then x^2;
else x^3; fi;
end;
```

In Sage, we write

```
sage: mysqcu = maple('proc(x) if x > 0 then x^2 else x^3 fi end') # optional - maple
sage: mysqcu(5)                                                  # optional - maple
25
sage: mysqcu(-5)                                                 # optional - maple
-125
```

More complicated programs should be put in a separate file and loaded.

```
class sage.interfaces.maple.Maple(maxread=100, script_subdirectory=None, server=None,
server_tmpdir=None, logfile=None, ulimit=None)
```

Bases: `sage.interfaces.expect.Expect`

Interface to the Maple interpreter.

Type `maple.[tab]` for a list of all the functions available from your Maple install. Type `maple.[tab]?` for Maple's help about a given function. Type `maple(...)` to create a new Maple object, and `maple.eval(...)` to run a string using Maple (and get the result back as a string).

**clear** (*var*)

Clear the variable named `var`.

To clear a Maple variable, you must assign 'itself' to itself. In Maple 'expr' prevents expr to be evaluated.

EXAMPLES:

```
sage: maple.set('xx', '2') # optional - maple
```

```
sage: maple.get('xx') # optional - maple
```

'2'

```
sage: maple.clear('xx')      # optional - maple
```

```
sage: maple.get('xx') # optional - maple
```

'xx'

completions ( $s$ )

Return all commands that complete the command starting with the string `s`. This is like typing `s[Ctrl-T]` in the maple interpreter.

EXAMPLES:

```
sage: c = maple.completions('di') # optional - maple
```

```
sage: 'divide' in c # optional - maple
```

True

```
console ()
```

Spawn a new Maple command-line session.

EXAMPLES:

```
sage: maple.console() # not tested
```

Maple 11 (IBM INTEL LINUX)

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< > Waterloo Maple Inc.

```
| Type ? for help.
```

 $\succ$ **cputime** ( $t=None$ )

Returns the amount of CPU time that the Maple session has used. If  $\tau$  is not None, then it returns the difference between the current CPU time and  $\tau$ .

EXAMPLES:

```
sage: t = maple.cputime() # optional - maple
```

```
sage: t = random_element(
# random; optional - maple
```

0.02

```
sage: x = maple('x')           # optional - maple
```

```
sage: maple.diff(x^2, x) # optional - maple
```

2.\*x

```
sage: maple.cputime(t)      # random; optional - maple
```

0.0

**expect ()**

Returns the pexpect object for this Maple session.

EXAMPLES:

```
sage: m = Maple()           # optional - maple
```

```
sage: m.expect() is None # optional - maple
```

```

True
sage: m._start()           # optional - maple
sage: m.expect()          # optional - maple
<pexpect.spawn instance at 0x...>
sage: m.quit()             # optional - maple

```

**get** (*var*)

Get the value of the variable *var*.

## EXAMPLES:

```

sage: maple.set('xx', '2') # optional - maple
sage: maple.get('xx')      # optional - maple
'2'

```

**help** (*str*)

Display Maple help about *str*.

This is the same as typing "?str" in the Maple console.

INPUT:

- *str* - a string to search for in the maple help system

## EXAMPLES:

```

sage: maple.help('digamma') #not tested
Psi - the Digamma and Polygamma functions
...

```

**load** (*package*)

Make a package of Maple procedures available in the interpreter.

INPUT:

- *package* - string

EXAMPLES: Some functions are unknown to Maple until you use with to include the appropriate package.

```

sage: maple.quit()      # reset maple; optional -- maple
sage: maple('partition(10)') # optional - maple
partition(10)
sage: maple('bell(10)')  # optional - maple
bell(10)
sage: maple.with_package('combinat') # optional - maple
sage: maple('partition(10)') # optional - maple
[[1, 1, 1, 1, 1, 1, 1, 1, 1, 1], [1, 1, 1, 1, 1, 1, 1, 1, 2], [1, 1, 1, 1, 1, 1, 2, 2], [1,
sage: maple('bell(10)')  # optional - maple
115975
sage: maple('fibonacci(10)') # optional - maple
55

```

**set** (*var*, *value*)

Set the variable *var* to the given value.

## EXAMPLES:

```

sage: maple.set('xx', '2') # optional - maple
sage: maple.get('xx')      # optional - maple
'2'

```

**source** (*s*)

Display the Maple source (if possible) about *s*. This is the same as returning the output produced by the following Maple commands:

```
interface(verboseproc=2): print(s)
```

INPUT:

- *s* - a string representing the function whose source code you want

EXAMPLES:

```
sage: maple.source('curry') #not tested
p -> subs('_X' = args[2 .. nargs], () -> p(_X, args))
```

**trait\_names** (*verbose=True, use\_disk\_cache=True*)

Returns a list of all the commands defined in Maple and optionally (per default) store them to disk.

EXAMPLES:

```
sage: c = maple.trait_names(use_disk_cache=False, verbose=False) # optional - maple
sage: len(c) > 100 # optional - maple
True
sage: 'dilog' in c # optional - maple
True
```

**with\_package** (*package*)

Make a package of Maple procedures available in the interpreter.

INPUT:

- *package* - string

EXAMPLES: Some functions are unknown to Maple until you use *with* to include the appropriate package.

```
sage: maple.quit() # reset maple; optional -- maple
sage: maple('partition(10)') # optional - maple
partition(10)
sage: maple('bell(10)') # optional - maple
bell(10)
sage: maple.with_package('combinat') # optional - maple
sage: maple('partition(10)') # optional - maple
[[1, 1, 1, 1, 1, 1, 1, 1, 1, 1], [1, 1, 1, 1, 1, 1, 1, 1, 2], [1, 1, 1, 1, 1, 1, 2, 2], [1,
sage: maple('bell(10)') # optional - maple
115975
sage: maple('fibonacci(10)') # optional - maple
55
```

```
class sage.interfaces.maple.MapleElement (parent, value, is_name=False, name=None)
```

Bases: `sage.interfaces.expect.ExpectElement`

**trait\_names** ()

EXAMPLES:

```
sage: a = maple(2) # optional - maple
sage: 'sin' in a.trait_names() # optional - maple
True
```

```
class sage.interfaces.maple.MapleFunction (parent, name)
```

Bases: `sage.interfaces.expect.ExpectFunction`

```
class sage.interfaces.maple.MapleFunctionElement (obj, name)
```

Bases: `sage.interfaces.expect.FunctionElement`

`sage.interfaces.maple.maple_console()`

Spawn a new Maple command-line session.

EXAMPLES:

```
sage: maple_console() #not tested
      |^/|      Maple 11 (IBM INTEL LINUX)
      ._|\\|    |/_|. Copyright (c) Maplesoft, a division of Waterloo Maple Inc. 2007
      \ MAPLE / All rights reserved. Maple is a trademark of
      <____> Waterloo Maple Inc.
      |      Type ? for help.
>
```

`sage.interfaces.maple.reduce_load_Maple()`

Returns the maple object created in `sage.interfaces.maple`.

EXAMPLES:

```
sage: from sage.interfaces.maple import reduce_load_Maple
sage: reduce_load_Maple()
Maple
```



## INTERFACE TO MATHEMATICA

The Mathematica interface will only work if Mathematica is installed on your computer with a command line interface that runs when you give the `math` command. The interface lets you send certain Sage objects to Mathematica, run Mathematica functions, import certain Mathematica expressions to Sage, or any combination of the above.

To send a Sage object `sobj` to Mathematica, call `mathematica(sobj)`. This exports the Sage object to Mathematica and returns a new Sage object wrapping the Mathematica expression/variable, so that you can use the Mathematica variable from within Sage. You can then call Mathematica functions on the new object; for example:

```
sage: mobj = mathematica(x^2-1)           # optional - mathematica
sage: mobj.Factor()                       # optional - mathematica
(-1 + x)*(1 + x)
```

In the above example the factorization is done using Mathematica's `Factor[]` function.

To see Mathematica's output you can simply print the Mathematica wrapper object. However if you want to import Mathematica's output back to Sage, call the Mathematica wrapper object's `sage()` method. This method returns a native Sage object:

```
sage: mobj = mathematica(x^2-1)           # optional - mathematica
sage: mobj2 = mobj.Factor(); mobj2        # optional - mathematica
(-1 + x)*(1 + x)
sage: mobj2.parent()                     # optional - mathematica
Mathematica
sage: sobj = mobj2.sage(); sobj           # optional - mathematica
(x - 1)*(x + 1)
sage: sobj.parent()                     # optional - mathematica
Symbolic Ring
```

If you want to run a Mathematica function and don't already have the input in the form of a Sage object, then it might be simpler to input a string to `mathematica(expr)`. This string will be evaluated as if you had typed it into Mathematica:

```
sage: mathematica('Factor[x^2-1]')        # optional - mathematica
(-1 + x)*(1 + x)
sage: mathematica('Range[3]')            # optional - mathematica
{1, 2, 3}
```

If you don't want Sage to go to the trouble of creating a wrapper for the Mathematica expression, then you can call `mathematica.eval(expr)`, which returns the result as a Mathematica `AsciiArtString` formatted string. If you want the result to be a string formatted like Mathematica's `InputForm`, call `repr(mobj)` on the wrapper object `mobj`. If you want a string formatted in Sage style, call `mobj._sage_repr()`:

```
sage: mathematica.eval('x^2 - 1')          # optional - mathematica
2
```

```
      -1 + x
sage: repr(mathematica('Range[3]'))      # optional - mathematica
'{1, 2, 3}'
sage: mathematica('Range[3]')._sage_repr() # optional - mathematica
'[1, 2, 3]'
```

Finally, if you just want to use a Mathematica command line from within Sage, the function `mathematica_console()` dumps you into an interactive command-line Mathematica session. This is an enhanced version of the usual Mathematica command-line, in that it provides readline editing and history (the usual one doesn't!)

## 23.1 Tutorial

We follow some of the tutorial from <http://library.wolfram.com/conferences/devconf99/withoff/Basic1.html/>.

For any of this to work you must buy and install the Mathematica program, and it must be available as the command `math` in your `PATH`.

### 23.1.1 Syntax

Now make 1 and add it to itself. The result is a Mathematica object.

```
sage: m = mathematica
sage: a = m(1) + m(1); a      # optional - mathematica
2
sage: a.parent()              # optional - mathematica
Mathematica
sage: m('1+1')                # optional - mathematica
2
sage: m(3)**m(50)              # optional - mathematica
717897987691852588770249
```

The following is equivalent to `Plus[2, 3]` in Mathematica:

```
sage: m = mathematica
sage: m(2).Plus(m(3))         # optional - mathematica
5
```

We can also compute  $7(2 + 3)$ .

```
sage: m(7).Times(m(2).Plus(m(3))) # optional - mathematica
35
sage: m('7(2+3)')               # optional - mathematica
35
```

### 23.1.2 Some typical input

We solve an equation and a system of two equations:

```
sage: eqn = mathematica('3x + 5 == 14') # optional - mathematica
sage: eqn                                # optional - mathematica
5 + 3*x == 14
sage: eqn.Solve('x')                    # optional - mathematica
```



```

{{x -> 3}}
sage: sys = mathematica('{x^2 - 3y == 3, 2x - y == 1}') # optional - mathematica
sage: print sys                                         # optional - mathematica
      2
      {x  - 3 y == 3, 2 x - y == 1}
sage: sys.Solve('{x, y}')                               # optional - mathematica
{{x -> 0, y -> -1}, {x -> 6, y -> 11}}

```

### 23.1.3 Assignments and definitions

If you assign the mathematica 5 to a variable *c* in Sage, this does not affect the *c* in Mathematica.

```

sage: c = m(5)                                         # optional - mathematica
sage: print m('b + c x')                               # optional - mathematica
      b + c x
sage: print m('b') + c*m('x')                         # optional - mathematica
      b + 5 x

```

The Sage interfaces changes Sage lists into Mathematica lists:

```

sage: m = mathematica
sage: eq1 = m('x^2 - 3y == 3')                        # optional - mathematica
sage: eq2 = m('2x - y == 1')                          # optional - mathematica
sage: v = m([eq1, eq2]); v                            # optional - mathematica
{x^2 - 3*y == 3, 2*x - y == 1}
sage: v.Solve(['x', 'y'])                             # optional - mathematica
{{x -> 0, y -> -1}, {x -> 6, y -> 11}}

```

### 23.1.4 Function definitions

Define mathematica functions by simply sending the definition to the interpreter.

```

sage: m = mathematica
sage: _ = mathematica('f[p_] = p^2');                 # optional - mathematica
sage: m('f[9]')                                       # optional - mathematica
81

```

### 23.1.5 Numerical Calculations

We find the  $x$  such that  $e^x - 3x = 0$ .

```

sage: e = mathematica('Exp[x] - 3x == 0') # optional - mathematica
sage: e.FindRoot(['x', 2])                 # optional - mathematica
{x -> 1.512134551657842}

```

Note that this agrees with what the PARI interpreter gp produces:

```

sage: gp('solve(x=1,2,exp(x)-3*x)')
1.512134551657842473896739678             # 32-bit
1.5121345516578424738967396780720387046   # 64-bit

```

Next we find the minimum of a polynomial using the two different ways of accessing Mathematica:

```

sage: mathematica('FindMinimum[x^3 - 6x^2 + 11x - 5, {x,3}]') # optional - mathematica
{0.6150998205402516, {x -> 2.5773502699629733}}
sage: f = mathematica('x^3 - 6x^2 + 11x - 5') # optional - mathematica
sage: f.FindMinimum(['x', 3]) # optional - mathematica
{0.6150998205402516, {x -> 2.5773502699629733}}

```

### 23.1.6 Polynomial and Integer Factorization

We factor a polynomial of degree 200 over the integers.

```

sage: R.<x> = PolynomialRing(ZZ)
sage: f = (x**100+17*x+5)*(x**100-5*x+20)
sage: f
x^200 + 12*x^101 + 25*x^100 - 85*x^2 + 315*x + 100
sage: g = mathematica(str(f)) # optional - mathematica
sage: print g # optional - mathematica
      2      100      101      200
100 + 315 x - 85 x + 25 x + 12 x + x
sage: g # optional - mathematica
100 + 315*x - 85*x^2 + 25*x^100 + 12*x^101 + x^200
sage: print g.Factor() # optional - mathematica
      100      100
(20 - 5 x + x ) (5 + 17 x + x )

```

We can also factor a multivariate polynomial:

```

sage: f = mathematica('x^6 + (-y - 2)*x^5 + (y^3 + 2*y)*x^4 - y^4*x^3') # optional - mathematica
sage: print f.Factor() # optional - mathematica
      3      2      3
x (x - y) (-2 x + x + y )

```

We factor an integer:

```

sage: n = mathematica(2434500) # optional - mathematica
sage: n.FactorInteger() # optional - mathematica
{{2, 2}, {3, 2}, {5, 3}, {541, 1}}
sage: n = mathematica(2434500) # optional - mathematica
sage: F = n.FactorInteger(); F # optional - mathematica
{{2, 2}, {3, 2}, {5, 3}, {541, 1}}
sage: F[1] # optional - mathematica
{2, 2}
sage: F[4] # optional - mathematica
{541, 1}

```

We can also load the ECM package and factoring using it:

```

sage: _ = mathematica.eval("<<NumberTheory`FactorIntegerECM`"); # optional - mathematica
sage: mathematica.FactorIntegerECM('932901*939321') # optional - mathematica
8396109

```

## 23.2 Long Input

The Mathematica interface reads in even very long input (using files) in a robust manner.

```

sage: t = '%s'%10^10000 # ten thousand character string.
sage: a = mathematica(t) # optional - mathematica
sage: a = mathematica.eval(t) # optional - mathematica

```

## 23.3 Loading and saving

Mathematica has an excellent `InputForm` function, which makes saving and loading Mathematica objects possible. The first examples test saving and loading to strings.

```

sage: x = mathematica(pi/2) # optional - mathematica
sage: print x # optional - mathematica
      Pi
      --
      2
sage: loads(dumps(x)) == x # optional - mathematica
True
sage: n = x.N(50) # optional - mathematica
sage: print n # optional - mathematica
      1.5707963267948966192313216916397514420985846996876
sage: loads(dumps(n)) == n # optional - mathematica
True

```

## 23.4 Complicated translations

The `mobj.sage()` method tries to convert a Mathematica object to a Sage object. In many cases, it will just work. In particular, it should be able to convert expressions entirely consisting of:

- numbers, i.e. integers, floats, complex numbers;
- functions and named constants also present in Sage, where:
  - Sage knows how to translate the function or constant's name from Mathematica's, or
  - the Sage name for the function or constant is trivially related to Mathematica's;
- symbolic variables whose names don't pathologically overlap with objects already defined in Sage.

This method will not work when Mathematica's output includes:

- strings;
- functions unknown to Sage;
- Mathematica functions with different parameters/parameter order to the Sage equivalent.

If you want to convert more complicated Mathematica expressions, you can instead call `mobj._sage_()` and supply a translation dictionary:

```

sage: m = mathematica('NewFn[x]') # optional - mathematica
sage: m._sage_(locals={'NewFn': sin}) # optional - mathematica
sin(x)

```

For more details, see the documentation for `._sage_()`.

OTHER Examples:

```
sage: def math_bessel_K(nu, x):
...     return mathematica(nu).BesselK(x).N(20)
...
sage: math_bessel_K(2, I)                                # optional - mathematica
0.180489972066962*I - 2.592886175491197                # 32-bit
-2.59288617549119697817 + 0.18048997206696202663*I    # 64-bit

sage: slist = [[1, 2], 3., 4 + I]
sage: mlist = mathematica(slist); mlist                # optional - mathematica
{{1, 2}, 3., 4 + I}
sage: slist2 = list(mlist); slist2                     # optional - mathematica
[[1, 2], 3., 4 + I]
sage: slist2[0]                                         # optional - mathematica
{1, 2}
sage: slist2[0].parent()                               # optional - mathematica
Mathematica
sage: slist3 = mlist.sage(); slist3                    # optional - mathematica
[[1, 2], 3.0, I + 4]

sage: mathematica('10.^80')                            # optional - mathematica
1.*^80
sage: mathematica('10.^80').sage()                     # optional - mathematica
1e+80
```

**AUTHORS:**

- William Stein (2005): first version
- Doug Cutrell (2006-03-01): Instructions for use under Cygwin/Windows.
- Felix Lawrence (2009-08-21): Added support for importing Mathematica lists and floats with exponents.

```
class sage.interfaces.mathematica.Mathematica(maxread=100, script_subdirectory=None,
                                              logfile=None, server=None,
                                              server_tmpdir=None)
```

Bases: `sage.interfaces.expect.Expect`

Interface to the Mathematica interpreter.

**chdir** (*dir*)

Change Mathematica's current working directory.

EXAMPLES:

```
sage: mathematica.chdir('/')                            # optional - mathematica
sage: mathematica('Directory[]')                       # optional - mathematica
"/"
```

**console** (*readline=True*)

**eval** (*code, strip=True, \*\*kwds*)

**get** (*var, ascii\_art=False*)

Get the value of the variable *var*.

**AUTHORS:**

- William Stein
- Kiran Kedlaya (2006-02-04): suggested using `InputForm`

**help** (*cmd*)

**set** (*var*, *value*)  
Set the variable *var* to the given value.

**trait\_names** ()

**class** sage.interfaces.mathematica.**MathematicaElement** (*parent*, *value*, *is\_name=False*,  
*name=None*)

Bases: sage.interfaces.expect.ExpectElement

**N** (\**args*)

EXAMPLES:

```
sage: mathematica('Pi').N(10)      # optional -- mathematica
3.1415926536
sage: mathematica('Pi').N(50)      # optional -- mathematica
3.14159265358979323846264338327950288419716939937511
```

**save\_image** (*filename*, *ImageSize=600*)

Save a mathematica graphics

INPUT:

- *filename* – string. The filename to save as. The extension determines the image file format.
- *ImageSize* – integer. The size of the resulting image.

EXAMPLES:

```
sage: P = mathematica('Plot[Sin[x],{x,-2Pi,4Pi}]') # optional - mathematica
sage: filename = tmp_filename()                  # optional - mathematica
sage: P.save(filename, ImageSize=800)             # optional - mathematica
```

**show** (*ImageSize=600*)

Show a mathematica expression immediately.

This method attempts to display the graphics immediately, without waiting for the currently running code (if any) to return to the command line. Be careful, calling it from within a loop will potentially launch a large number of external viewer programs.

INPUT:

- *ImageSize* – integer. The size of the resulting image.

OUTPUT:

This method does not return anything. Use `save()` if you want to save the figure as an image.

EXAMPLES:

```
sage: P = mathematica('Plot[Sin[x],{x,-2Pi,4Pi}]') # optional - mathematica
sage: show(P)                                     # optional - mathematica
sage: P.show(ImageSize=800)                       # optional - mathematica
sage: Q = mathematica('Sin[x Cos[y]]/Sqrt[1-x^2]') # optional - mathematica
sage: show(Q)                                     # optional - mathematica
<html><div class="math">\frac{\sin (x \cos (y))}{\sqrt{1-x^2}}</div></html>
```

**str** ()

**class** sage.interfaces.mathematica.**MathematicaFunction** (*parent*, *name*)

Bases: sage.interfaces.expect.ExpectFunction

**class** sage.interfaces.mathematica.**MathematicaFunctionElement** (*obj*, *name*)

Bases: sage.interfaces.expect.FunctionElement

sage.interfaces.mathematica.**clean\_output** (*s*)

`sage.interfaces.mathematica.mathematica_console(readline=True)`

`sage.interfaces.mathematica.reduce_load(X)`

## INTERFACE TO MATLAB

According to their website, MATLAB is “a high-level language and interactive environment that enables you to perform computationally intensive tasks faster than with traditional programming languages such as C, C++, and Fortran.”

The commands in this section only work if you have the “matlab” interpreter installed and available in your PATH. It’s not necessary to install any special Sage packages.

EXAMPLES:

```
sage: matlab.eval('2+2')           # optional - matlab
'\nans =\n\n      4\n'
```

```
sage: a = matlab(10)               # optional - matlab
sage: a**10                        # optional - matlab
      1.0000e+10
```

AUTHORS:

- William Stein (2006-10-11)

### 24.1 Tutorial

EXAMPLES:

```
sage: matlab('4+10')              # optional - matlab
14
sage: matlab('date')              # optional - matlab; random output
18-Oct-2006
sage: matlab('5*10 + 6')          # optional - matlab
56
sage: matlab('(6+6)/3')           # optional - matlab
4
sage: matlab('9')^2               # optional - matlab
81
sage: a = matlab(10); b = matlab(20); c = matlab(30)    # optional - matlab
sage: avg = (a+b+c)/3 ; avg       # optional - matlab
20
sage: parent(avg)                 # optional - matlab
Matlab
```

```
sage: my_scalar = matlab('3.1415') # optional - matlab
sage: my_scalar                    # optional - matlab
3.1415
sage: my_vector1 = matlab('[1,5,7]') # optional - matlab
```

```

sage: my_vector1                                # optional - matlab
1      5      7
sage: my_vector2 = matlab(' [1;5;7]')          # optional - matlab
sage: my_vector2                                # optional - matlab
1
5
7
sage: my_vector1 * my_vector2                  # optional - matlab
75

sage: row_vector1 = matlab(' [1 2 3]')          # optional - matlab
sage: row_vector2 = matlab(' [3 2 1]')          # optional - matlab
sage: matrix_from_row_vec = matlab(' [%s; %s]'%(row_vector1.name(), row_vector2.name())) # optional - matlab
sage: matrix_from_row_vec                      # optional - matlab
1      2      3
3      2      1

sage: column_vector1 = matlab(' [1;3]')         # optional - matlab
sage: column_vector2 = matlab(' [2;8]')         # optional - matlab
sage: matrix_from_col_vec = matlab(' [%s %s]%(column_vector1.name(), column_vector2.name())') # optional - matlab
sage: matrix_from_col_vec                     # optional - matlab
1      2
3      8

sage: my_matrix = matlab(' [8, 12, 19; 7, 3, 2; 12, 4, 23; 8, 1, 1]') # optional - matlab
sage: my_matrix                               # optional - matlab
8      12      19
7      3      2
12     4      23
8      1      1

sage: combined_matrix = matlab(' [%s, %s]%(my_matrix.name(), my_matrix.name())')
sage: combined_matrix                         # optional - matlab
8      12      19      8      12      19
7      3      2      7      3      2
12     4      23     12     4      23
8      1      1      8      1      1

sage: tm = matlab(' 0.5:2:10')                # optional - matlab
sage: tm                                       # optional - matlab
0.5000    2.5000    4.5000    6.5000    8.5000

sage: my_vector1 = matlab(' [1,5,7]')          # optional - matlab
sage: my_vector1(1)                          # optional - matlab
1
sage: my_vector1(2)                          # optional - matlab
5
sage: my_vector1(3)                          # optional - matlab
7

```

Matrix indexing works as follows:

```

sage: my_matrix = matlab(' [8, 12, 19; 7, 3, 2; 12, 4, 23; 8, 1, 1]') # optional - matlab
sage: my_matrix(3,2)                                         # optional - matlab
4

```

Setting using parenthesis cannot work (because of how the Python language works). Use square brackets or the set



function:

```
sage: my_matrix = matlab('[8, 12, 19; 7, 3, 2; 12, 4, 23; 8, 1, 1]') # optional - matlab
sage: my_matrix.set(2,3, 1999) # optional - matlab
sage: my_matrix # optional - matlab
      8      12      19
      7       3     1999
     12       4      23
      8       1       1
```

```
class sage.interfaces.matlab.Matlab(maxread=100, script_subdirectory=None, logfile=None,
                                     server=None, server_tmpdir=None)
```

Bases: `sage.interfaces.expect.Expect`

Interface to the Matlab interpreter.

EXAMPLES:

```
sage: a = matlab('[ 1, 1, 2; 3, 5, 8; 13, 21, 33 ]') # optional - matlab
sage: b = matlab('[ 1; 3; 13]') # optional - matlab
sage: c = a * b # optional - matlab
sage: print c # optional - matlab
      30
     122
     505
```

**chdir** (*directory*)

Change MATLAB's current working directory.

EXAMPLES:

```
sage: matlab.chdir('/') # optional - matlab
sage: matlab.pwd() # optional - matlab
/
```

**console** ()

**get** (*var*)

Get the value of the variable *var*.

EXAMPLES:

```
sage: s = matlab.eval('a = 2') # optional - matlab
sage: matlab.get('a') # optional - matlab
      2
```

**sage2matlab\_matrix\_string** (*A*)

Return an matlab matrix from a Sage matrix.

INPUT: A Sage matrix with entries in the rationals or reals.

OUTPUT: A string that evaluates to an Matlab matrix.

EXAMPLES:

```
sage: M33 = MatrixSpace(QQ,3,3)
sage: A = M33([1,2,3,4,5,6,7,8,0])
sage: matlab.sage2matlab_matrix_string(A) # optional - matlab
'[1, 2, 3; 4, 5, 6; 7, 8, 0]'
```

AUTHOR:

•David Joyner and William Stein

**set** (*var*, *value*)  
Set the variable *var* to the given value.

**strip\_answer** (*s*)  
Returns the string *s* with Matlab's answer prompt removed.

EXAMPLES:

```
sage: s = '\nans =\n\n      2\n'
sage: matlab.strip_answer(s)
'      2'
```

**version** ()

**whos** ()

**class** sage.interfaces.matlab.**MatlabElement** (*parent*, *value*, *is\_name=False*, *name=None*)  
Bases: sage.interfaces.expect.ExpectElement

**set** (*i*, *j*, *x*)

sage.interfaces.matlab.**matlab\_console** ()

This requires that the optional matlab program be installed and in your PATH, but no optional Sage packages need be installed.

EXAMPLES:

```
sage: matlab_console()           # optional - matlab; not tested
                                   < M A T L A B >
                                   Copyright 1984-2006 The MathWorks, Inc.
```

```
...
>> 2+3
```

```
ans =
```

```
5
```

```
quit
```

Typing quit exits the matlab console and returns you to Sage. matlab, like Sage, remembers its history from one session to another.

sage.interfaces.matlab.**matlab\_version** ()  
Return the version of Matlab installed.

EXAMPLES:

```
sage: matlab_version()          # random; optional - matlab
'7.2.0.283 (R2006a)'
```

sage.interfaces.matlab.**reduce\_load\_Matlab** ()

## PEXPECT INTERFACE TO MAXIMA

Maxima is a free GPL'd general purpose computer algebra system whose development started in 1968 at MIT. It contains symbolic manipulation algorithms, as well as implementations of special functions, including elliptic functions and generalized hypergeometric functions. Moreover, Maxima has implementations of many functions relating to the invariant theory of the symmetric group  $S_n$ . (However, the commands for group invariants, and the corresponding Maxima documentation, are in French.) For many links to Maxima documentation see <http://maxima.sourceforge.net/documentation.html>.

AUTHORS:

- William Stein (2005-12): Initial version
- David Joyner: Improved documentation
- William Stein (2006-01-08): Fixed bug in parsing
- William Stein (2006-02-22): comparisons (following suggestion of David Joyner)
- William Stein (2006-02-24): *greatly* improved robustness by adding sequence numbers to IO bracketing in `_eval_line`
- Robert Bradshaw, Nils Bruin, Jean-Pierre Flori (2010,2011): Binary library interface

This is the interface used by the maxima object:

```
sage: type(maxima)
<class 'sage.interfaces.maxima.Maxima'>
```

If the string “error” (case insensitive) occurs in the output of anything from Maxima, a `RuntimeError` exception is raised.

EXAMPLES: We evaluate a very simple expression in Maxima.

```
sage: maxima('3 * 5')
15
```

We factor  $x^5 - y^5$  in Maxima in several different ways. The first way yields a Maxima object.

```
sage: F = maxima.factor('x^5 - y^5')
sage: F
-(y-x) * (y^4+x*y^3+x^2*y^2+x^3*y+x^4)
sage: type(F)
<class 'sage.interfaces.maxima.MaximaElement'>
```

Note that Maxima objects can also be displayed using “ASCII art”; to see a normal linear representation of any Maxima object `x`. Just use the print command: use `str(x)`.

```
sage: print F
```

$$-(y-x) \left( y^4 + x y^3 + x^2 y^2 + x^3 y + x^4 \right)$$

You can always use `repr(x)` to obtain the linear representation of an object. This can be useful for moving maxima data to other systems.

```
sage: repr(F)
'-(y-x)*(y^4+x*y^3+x^2*y^2+x^3*y+x^4)'
```

```
sage: F.str()
'-(y-x)*(y^4+x*y^3+x^2*y^2+x^3*y+x^4)'
```

The `maxima.eval` command evaluates an expression in maxima and returns the result as a *string* not a maxima object.

```
sage: print maxima.eval('factor(x^5 - y^5)')
-(y-x)*(y^4+x*y^3+x^2*y^2+x^3*y+x^4)
```

We can create the polynomial  $f$  as a Maxima polynomial, then call the factor method on it. Notice that the notation `f.factor()` is consistent with how the rest of Sage works.

```
sage: f = maxima('x^5 - y^5')
sage: f^2
(x^5-y^5)^2
sage: f.factor()
-(y-x)*(y^4+x*y^3+x^2*y^2+x^3*y+x^4)
```

Control-C interruption works well with the maxima interface, because of the excellent implementation of maxima. For example, try the following sum but with a much bigger range, and hit control-C.

```
sage: maxima('sum(1/x^2, x, 1, 10)')
1968329/1270080
```

## 25.1 Tutorial

We follow the tutorial at <http://maxima.sourceforge.net/docs/intromax/intromax.html>.

```
sage: maxima('1/100 + 1/101')
201/10100
```

```
sage: a = maxima('(1 + sqrt(2))^5'); a
(sqrt(2)+1)^5
sage: a.expand()
29*sqrt(2)+41
```

```
sage: a = maxima('(1 + sqrt(2))^5')
sage: float(a)
82.01219330881975
sage: a.numer()
82.01219330881975
```

```
sage: maxima.eval('fpprec : 100')
'100'
```

```

sage: a.bfloat()
8.2012129330881975641524897300208124427852048438593149412212371240173124187540110412666123849550160561

sage: maxima('100!')
9332621544394415268169923885626670049071596826438162146859296389521759999322991560894146397615651828

sage: f = maxima('(x + 3*y + x^2*y)^3')
sage: f.expand()
x^6*y^3+9*x^4*y^3+27*x^2*y^3+27*y^3+3*x^5*y^2+18*x^3*y^2+27*x*y^2+3*x^4*y+9*x^2*y+x^3
sage: f.subst('x=5/z')
(5/z+25*y/z^2+3*y)^3
sage: g = f.subst('x=5/z')
sage: h = g.ratsimp(); h
(27*y^3*z^6+135*y^2*z^5+(675*y^3+225*y)*z^4+(2250*y^2+125)*z^3+(5625*y^3+1875*y)*z^2+9375*y^2*z+15625)*z^3
sage: h.factor()
(3*y*z^2+5*z+25*y)^3/z^6

sage: eqn = maxima(['a+b*c=1', 'b-a*c=0', 'a+b=5'])
sage: s = eqn.solve(['a,b,c']); s
[[a=(25*sqrt(79)*%i+25)/(6*sqrt(79)*%i-34),b=(5*sqrt(79)*%i+5)/(sqrt(79)*%i+11),c=(sqrt(79)*%i+1)/10]]

```

Here is an example of solving an algebraic equation:

```

sage: maxima('x^2+y^2=1').solve('y')
[y=-sqrt(1-x^2),y=sqrt(1-x^2)]
sage: maxima('x^2 + y^2 = (x^2 - y^2)/sqrt(x^2 + y^2)').solve('y')
[y=-sqrt((-y^2-x^2)*sqrt(y^2+x^2)+x^2),y=sqrt((-y^2-x^2)*sqrt(y^2+x^2)+x^2)]

```

You can even nicely typeset the solution in latex:

```

sage: latex(s)
\left[ \left[ a=\frac{25\sqrt{79}i+25}{6\sqrt{79}i-34} , \quad b=\frac{5\sqrt{79}i+5}{\sqrt{79}i+11}, c=\frac{\sqrt{79}i+1}{10} \right] \right]

```

To have the above appear onscreen via `xdvi`, type `view(s)`. (TODO: For OS X should create pdf output and use preview instead?)

```

sage: e = maxima('sin(u + v) * cos(u)^3'); e
cos(u)^3*sin(v+u)
sage: f = e.trigexpand(); f
cos(u)^3*(cos(u)*sin(v)+sin(u)*cos(v))
sage: f.trigreduce()
(sin(v+4*u)+sin(v-2*u))/8+(3*sin(v+2*u)+3*sin(v))/8
sage: w = maxima('3 + k*i')
sage: f = w^2 + maxima('%e')^w
sage: f.realpart()
%e^3*cos(k)-k^2+9

sage: f = maxima('x^3 * %e^(k*x) * sin(w*x)'); f
x^3*%e^(k*x)*sin(w*x)
sage: f.diff('x')
k*x^3*%e^(k*x)*sin(w*x)+3*x^2*%e^(k*x)*sin(w*x)+w*x^3*%e^(k*x)*cos(w*x)
sage: f.integrate('x')
(((k*w^6+3*k^3*w^4+3*k^5*w^2+k^7)*x^3+(3*w^6+3*k^2*w^4-3*k^4*w^2-3*k^6)*x^2+(-18*k*w^4-12*k^3*w^2+6*k^5)*x+k^7)*%e^(k*x))/k^7

sage: f = maxima('1/x^2')
sage: f.integrate('x', 1, 'inf')

```

```
1
sage: g = maxima('f/sinh(k*x)^4')
sage: g.taylor('x', 0, 3)
f/(k^4*x^4)-2*f/(3*k^2*x^2)+11*f/45-62*k^2*f*x^2/945
```

```
sage: maxima.taylor('asin(x)', 'x', 0, 10)
x+x^3/6+3*x^5/40+5*x^7/112+35*x^9/1152
```

## 25.2 Examples involving matrices

We illustrate computing with the matrix whose  $i, j$  entry is  $i/j$ , for  $i, j = 1, \dots, 4$ .

```
sage: f = maxima.eval('f[i,j] := i/j')
sage: A = maxima('genmatrix(f,4,4)'); A
matrix([[1, 1/2, 1/3, 1/4], [2, 1, 2/3, 1/2], [3, 3/2, 1, 3/4], [4, 2, 4/3, 1]])
sage: A.determinant()
0
sage: A.echelon()
matrix([[1, 1/2, 1/3, 1/4], [0, 0, 0, 0], [0, 0, 0, 0], [0, 0, 0, 0]])
sage: A.eigenvalues()
[[0, 4], [3, 1]]
sage: A.eigenvectors()
[[[0, 4], [3, 1]], [[1, 0, 0, -4], [0, 1, 0, -2], [0, 0, 1, -4/3]], [[1, 2, 3, 4]]]
```

We can also compute the echelon form in Sage:

```
sage: B = matrix(QQ, A)
sage: B.echelon_form()
[ 1 1/2 1/3 1/4]
[ 0  0  0  0]
[ 0  0  0  0]
[ 0  0  0  0]
sage: B.charpoly('x').factor()
(x - 4) * x^3
```

## 25.3 Laplace Transforms

We illustrate Laplace transforms:

```
sage: _ = maxima.eval("f(t) := t*sin(t)")
sage: maxima("laplace(f(t),t,s)")
2*s/(s^2+1)^2

sage: maxima("laplace(delta(t-3),t,s)") #Dirac delta function
%e^-(3*s)

sage: _ = maxima.eval("f(t) := exp(t)*sin(t)")
sage: maxima("laplace(f(t),t,s)")
1/(s^2-2*s+2)
```



You can formally evaluate sums (note the `nusum` command):

```
sage: S = maxima('nusum(exp(1+2*i/n), i, 1, n)')
sage: print S
```

$$\frac{e^{\frac{2}{n} + 3}}{e^{\frac{1}{n}} - 1} - \frac{e^{\frac{2}{n} + 1}}{e^{\frac{1}{n}} + 1}$$

We formally compute the limit as  $n \rightarrow \infty$  of  $2S/n$  as follows:

```
sage: T = S*maxima('2/n')
sage: T.tlimit('n', 'inf')
%e^3-%e
```

## 25.6 Miscellaneous

Obtaining digits of  $\pi$ :

```
sage: maxima.eval('fpprec : 100')
'100'
sage: maxima(pi).bfloat()
3.141592653589793238462643383279502884197169399375105820974944592307816406286208998628034825342117068
```

Defining functions in maxima:

```
sage: maxima.eval('fun[a] := a^2')
'fun[a]:=a^2'
sage: maxima('fun[10]')
100
```

## 25.7 Interactivity

Unfortunately maxima doesn't seem to have a non-interactive mode, which is needed for the Sage interface. If any Sage call leads to maxima interactively answering questions, then the questions can't be answered and the maxima session may hang. See the discussion at <http://www.ma.utexas.edu/pipermail/maxima/2005/011061.html> for some ideas about how to fix this problem. An example that illustrates this problem is `maxima.eval('integrate(exp(a*x), x, 0, inf)')`.

## 25.8 Latex Output

To TeX a maxima object do this:

```
sage: latex(maxima('sin(u) + sinh(v^2)'))
\sinh v^2+\sin u
```

Here's another example:

```
sage: g = maxima('exp(3*i*x)/(6*i) + exp(i*x)/(2*i) + c')
sage: latex(g)
-{{i\,e^{3\,i\,x}}\over{6}}-{{i\,e^{i\,x}}\over{2}}+c
```



**Note:** Using `maxima.eval` for long input is much less robust, and is not recommended.

EXAMPLES:

```
sage: maxima.set('xxxxx', '2')
sage: maxima.get('xxxxx')
'2'
sage: maxima.clear('xxxxx')
sage: maxima.get('xxxxx')
'xxxxx'
```

**get** (*var*)

Get the string value of the variable *var*.

EXAMPLES:

```
sage: maxima.set('xxxxx', '2')
sage: maxima.get('xxxxx')
'2'
```

**lisp** (*cmd*)

Send a lisp command to Maxima.

---

**Note:** The output of this command is very raw - not pretty.

---

EXAMPLES:

```
sage: maxima.lisp("(+ 2 17)")    # random formatted output
:lisp (+ 2 17)
19
(
```

**set** (*var*, *value*)

Set the variable *var* to the given value.

INPUT:

- *var* - string
- *value* - string

EXAMPLES:

```
sage: maxima.set('xxxxx', '2')
sage: maxima.get('xxxxx')
'2'
```

**set\_seed** (*seed=None*)

[http://maxima.sourceforge.net/docs/manual/maxima\\_10.html](http://maxima.sourceforge.net/docs/manual/maxima_10.html) `make_random_state` (*n*) returns a new random state object created from an integer seed value equal to *n* modulo  $2^{32}$ . *n* may be negative.

EXAMPLES:

```
sage: m = Maxima()
sage: m.set_seed(1)
1
sage: [m.random(100) for i in range(5)]
[45, 39, 24, 68, 63]
```

**class** `sage.interfaces.maxima.MaximaElement` (*parent*, *value*, *is\_name=False*, *name=None*)

Bases: `sage.interfaces.maxima_abstract.MaximaAbstractElement`,  
`sage.interfaces.expect.ExpectElement`

Element of Maxima through Pexpect interface.

## EXAMPLES:

Elements of this class should not be created directly. The targeted parent should be used instead:

```
sage: maxima(3)
3
sage: maxima(cos(x)+e^234)
cos(_SAGE_VAR_x)+%e^234
```

**display2d** (*onscreen=True*)

Return the 2d string representation of this Maxima object.

## EXAMPLES:

```
sage: F = maxima('x^5 - y^5').factor()
sage: F.display2d()
      4      3      2      2      3      4
      - (y - x) (y + x y + x y + x y + x )
```

**class** sage.interfaces.maxima.**MaximaElementFunction** (*parent, name, defn, args, latex*)

Bases: sage.interfaces.maxima.MaximaElement, sage.interfaces.maxima\_abstract.MaximaAbstractFunction

Maxima user-defined functions.

## EXAMPLES:

Elements of this class should not be created directly. The method function of the targeted parent should be used instead:

```
sage: maxima.function('x,y','h(x)*y')
h(x)*y
```

**class** sage.interfaces.maxima.**MaximaFunction** (*parent, name*)

Bases: sage.interfaces.maxima\_abstract.MaximaAbstractFunction, sage.interfaces.expect.ExpectFunction

**class** sage.interfaces.maxima.**MaximaFunctionElement** (*obj, name*)

Bases: sage.interfaces.maxima\_abstract.MaximaAbstractFunctionElement, sage.interfaces.expect.FunctionElement

sage.interfaces.maxima.**is\_MaximaElement** (*x*)

Returns True if x is of type MaximaElement.

## EXAMPLES:

```
sage: from sage.interfaces.maxima import is_MaximaElement
sage: m = maxima(1)
sage: is_MaximaElement(m)
True
sage: is_MaximaElement(1)
False
```

sage.interfaces.maxima.**reduce\_load\_Maxima** ()

Unpickle a Maxima Pexpect interface.

## EXAMPLES:

```
sage: from sage.interfaces.maxima import reduce_load_Maxima
sage: reduce_load_Maxima()
Maxima
```

sage.interfaces.maxima.**reduce\_load\_Maxima\_function** (*parent, defn, args, latex*)

Unpickle a Maxima function.

EXAMPLES:

```
sage: from sage.interfaces.maxima import reduce_load_Maxima_function
sage: f = maxima.function('x,y','sin(x+y)')
sage: _, args = f.__reduce__()
sage: g = reduce_load_Maxima_function(*args)
sage: g == f
True
```

## ABSTRACT INTERFACE TO MAXIMA

Maxima is a free GPL'd general purpose computer algebra system whose development started in 1968 at MIT. It contains symbolic manipulation algorithms, as well as implementations of special functions, including elliptic functions and generalized hypergeometric functions. Moreover, Maxima has implementations of many functions relating to the invariant theory of the symmetric group  $S_n$ . (However, the commands for group invariants, and the corresponding Maxima documentation, are in French.) For many links to Maxima documentation see <http://maxima.sourceforge.net/docs.shtml/>.

### AUTHORS:

- William Stein (2005-12): Initial version
- David Joyner: Improved documentation
- William Stein (2006-01-08): Fixed bug in parsing
- William Stein (2006-02-22): comparisons (following suggestion of David Joyner)
- William Stein (2006-02-24): *greatly* improved robustness by adding sequence numbers to IO bracketing in `_eval_line`
- Robert Bradshaw, Nils Bruin, Jean-Pierre Flori (2010,2011): Binary library interface

This is an abstract class implementing the functions shared between the Pexpect and library interfaces to Maxima.

**class** `sage.interfaces.maxima_abstract.MaximaAbstract` (*name='maxima\_abstract'*)

Bases: `sage.interfaces.interface.Interface`

Abstract interface to Maxima.

### INPUT:

- *name* - string

OUTPUT: the interface

### EXAMPLES:

This class should not be instantiated directly, but through its subclasses `Maxima` (Pexpect interface) or `MaximaLib` (library interface):

```
sage: m = Maxima()
sage: from sage.interfaces.maxima_abstract import MaximaAbstract
sage: isinstance(m, MaximaAbstract)
True
```

**chdir** (*dir*)

Change Maxima's current working directory.

### INPUT:

- dir - string

OUTPUT: none

EXAMPLES:

```
sage: maxima.chdir('/')
```

**completions** (*s*, *verbose=True*)

Return all commands that complete the command starting with the string *s*. This is like typing *s*[tab] in the Maxima interpreter.

INPUT:

- s - string
- verbose - boolean (default: True)

OUTPUT: array of strings

EXAMPLES:

```
sage: sorted(maxima.completions('gc', verbose=False))
['gcd', 'gcdex', 'gcfactor', 'gctime']
```

**console** ()

Start the interactive Maxima console. This is a completely separate maxima session from this interface. To interact with this session, you should instead use `maxima.interact()`.

INPUT: none

OUTPUT: none

EXAMPLES:

```
sage: maxima.console()           # not tested (since we can't)
Maxima 5.34.1 http://maxima.sourceforge.net
Using Lisp ECL 13.5.1
Distributed under the GNU Public License. See the file COPYING.
Dedicated to the memory of William Schelter.
This is a development version of Maxima. The function bug_report()
provides bug reporting information.
(%i1)

sage: maxima.interact()          # this is not tested either
--> Switching to Maxima <--
maxima: 2+2
4
maxima:
--> Exiting back to Sage <--
```

**cputime** (*t=None*)

Returns the amount of CPU time that this Maxima session has used.

INPUT:

- t - float (default: None); If var{t} is not None, then it returns the difference between the current CPU time and var{t}.

OUTPUT: float

**EXAMPLES:** `sage: t = maxima.cputime()` `sage: _ = maxima.de_solve('diff(y,x,2) + 3*x = y', ['x','y'], [1,1,1])` `sage: maxima.cputime(t)` # output random 0.568913

**de\_solve** (*de*, *vars*, *ics*=None)

Solves a 1st or 2nd order ordinary differential equation (ODE) in two variables, possibly with initial conditions.

INPUT:

- **de** - a string representing the ODE
- **vars** - a list of strings representing the two variables.
- **ics** - a triple of numbers [a,b1,b2] representing  $y(a)=b1$ ,  $y'(a)=b2$

EXAMPLES:

```
sage: maxima.de_solve('diff(y,x,2) + 3*x = y', ['x','y'], [1,1,1])
y=3*x-2*e^(x-1)
sage: maxima.de_solve('diff(y,x,2) + 3*x = y', ['x','y'])
y=%k1*e^x+%k2*e^-x+3*x
sage: maxima.de_solve('diff(y,x) + 3*x = y', ['x','y'])
y=(%c-3*(-x-1))*e^-x)*e^x
sage: maxima.de_solve('diff(y,x) + 3*x = y', ['x','y'], [1,1])
y=-e^-1*(5*e^x-3*e*x-3*e)
```

**de\_solve\_laplace** (*de*, *vars*, *ics*=None)

Solves an ordinary differential equation (ODE) using Laplace transforms.

INPUT:

- **de** - a string representing the ODE (e.g., **de** = "diff(f(x),x,2)=diff(f(x),x)+sin(x)")
- **vars** - a list of strings representing the variables (e.g., **vars** = ["x","f"])
- **ics** - a list of numbers representing initial conditions, with symbols allowed which are represented by strings (eg,  $f(0)=1$ ,  $f'(0)=2$  is **ics** = [0,1,2])

EXAMPLES:

```
sage: maxima.clear('x'); maxima.clear('f')
sage: maxima.de_solve_laplace("diff(f(x),x,2) = 2*diff(f(x),x)-f(x)", ["x","f"], [0,1,2])
f(x)=x*e^x+e^x

sage: maxima.clear('x'); maxima.clear('f')
sage: f = maxima.de_solve_laplace("diff(f(x),x,2) = 2*diff(f(x),x)-f(x)", ["x","f"])
sage: f
f(x)=x*e^x*(('at('diff(f(x),x,1),x=0))-f(0)*x*e^x+f(0)*e^x
sage: print f
```

$$f(x) = x e^x \left( \frac{d}{dx} (f(x)) - f(0) \right) - f(0) x e^x + f(0) e^x$$

$$!x = 0$$

---

**Note:** The second equation sets the values of  $f(0)$  and  $f'(0)$  in Maxima, so subsequent ODEs involving these variables will have these initial conditions automatically imposed.

---

**demo** (*s*)

Run Maxima's demo for *s*.

INPUT:

- **s** - string

OUTPUT: none

EXAMPLES:

```
sage: maxima.demo('cf') # not tested
read and interpret file: ../local/share/maxima/5.34.1/demo/cf.dem
```

At the '\_' prompt, type ';' and <enter> to get next demonstration.  
frac1:cf([1,2,3,4])  
...

**describe**(s)

Return Maxima's help for s.

INPUT:

- s - string

OUTPUT:

Maxima's help for s

EXAMPLES:

```
sage: maxima.help('gcd')
-- Function: gcd (<p_1>, <p_2>, <x_1>, ...)
...
```

**example**(s)

Return Maxima's examples for s.

INPUT:

- s - string

OUTPUT:

Maxima's examples for s

EXAMPLES:

```
sage: maxima.example('arrays')
a[n]:=n*a[n-1]

a := n a
n      n - 1

a[0]:1
a[5]

120

a[n]:=n
a[6]

6

a[4]

24
done
```

**function**(args, defn, rep=None, latex=None)

Return the Maxima function with given arguments and definition.

INPUT:

- args - a string with variable names separated by commas
- defn - a string (or Maxima expression) that defines a function of the arguments in Maxima.
- rep - an optional string; if given, this is how the function will print.



OUTPUT: Maxima function

EXAMPLES:

```
sage: f = maxima.function('x', 'sin(x)')
sage: f(3.2) # abs tol 2e-16
-0.058374143427579909
sage: f = maxima.function('x,y', 'sin(x)+cos(y)')
sage: f(2, 3.5) # abs tol 2e-16
sin(2)-0.9364566872907963
sage: f
sin(x)+cos(y)

sage: g = f.integrate('z')
sage: g
(cos(y)+sin(x))*z
sage: g(1,2,3)
3*(cos(2)+sin(1))
```

The function definition can be a Maxima object:

```
sage: an_expr = maxima('sin(x)*gamma(x)')
sage: t = maxima.function('x', an_expr)
sage: t
gamma(x)*sin(x)
sage: t(2)
sin(2)
sage: float(t(2))
0.9092974268256817
sage: loads(t.dumps())
gamma(x)*sin(x)
```

**help**(s)

Return Maxima's help for s.

INPUT:

- s - string

OUTPUT:

Maxima's help for s

EXAMPLES:

```
sage: maxima.help('gcd')
-- Function: gcd (<p_1>, <p_2>, <x_1>, ...)
...
```

**plot2d**(\*args)

Plot a 2d graph using Maxima / gnuplot.

```
maxima.plot2d(f, '[var, min, max]', options)
```

INPUT:

- f** - a string representing a function (such as f="sin(x)") [var, xmin, xmax]
- options** - an optional string representing plot2d options in gnuplot format

EXAMPLES:

```

sage: maxima.plot2d('sin(x)', '[x,-5,5]') # not tested
sage: opts = '[gnuplot_term, ps], [gnuplot_out_file, "sin-plot.eps"]'
sage: maxima.plot2d('sin(x)', '[x,-5,5]', opts) # not tested

```

The eps file is saved in the current directory.

**plot2d\_parametric** (*r*, *var*, *trange*, *nticks*=50, *options*=None)

Plot  $r = [x(t), y(t)]$  for  $t = tmin...tmax$  using gnuplot with options.

INPUT:

- **r** - a string representing a function (such as `r="[x(t),y(t)]"`)
- **var** - a string representing the variable (such as `var="t"`)
- **trange** - `[tmin, tmax]` are numbers with `tmin < tmax`
- **nticks** - int (default: 50)
- **options** - an optional string representing plot2d options in gnuplot format

EXAMPLES:

```

sage: maxima.plot2d_parametric(["sin(t)", "cos(t)"], "t", [-3.1, 3.1]) # not tested

```

```

sage: opts = '[gnuplot_preamble, "set nokey"], [gnuplot_term, ps], [gnuplot_out_file, "circle.eps"]'

```

```

sage: maxima.plot2d_parametric(["sin(t)", "cos(t)"], "t", [-3.1, 3.1], options=opts) # not tested

```

The eps file is saved to the current working directory.

Here is another fun plot:

```

sage: maxima.plot2d_parametric(["sin(5*t)", "cos(11*t)"], "t", [0, 2*pi()], nticks=400) # not tested

```

**plot3d** (\*args)

Plot a 3d graph using Maxima / gnuplot.

`maxima.plot3d(f, '[x, xmin, xmax]', '[y, ymin, ymax]', '[grid, nx, ny]', options)`

INPUT:

- **f** - a string representing a function (such as `f="sin(x)"`) [*var*, *min*, *max*]
- **args** should be of the form `'[x, xmin, xmax]', '[y, ymin, ymax]', '[grid, nx, ny]', options`

EXAMPLES:

```

sage: maxima.plot3d('1 + x^3 - y^2', '[x,-2,2]', '[y,-2,2]', '[grid,12,12]') # not tested

```

```

sage: maxima.plot3d('sin(x)*cos(y)', '[x,-2,2]', '[y,-2,2]', '[grid,30,30]') # not tested

```

```

sage: opts = '[gnuplot_term, ps], [gnuplot_out_file, "sin-plot.eps"]'

```

```

sage: maxima.plot3d('sin(x+y)', '[x,-5,5]', '[y,-1,1]', opts) # not tested

```

The eps file is saved in the current working directory.

**plot3d\_parametric** (*r*, *vars*, *urange*, *vrange*, *options*=None)

Plot a 3d parametric graph with  $r=(x,y,z)$ ,  $x = x(u,v)$ ,  $y = y(u,v)$ ,  $z = z(u,v)$ , for  $u = u_{min}...u_{max}$ ,  $v = v_{min}...v_{max}$  using gnuplot with options.

INPUT:

- **x, y, z** - a string representing a function (such as `x="u2+v2", ...`) *vars* is a list or two strings representing variables (such as `vars=["u","v"]`)
- **urange** - `[umin, umax]`

•**vrange** - [vmin, vmax] are lists of numbers with  $u_{\min}$   $u_{\max}$ ,  $v_{\min}$   $v_{\max}$

•**options** - optional string representing **plot2d** options in gnuplot format

OUTPUT: displays a plot on screen or saves to a file

EXAMPLES:

```
sage: maxima.plot3d_parametric(["v*sin(u)", "v*cos(u)", "v"], ["u", "v"], [-3.2, 3.2], [0, 3])
sage: opts = '[gnuplot_term, ps], [gnuplot_out_file, "sin-cos-plot.eps"]'
sage: maxima.plot3d_parametric(["v*sin(u)", "v*cos(u)", "v"], ["u", "v"], [-3.2, 3.2], [0, 3], opts)
```

The eps file is saved in the current working directory.

Here is a torus:

```
sage: _ = maxima.eval("expr_1: cos(y)*(10.0+6*cos(x)); expr_2: sin(y)*(10.0+6*cos(x)); expr_3: 10.0")
sage: maxima.plot3d_parametric(["expr_1", "expr_2", "expr_3"], ["x", "y"], [0, 6], [0, 6]) # not tested
```

Here is a Mobius strip:

```
sage: x = "cos(u)*(3 + v*cos(u/2))"
sage: y = "sin(u)*(3 + v*cos(u/2))"
sage: z = "v*sin(u/2)"
sage: maxima.plot3d_parametric([x, y, z], ["u", "v"], [-3.1, 3.2], [-1/10, 1/10]) # not tested
```

**plot\_list** (ptsx, ptsy, options=None)

Plots a curve determined by a sequence of points.

INPUT:

•**ptsx** - [x1,...,xn], where the xi and yi are real,

•**ptsy** - [y1,...,yn]

•**options** - a string representing maxima **plot2d** options.

The points are (x1,y1), (x2,y2), etc.

This function requires maxima 5.9.2 or newer.

---

**Note:** More than 150 points can sometimes lead to the program hanging. Why?

---

EXAMPLES:

```
sage: zeta_ptsx = [ (pari(1/2 + i*I/10).zeta().real()).precision(1) for i in range (70,150)]
sage: zeta_ptsy = [ (pari(1/2 + i*I/10).zeta().imag()).precision(1) for i in range (70,150)]
sage: maxima.plot_list(zeta_ptsx, zeta_ptsy) # not tested
sage: opts='[gnuplot_preamble, "set nokey"], [gnuplot_term, ps], [gnuplot_out_file, "zeta.eps"]
sage: maxima.plot_list(zeta_ptsx, zeta_ptsy, opts) # not tested
```

**plot\_multilist** (pts\_list, options=None)

Plots a list of list of points pts\_list=[pts1,pts2,...,ptsn], where each ptsi is of the form [[x1,y1],...,[xn,yn]]  
x's must be integers and y's reals options is a string representing maxima **plot2d** options.

INPUT:

•**pts\_list** - list of points; each point must be of the form [x,y] where x is an integer and y is a real

•**var** - string; representing Maxima's **plot2d** options

Requires maxima 5.9.2 at least.

---

**Note:** More than 150 points can sometimes lead to the program hanging.

---

## EXAMPLES:

```

sage: xx = [ i/10.0 for i in range (-10,10)]
sage: yy = [ i/10.0 for i in range (-10,10)]
sage: x0 = [ 0 for i in range (-10,10)]
sage: y0 = [ 0 for i in range (-10,10)]
sage: zeta_ptsx1 = [ (pari(1/2+i*I/10).zeta().real()).precision(1) for i in range (10)]
sage: zeta_ptsy1 = [ (pari(1/2+i*I/10).zeta().imag()).precision(1) for i in range (10)]
sage: maxima.plot_multilist([[zeta_ptsx1,zeta_ptsy1],[xx,y0],[x0,yy]]) # not tested
sage: zeta_ptsx1 = [ (pari(1/2+i*I/10).zeta().real()).precision(1) for i in range (10,150)]
sage: zeta_ptsy1 = [ (pari(1/2+i*I/10).zeta().imag()).precision(1) for i in range (10,150)]
sage: maxima.plot_multilist([[zeta_ptsx1,zeta_ptsy1],[xx,y0],[x0,yy]]) # not tested
sage: opts='[gnuplot_preamble, "set nokey"]'
sage: maxima.plot_multilist([[zeta_ptsx1,zeta_ptsy1],[xx,y0],[x0,yy]],opts) # not tested

```

**solve\_linear** (*eqns*, *vars*)

Wraps maxima's linsolve.

## INPUT:

- *eqns* - a list of *m* strings; each representing a linear question in  $m = n$  variables
- *vars* - a list of *n* strings; each representing a variable

## EXAMPLES:

```

sage: eqns = ["x + z = y", "2*a*x - y = 2*a^2", "y - 2*z = 2"]
sage: vars = ["x", "y", "z"]
sage: maxima.solve_linear(eqns, vars)
[x=a+1, y=2*a, z=a-1]

```

**trait\_names** (*verbose=True*, *use\_disk\_cache=True*)

Return all Maxima commands, which is useful for tab completion.

## INPUT:

- *verbose* - boolean (default: True)
- *use\_disk\_cache* - boolean (default: True); if set to True, try to read cached result from disk

OUTPUT: array of strings

## EXAMPLES:

```

sage: t = maxima.trait_names(verbose=False)
sage: 'gcd' in t
True
sage: len(t) # random output
1840

```

**unit\_quadratic\_integer** (*n*)

Finds a unit of the ring of integers of the quadratic number field  $\mathbf{Q}(\sqrt{n})$ ,  $n > 1$ , using the qunit maxima command.

## INPUT:

- *n* - an integer

## EXAMPLES:

```

sage: u = maxima.unit_quadratic_integer(101); u
a + 10
sage: u.parent()

```

```

Number Field in a with defining polynomial x^2 - 101
sage: u = maxima.unit_quadratic_integer(13)
sage: u
5*a + 18
sage: u.parent()
Number Field in a with defining polynomial x^2 - 13

```

**version()**

Return the version of Maxima that Sage includes.

INPUT: none

OUTPUT: none

EXAMPLES:

```

sage: maxima.version()
'5.35.1'

```

```

class sage.interfaces.maxima_abstract.MaximaAbstractElement (parent, value,
                                                             is_name=False,
                                                             name=None)

```

Bases: `sage.interfaces.interface.InterfaceElement`

Element of Maxima through an abstract interface.

EXAMPLES:

Elements of this class should not be created directly. The targeted parent of a concrete inherited class should be used instead:

```

sage: from sage.interfaces.maxima_lib import maxima_lib
sage: xp = maxima(x)
sage: type(xp)
<class 'sage.interfaces.maxima.MaximaElement'>
sage: x1 = maxima_lib(x)
sage: type(x1)
<class 'sage.interfaces.maxima_lib.MaximaLibElement'>

```

**bool()**

Convert self into a boolean.

INPUT: none

OUTPUT: boolean

EXAMPLES:

```

sage: maxima(0).bool()
False
sage: maxima(1).bool()
True

```

**comma(args)**

Form the expression that would be written 'self, args' in Maxima.

INPUT:

- args - string

OUTPUT: Maxima object

EXAMPLES:

```
sage: maxima('sqrt(2) + I').comma('numer')
I+1.41421356237309...
sage: maxima('sqrt(2) + I*a').comma('a=5')
5*I+sqrt(2)
```

**derivative** (*var='x', n=1*)

Return the n-th derivative of self.

INPUT:

- var - variable (default: 'x')
- n - integer (default: 1)

OUTPUT: n-th derivative of self with respect to the variable var

EXAMPLES:

```
sage: f = maxima('x^2')
sage: f.diff()
2*x
sage: f.diff('x')
2*x
sage: f.diff('x', 2)
2
sage: maxima('sin(x^2)').diff('x', 4)
16*x^4*sin(x^2)-12*sin(x^2)-48*x^2*cos(x^2)

sage: f = maxima('x^2 + 17*y^2')
sage: f.diff('x')
34*y*diff(y,x,1)+2*x
sage: f.diff('y')
34*y
```

**diff** (*var='x', n=1*)

Return the n-th derivative of self.

INPUT:

- var - variable (default: 'x')
- n - integer (default: 1)

OUTPUT: n-th derivative of self with respect to the variable var

EXAMPLES:

```
sage: f = maxima('x^2')
sage: f.diff()
2*x
sage: f.diff('x')
2*x
sage: f.diff('x', 2)
2
sage: maxima('sin(x^2)').diff('x', 4)
16*x^4*sin(x^2)-12*sin(x^2)-48*x^2*cos(x^2)

sage: f = maxima('x^2 + 17*y^2')
sage: f.diff('x')
34*y*diff(y,x,1)+2*x
sage: f.diff('y')
34*y
```

**dot** (*other*)Implements the notation `self . other`.

INPUT:

- `other` - matrix; argument to `dot`.

OUTPUT: Maxima matrix

EXAMPLES:

```
sage: A = maxima('matrix ([a1],[a2])')
sage: B = maxima('matrix ([b1, b2])')
sage: A.dot(B)
matrix([a1*b1, a1*b2], [a2*b1, a2*b2])
```

**imag** ()

Return the imaginary part of this Maxima element.

INPUT: none

OUTPUT: Maxima real

EXAMPLES:

```
sage: maxima('2 + (2/3)*%i').imag()
2/3
```

**integral** (*var='x', min=None, max=None*)Return the integral of `self` with respect to the variable `x`.

INPUT:

- `var` - variable
- `min` - default: None
- `max` - default: None

OUTPUT:

- the definite integral if `xmin` is not None
- an indefinite integral otherwise

EXAMPLES:

```
sage: maxima('x^2+1').integral()
x^3/3+x
sage: maxima('x^2+ 1 + y^2').integral('y')
y^3/3+x^2*y+y
sage: maxima('x / (x^2+1)').integral()
log(x^2+1)/2
sage: maxima('1/(x^2+1)').integral()
atan(x)
sage: maxima('1/(x^2+1)').integral('x', 0, infinity)
%pi/2
sage: maxima('x/(x^2+1)').integral('x', -1, 1)
0

sage: f = maxima('exp(x^2)').integral('x',0,1); f
-sqrt(%pi)*%i*erf(%i)/2
sage: f.numer()
1.46265174590718...
```

**integrate** (*var='x', min=None, max=None*)

Return the integral of self with respect to the variable x.

INPUT:

- var - variable
- min - default: None
- max - default: None

OUTPUT:

- the definite integral if xmin is not None
- an indefinite integral otherwise

EXAMPLES:

```
sage: maxima('x^2+1').integral()
x^3/3+x
sage: maxima('x^2+ 1 + y^2').integral('y')
y^3/3+x^2*y+y
sage: maxima('x / (x^2+1)').integral()
log(x^2+1)/2
sage: maxima('1/(x^2+1)').integral()
atan(x)
sage: maxima('1/(x^2+1)').integral('x', 0, infinity)
%pi/2
sage: maxima('x/(x^2+1)').integral('x', -1, 1)
0

sage: f = maxima('exp(x^2)').integral('x',0,1); f
-sqrt(%pi)*%i*erf(%i)/2
sage: f.numer()
1.46265174590718...
```

**nintegral** (*var='x', a=0, b=1, desired\_relative\_error='1e-8', maximum\_num\_subintervals=200*)

Return a numerical approximation to the integral of self from a to b.

INPUT:

- var - variable to integrate with respect to
- a - lower endpoint of integration
- b - upper endpoint of integration
- desired\_relative\_error** - (default: '1e-8') the desired relative error
- maximum\_num\_subintervals** - (default: 200) maxima number of subintervals

OUTPUT:

- approximation to the integral
- estimated absolute error of the** approximation
- the number of integrand evaluations
- an error code:
  - 0 - no problems were encountered
  - 1 - too many subintervals were done
  - 2 - excessive roundoff error



- 3 - extremely bad integrand behavior
- 4 - failed to converge
- 5 - integral is probably divergent or slowly convergent
- 6 - the input is invalid

EXAMPLES:

```
sage: maxima('exp(-sqrt(x))').nintegral('x', 0, 1)
(0.5284822353142306, 4.16331413788384...e-11, 231, 0)
```

Note that GP also does numerical integration, and can do so to very high precision very quickly:

```
sage: gp('intnum(x=0,1,exp(-sqrt(x)))')
0.5284822353142307136179049194          # 32-bit
0.52848223531423071361790491935415653022 # 64-bit
sage: _ = gp.set_precision(80)
sage: gp('intnum(x=0,1,exp(-sqrt(x)))')
0.52848223531423071361790491935415653021675547587292866196865279321015401702040079
```

**numer()**

Return numerical approximation to self as a Maxima object.

INPUT: none

OUTPUT: Maxima object

EXAMPLES:

```
sage: a = maxima('sqrt(2)').numer(); a
1.41421356237309...
sage: type(a)
<class 'sage.interfaces.maxima.MaximaElement'>
```

**partial\_fraction\_decomposition(var='x')**

Return the partial fraction decomposition of self with respect to the variable var.

INPUT:

- var - string

OUTPUT: Maxima object

EXAMPLES:

```
sage: f = maxima('1/((1+x)*(x-1))')
sage: f.partial_fraction_decomposition('x')
1/(2*(x-1))-1/(2*(x+1))
sage: print f.partial_fraction_decomposition('x')
      1          1
----- - -----
  2 (x - 1)    2 (x + 1)
```

**real()**

Return the real part of this Maxima element.

INPUT: none

OUTPUT: Maxima real

EXAMPLES:

```
sage: maxima('2 + (2/3)*%i').real()
2
```

**str()**  
Return string representation of this Maxima object.

INPUT: none

OUTPUT: string

EXAMPLES:

```
sage: maxima('sqrt(2) + 1/3').str()
'sqrt(2)+1/3'
```

**subst(val)**  
Substitute a value or several values into this Maxima object.

INPUT:

- val - string representing substitution(s) to perform

OUTPUT: Maxima object

EXAMPLES:

```
sage: maxima('a^2 + 3*a + b').subst('b=2')
a^2+3*a+2
```

```
sage: maxima('a^2 + 3*a + b').subst('a=17')
b+340
```

```
sage: maxima('a^2 + 3*a + b').subst('a=17, b=2')
342
```

**trait\_names(verbose=False)**  
Return all Maxima commands, which is useful for tab completion.

INPUT:

- verbose - boolean

OUTPUT: list of strings

EXAMPLES:

```
sage: m = maxima(2)
```

```
sage: 'gcd' in m.trait_names()
True
```

```
class sage.interfaces.maxima_abstract.MaximaAbstractElementFunction(parent,
                                                                    name, defn,
                                                                    args, latex)
```

Bases: `sage.interfaces.maxima_abstract.MaximaAbstractElement`

Create a Maxima function with the parent `parent`, name `name`, definition `defn`, arguments `args` and latex representation `latex`.

INPUT:

- parent - an instance of a concrete Maxima interface
- name - string
- defn - string
- args - string; comma separated names of arguments

- `latex` - string

OUTPUT: Maxima function

EXAMPLES:

```
sage: maxima.function('x,y','e^cos(x)')
e^cos(x)
```

**arguments** (*split=True*)

Returns the arguments of this Maxima function.

INPUT:

- `split` - boolean; if True return a tuple of strings, otherwise return a string of comma-separated arguments

OUTPUT:

- a string if `split` is False
- a list of strings if `split` is True

EXAMPLES:

```
sage: f = maxima.function('x,y','sin(x+y)')
sage: f.arguments()
['x', 'y']
sage: f.arguments(split=False)
'x,y'
sage: f = maxima.function('', 'sin(x)')
sage: f.arguments()
[]
```

**definition** ()

Returns the definition of this Maxima function as a string.

INPUT: none

OUTPUT: string

EXAMPLES:

```
sage: f = maxima.function('x,y','sin(x+y)')
sage: f.definition()
'sin(x+y)'
```

**integral** (*var*)

Returns the integral of self with respect to the variable `var`.

INPUT:

- `var` - a variable

OUTPUT: Maxima function

Note that `integrate` is an alias of `integral`.

EXAMPLES:

```
sage: x,y = var('x,y')
sage: f = maxima.function('x','sin(x)')
sage: f.integral(x)
-cos(x)
sage: f.integral(y)
sin(x)*y
```

**integrate** (*var*)

Returns the integral of self with respect to the variable *var*.

INPUT:

- *var* - a variable

OUTPUT: Maxima function

Note that `integrate` is an alias of `integral`.

EXAMPLES:

```
sage: x,y = var('x,y')
sage: f = maxima.function('x','sin(x)')
sage: f.integral(x)
-cos(x)
sage: f.integral(y)
sin(x)*y
```

**class** `sage.interfaces.maxima_abstract.MaximaAbstractFunction` (*parent, name*)

Bases: `sage.interfaces.interface.InterfaceFunction`

**class** `sage.interfaces.maxima_abstract.MaximaAbstractFunctionElement` (*obj, name*)

Bases: `sage.interfaces.interface.InterfaceFunctionElement`

`sage.interfaces.maxima_abstract.maxima_console()`

Spawn a new Maxima command-line session.

EXAMPLES:

```
sage: from sage.interfaces.maxima_abstract import maxima_console
sage: maxima_console() # not tested
Maxima 5.34.1 http://maxima.sourceforge.net
...
```

`sage.interfaces.maxima_abstract.maxima_version()`

Return Maxima version.

Currently this calls a new copy of Maxima.

EXAMPLES:

```
sage: from sage.interfaces.maxima_abstract import maxima_version
sage: maxima_version()
'5.35.1'
```

`sage.interfaces.maxima_abstract.reduce_load_MaximaAbstract_function` (*parent,*  
*defn,*  
*args,*  
*latex*)

Unpickle a Maxima function.

EXAMPLES:

```
sage: from sage.interfaces.maxima_abstract import reduce_load_MaximaAbstract_function
sage: f = maxima.function('x,y','sin(x+y)')
sage: _, args = f.__reduce__()
sage: g = reduce_load_MaximaAbstract_function(*args)
sage: g == f
True
```

## LIBRARY INTERFACE TO MAXIMA

Maxima is a free GPL'd general purpose computer algebra system whose development started in 1968 at MIT. It contains symbolic manipulation algorithms, as well as implementations of special functions, including elliptic functions and generalized hypergeometric functions. Moreover, Maxima has implementations of many functions relating to the invariant theory of the symmetric group  $S_n$ . (However, the commands for group invariants, and the corresponding Maxima documentation, are in French.) For many links to Maxima documentation, see <http://maxima.sourceforge.net/documentation.html>.

### AUTHORS:

- William Stein (2005-12): Initial version
- David Joyner: Improved documentation
- William Stein (2006-01-08): Fixed bug in parsing
- William Stein (2006-02-22): comparisons (following suggestion of David Joyner)
- William Stein (2006-02-24): *greatly* improved robustness by adding sequence numbers to IO bracketing in `_eval_line`
- Robert Bradshaw, Nils Bruin, Jean-Pierre Flori (2010,2011): Binary library interface

For this interface, Maxima is loaded into ECL which is itself loaded as a C library in Sage. Translations between Sage and Maxima objects (which are nothing but wrappers to ECL objects) is made as much as possible directly, but falls back to the string based conversion used by the classical Maxima Pexpect interface in case no new implementation has been made.

This interface is the one used for calculus by Sage and is accessible as *maxima\_calculus*:

```
sage: maxima_calculus
Maxima_lib
```

Only one instance of this interface can be instantiated, so the user should not try to instantiate another one, which is anyway set to raise an error:

```
sage: from sage.interfaces.maxima_lib import MaximaLib
sage: MaximaLib()
Traceback (most recent call last):
...
RuntimeError: Maxima interface in library mode can only be instantiated once
```

```
class sage.interfaces.maxima_lib.MaximaLib
    Bases: sage.interfaces.maxima_abstract.MaximaAbstract
    Interface to Maxima as a Library.
    INPUT: none
```

OUTPUT: Maxima interface as a Library

EXAMPLES:

```
sage: from sage.interfaces.maxima_lib import MaximaLib, maxima_lib
sage: isinstance(maxima_lib, MaximaLib)
True
```

Only one such interface can be instantiated:

```
sage: MaximaLib()
Traceback (most recent call last):
...
RuntimeError: Maxima interface in library mode can only
be instantiated once
```

**clear** (*var*)

Clear the variable named *var*.

INPUT:

- *var* - string

OUTPUT: none

EXAMPLES:

```
sage: from sage.interfaces.maxima_lib import maxima_lib
sage: maxima_lib.set('xxxxx', '2')
sage: maxima_lib.get('xxxxx')
'2'
sage: maxima_lib.clear('xxxxx')
sage: maxima_lib.get('xxxxx')
'xxxxx'
```

**eval** (*line*, *locals=None*, *reformat=True*, *\*\*kws*)

Evaluate the line in Maxima.

INPUT:

- *line* - string; text to evaluate
- *locals* - None (ignored); this is used for compatibility with the Sage notebook's generic system interface.
- *reformat* - boolean; whether to strip output or not
- *\*\*kws* - All other arguments are currently ignored.

OUTPUT: string representing Maxima output

EXAMPLES:

```
sage: from sage.interfaces.maxima_lib import maxima_lib
sage: maxima_lib._eval_line('1+1')
'2'
sage: maxima_lib._eval_line('1+1;')
'2'
sage: maxima_lib._eval_line('1+1$')
''
sage: maxima_lib._eval_line('randvar : cos(x)+sin(y)$')
''
sage: maxima_lib._eval_line('randvar')
'sin(y)+cos(x)'
```

**get** (*var*)Get the string value of the variable *var*.

INPUT:

•*var* - string

OUTPUT: string

EXAMPLES:

```
sage: from sage.interfaces.maxima_lib import maxima_lib
sage: maxima_lib.set('xxxxx', '2')
sage: maxima_lib.get('xxxxx')
'2'
```

**lisp** (*cmd*)

Send a lisp command to maxima.

INPUT:

•*cmd* - string

OUTPUT: ECL object

---

**Note:** The output of this command is very raw - not pretty.

---

EXAMPLES:

```
sage: from sage.interfaces.maxima_lib import maxima_lib
sage: maxima_lib.lisp("(+ 2 17)")
<ECL: 19>
```

**set** (*var*, *value*)Set the variable *var* to the given value.

INPUT:

•*var* - string•*value* - string

OUTPUT: none

EXAMPLES:

```
sage: from sage.interfaces.maxima_lib import maxima_lib
sage: maxima_lib.set('xxxxx', '2')
sage: maxima_lib.get('xxxxx')
'2'
```

**sr\_integral** (*\*args*)

Helper function to wrap calculus use of Maxima's integration.

TESTS:

```
sage: a,b=var('a,b')
sage: integrate(1/(x^3 * (a+b*x)^(1/3)), x)
Traceback (most recent call last):
...
ValueError: Computation failed since Maxima requested additional
constraints; using the 'assume' command before evaluation
*may* help (example of legal syntax is 'assume(a>0)', see
'assume?' for more details)
```

```

Is a positive or negative?
sage: assume(a>0)
sage: integrate(1/(x^3 * (a+b*x)^(1/3)), x)
2/9*sqrt(3)*b^2*arctan(1/3*sqrt(3)*(2*(b*x + a)^(1/3) + a^(1/3))/a^(1/3))/a^(7/3) - 1/9*b^2*
sage: var('x, n')
(x, n)
sage: integral(x^n, x)
Traceback (most recent call last):
...
ValueError: Computation failed since Maxima requested additional
constraints; using the 'assume' command before evaluation
*may* help (example of legal syntax is 'assume(n>0)',
see 'assume?' for more details)
Is n equal to -1?
sage: assume(n+1>0)
sage: integral(x^n, x)
x^(n + 1)/(n + 1)
sage: forget()
sage: assumptions() # Check the assumptions really were forgotten
[]

```

Make sure the `abs_integrate` package is being used, [trac ticket #11483](#). The following are examples from the Maxima `abs_integrate` documentation:

```

sage: integrate(abs(x), x)
1/2*x*abs(x)

sage: integrate(sgn(x) - sgn(1-x), x)
abs(x - 1) + abs(x)

sage: integrate(1 / (1 + abs(x-5)), x, -5, 6)
log(11) + log(2)

sage: integrate(1/(1 + abs(x)), x)
1/2*(log(x + 1) + log(-x + 1))*sgn(x) + 1/2*log(x + 1) - 1/2*log(-x + 1)

sage: integrate(cos(x + abs(x)), x)
-1/2*x*sgn(x) + 1/4*(sgn(x) + 1)*sin(2*x) + 1/2*x

```

The last example relies on the following simplification:

```

sage: maxima("realpart(signum(x))")
signum(x)

```

An example from sage-support thread [e641001f8b8d1129](#):

```

sage: f = e^(-x^2/2)/sqrt(2*pi) * sgn(x-1)
sage: integrate(f, x, -Infinity, Infinity)
-erf(1/2*sqrt(2))

```

From [trac ticket #8624](#):

```

sage: integral(abs(cos(x))*sin(x), (x, pi/2, pi))
1/2

sage: integrate(sqrt(x + sqrt(x)), x).canonicalize_radical()
1/12*((8*x - 3)*x^(1/4) + 2*x^(3/4))*sqrt(sqrt(x) + 1) + 1/8*log(sqrt(sqrt(x) + 1) + x^(1/4))

```

And [trac ticket #11594](#):



```
sage: integrate(abs(x^2 - 1), x, -2, 2)
4
```

This definite integral returned zero (incorrectly) in at least Maxima 5.23. The correct answer is now given ([trac ticket #11591](#)):

```
sage: f = (x^2)*exp(x) / (1+exp(x))^2
sage: integrate(f, (x, -infinity, infinity))
1/3*pi^2
```

Sometimes one needs different simplification settings, such as `radexpand`, to compute an integral (see [trac ticket #10955](#)):

```
sage: f = sqrt(x + 1/x^2)
sage: maxima = sage.calculus.calculus.maxima
sage: maxima('radexpand')
true
sage: integrate(f, x)
integrate(sqrt(x + 1/x^2), x)
sage: maxima('radexpand: all')
all
sage: g = integrate(f, x); g
2/3*sqrt(x^3 + 1) - 1/3*log(sqrt(x^3 + 1) + 1) + 1/3*log(sqrt(x^3 + 1) - 1)
sage: (f - g.diff(x)).canonicalize_radical()
0
sage: maxima('radexpand: true')
true
```

The following integral was computed incorrectly in versions of Maxima before 5.27 (see [trac ticket #12947](#)):

```
sage: a = integrate(x*cos(x^3), (x, 0, 1/2)).n()
sage: a.real()
0.124756040961038
sage: a.imag().abs() < 3e-17
True
```

**sr\_limit** (*expr*, *v*, *a*, *dir=None*)

Helper function to wrap calculus use of Maxima's limits.

TESTS:

```
sage: f = (1+1/x)^x
sage: limit(f, x = oo)
e
sage: limit(f, x = 5)
7776/3125
sage: limit(f, x = 1.2)
2.06961575467...
sage: var('a')
```

*a*

```
sage: limit(x^a, x=0)
```

```
Traceback (most recent call last):
```

```
...
```

```
ValueError: Computation failed since Maxima requested additional
constraints; using the 'assume' command before evaluation
*may* help (example of legal syntax is 'assume(a>0)', see 'assume?'
for more details)
Is a positive, negative or zero?
sage: assume(a>0)
```

```

sage: limit(x^a,x=0)
Traceback (most recent call last):
...
ValueError: Computation failed ...
Is a an integer?
sage: assume(a,'integer')
sage: assume(a,'even') # Yes, Maxima will ask this too
sage: limit(x^a,x=0)
0
sage: forget()
sage: assumptions() # check the assumptions were really forgotten
[]

```

The second limit below was computed incorrectly prior to Maxima 5.24 ([trac ticket #10868](#)):

```

sage: f(n) = 2 + 1/factorial(n)
sage: limit(f(n), n=infinity)
2
sage: limit(1/f(n), n=infinity)
1/2

```

The limit below was computed incorrectly prior to Maxima 5.30 (see [trac ticket #13526](#)):

```

sage: n = var('n')
sage: l = (3^n + (-2)^n) / (3^(n+1) + (-2)^(n+1))
sage: l.limit(n=oo)
1/3

```

The following limit computation used to incorrectly return 0 or infinity, depending on the domain (see [trac ticket #15033](#)):

```

sage: m = sage.calculus.calculus.maxima
sage: _ = m.eval('domain: real') # much faster than 'domain: complex'
sage: limit(gamma(x + 1/2)/(sqrt(x)*gamma(x)), x=infinity)
1
sage: _ = m.eval('domain: complex')

```

**sr\_sum(\*args)**

Helper function to wrap calculus use of Maxima's summation.

TESTS:

Check that [trac ticket #16224](#) is fixed:

```

sage: k = var('k')
sage: sum(x^(2*k)/factorial(2*k), k, 0, oo).canonicalize_radical()
cosh(x)

```

```

sage: x, y, k, n = var('x, y, k, n')
sage: sum(binomial(n,k) * x^k * y^(n-k), k, 0, n)
(x + y)^n
sage: q, a = var('q, a')
sage: sum(a*q^k, k, 0, oo)
Traceback (most recent call last):
...
ValueError: Computation failed since Maxima requested additional
constraints; using the 'assume' command before evaluation *may* help
(example of legal syntax is 'assume(abs(q)-1>0)', see 'assume?'
for more details)
Is abs(q)-1 positive, negative or zero?

```

```

sage: assume(q > 1)
sage: sum(a*q^k, k, 0, oo)
Traceback (most recent call last):
...
ValueError: Sum is divergent.
sage: forget()
sage: assume(abs(q) < 1)
sage: sum(a*q^k, k, 0, oo)
-a/(q - 1)
sage: forget()
sage: assumptions() # check the assumptions were really forgotten
[]

```

Taking the sum of all natural numbers informs us that the sum is divergent. Maxima (before 5.29.1) used to ask questions about  $m$ , leading to a different error (see [trac ticket #11990](#)):

```

sage: m = var('m')
sage: sum(m, m, 0, infinity)
Traceback (most recent call last):
...
ValueError: Sum is divergent.

```

An error with an infinite sum in Maxima (before 5.30.0, see [trac ticket #13712](#)):

```

sage: n = var('n')
sage: sum(1/((2*n-1)^2*(2*n+1)^2*(2*n+3)^2), n, 0, oo)
3/256*pi^2

```

Maxima correctly detects division by zero in a symbolic sum (see [trac ticket #11894](#)):

```

sage: sum(1/(m^4 + 2*m^3 + 3*m^2 + 2*m)^2, m, 0, infinity)
Traceback (most recent call last):
...
RuntimeError: ECL says: Error executing code in Maxima: Zero to negative power computed.

```

Similar situation for [trac ticket #12410](#):

```

sage: x = var('x')
sage: sum(1/x*(-1)^x, x, 0, oo)
Traceback (most recent call last):
...
RuntimeError: ECL says: Error executing code in Maxima: Zero to negative power computed.

```

**sr\_tlimit** (*expr*, *v*, *a*, *dir=None*)

Helper function to wrap calculus use of Maxima's Taylor series limits.

TESTS:

```

sage: f = (1+1/x)^x
sage: limit(f, x = I, taylor=True)
(-I + 1)^I

```

**class** sage.interfaces.maxima\_lib.**MaximaLibElement** (*parent*, *value*, *is\_name=False*,  
*name=None*)

Bases: sage.interfaces.maxima\_abstract.MaximaAbstractElement

Element of Maxima through library interface.

EXAMPLES:

Elements of this class should not be created directly. The targeted parent should be used instead:

```
sage: from sage.interfaces.maxima_lib import maxima_lib
sage: maxima_lib(4)
4
sage: maxima_lib(log(x))
log(_SAGE_VAR_x)
```

**display2d** (*onscreen=True*)

Return the 2d representation of this Maxima object.

INPUT:

- *onscreen* - boolean (default: True); whether to print or return

OUTPUT:

The representation is printed if *onscreen* is set to True and returned as a string otherwise.

EXAMPLES:

```
sage: from sage.interfaces.maxima_lib import maxima_lib
sage: F = maxima_lib('x^5 - y^5').factor()
sage: F.display2d()
```

$$- (y - x) (y^4 + x y^3 + x^2 y^2 + x^3 y + x^4)$$

**ecl** ()

Return the underlying ECL object of this MaximaLib object.

INPUT: none

OUTPUT: ECL object

EXAMPLES:

```
sage: from sage.interfaces.maxima_lib import maxima_lib
sage: maxima_lib(x+cos(19)).ecl()
<ECL: ((MPLUS SIMP) ((%COS SIMP) 19) |$_SAGE_VAR_x|)>
```

**to\_poly\_solve** (*vars*, *options=''*)

Use Maxima's `to_poly_solver` package.

INPUT:

- *vars* - symbolic expressions
- *options* - string (default='')

OUTPUT: Maxima object

EXAMPLES:

The `zXXX` below are names for arbitrary integers and subject to change:

```
sage: from sage.interfaces.maxima_lib import maxima_lib
sage: sol = maxima_lib(sin(x) == 0).to_poly_solve(x)
sage: sol.sage()
[[x == pi*z54]]
```

**class** `sage.interfaces.maxima_lib.MaximaLibElementFunction` (*parent*, *name*, *defn*, *args*,  
*latex*)

Bases: `sage.interfaces.maxima_lib.MaximaLibElement`, `sage.interfaces.maxima_abstract.Maxima`

Create a Maxima function. See `MaximaAbstractElementFunction` for full documentation.

TESTS:

```
sage: from sage.interfaces.maxima_abstract import MaximaAbstractElementFunction
sage: MaximaAbstractElementFunction == loads(dumps(MaximaAbstractElementFunction))
True
sage: f = maxima.function('x,y','sin(x+y)')
sage: f == loads(dumps(f))
True
```

```
class sage.interfaces.maxima_lib.MaximaLibFunction(parent, name)
    Bases: sage.interfaces.maxima_abstract.MaximaAbstractFunction
```

```
class sage.interfaces.maxima_lib.MaximaLibFunctionElement(obj, name)
    Bases: sage.interfaces.maxima_abstract.MaximaAbstractFunctionElement
```

```
sage.interfaces.maxima_lib.dummy_integrate(expr)
```

We would like to simply tie Maxima's integrate to sage.calculus.calculus.dummy\_integrate, but we're being imported there so to avoid circularity we define it here.

INPUT:

- expr - ECL object; a Maxima %INTEGRATE expression

OUTPUT: symbolic expression

EXAMPLES:

```
sage: from sage.interfaces.maxima_lib import maxima_lib, dummy_integrate
sage: f = maxima_lib('f(x)').integrate('x')
sage: f.ecl()
<ECL: ((%INTEGRATE SIMP) (($F SIMP) $X) $X)>
sage: dummy_integrate(f.ecl())
integrate(f(x), x)

sage: f = maxima_lib('f(x)').integrate('x', 0, 10)
sage: f.ecl()
<ECL: ((%INTEGRATE SIMP) (($F SIMP) $X) $X 0 10)>
sage: dummy_integrate(f.ecl())
integrate(f(x), x, 0, 10)
```

```
sage.interfaces.maxima_lib.is_MaximaLibElement(x)
```

Returns True if x is of type MaximaLibElement.

EXAMPLES:

```
sage: from sage.interfaces.maxima_lib import maxima_lib, is_MaximaLibElement
sage: m = maxima_lib(1)
sage: is_MaximaLibElement(m)
True
sage: is_MaximaLibElement(1)
False
```

```
sage.interfaces.maxima_lib.max_at_to_sage(expr)
```

Special conversion rule for AT expressions.

INPUT:

- expr - ECL object; a Maxima AT expression

OUTPUT: symbolic expression

EXAMPLES:

```
sage: from sage.interfaces.maxima_lib import maxima_lib, max_at_to_sage
sage: a=maxima_lib("'at(f(x,y,z),[x=1,y=2,z=3])")
sage: a
'at(f(x,y,z),[x=1,y=2,z=3])
sage: max_at_to_sage(a.ecl())
f(1, 2, 3)
sage: a=maxima_lib("'at(f(x,y,z),x=1)")
sage: a
'at(f(x,y,z),x=1)
sage: max_at_to_sage(a.ecl())
f(1, y, z)
```

`sage.interfaces.maxima_lib.max_to_sr(expr)`

Convert a Maxima object into a symbolic expression.

INPUT:

- expr - ECL object

OUTPUT: symbolic expression

EXAMPLES:

```
sage: from sage.interfaces.maxima_lib import maxima_lib, max_to_sr
sage: f = maxima_lib('f(x)')
sage: f.ecl()
<ECL: (($F SIMP) $X)>
sage: max_to_sr(f.ecl())
f(x)
```

TESTS:

```
sage: from sage.interfaces.maxima_lib import sr_to_max, max_to_sr
sage: f = function('f',x).diff()
sage: bool(max_to_sr(sr_to_max(f)) == f)
True
```

`sage.interfaces.maxima_lib.max_to_string(s)`

Return the Maxima string corresponding to this ECL object.

INPUT:

- s - ECL object

OUTPUT: string

EXAMPLES:

```
sage: from sage.interfaces.maxima_lib import maxima_lib, max_to_string
sage: ecl = maxima_lib(cos(x)).ecl()
sage: max_to_string(ecl)
'cos(_SAGE_VAR_x)'
```

`sage.interfaces.maxima_lib.mdiffto_sage(expr)`

Special conversion rule for %DERIVATIVE expressions.

INPUT:

- expr - ECL object; a Maxima %DERIVATIVE expression

OUTPUT: symbolic expression

EXAMPLES:

```

sage: from sage.interfaces.maxima_lib import maxima_lib, mdiff_to_sage
sage: f = maxima_lib('f(x)').diff('x', 4)
sage: f.ecl()
<ECL: ((%DERIVATIVE SIMP) (($F SIMP) $X) $X 4)>
sage: mdiff_to_sage(f.ecl())
D[0, 0, 0, 0](f)(x)

```

`sage.interfaces.maxima_lib.mlist_to_sage(expr)`

Special conversion rule for MLIST expressions.

INPUT:

- `expr` - ECL object; a Maxima MLIST expression (i.e., a list)

OUTPUT: a Python list of converted expressions.

EXAMPLES:

```

sage: from sage.interfaces.maxima_lib import maxima_lib, mlist_to_sage
sage: L=maxima_lib("[1,2,3]")
sage: L.ecl()
<ECL: ((MLIST SIMP) 1 2 3)>
sage: mlist_to_sage(L.ecl())
[1, 2, 3]

```

`sage.interfaces.maxima_lib.mqapply_to_sage(expr)`

Special conversion rule for MQAPPLY expressions.

INPUT:

- `expr` - ECL object; a Maxima MQAPPLY expression

OUTPUT: symbolic expression

MQAPPLY is used for function as `li[x](y)` and `psi[x](y)`.

EXAMPLES:

```

sage: from sage.interfaces.maxima_lib import maxima_lib, mqapply_to_sage
sage: c = maxima_lib('li[2](3)')
sage: c.ecl()
<ECL: ((MQAPPLY SIMP) (($LI SIMP ARRAY) 2) 3)>
sage: mqapply_to_sage(c.ecl())
polylog(2, 3)

```

`sage.interfaces.maxima_lib.mrat_to_sage(expr)`

Convert a Maxima MRAT expression to Sage SR.

INPUT:

- `expr` - ECL object; a Maxima MRAT expression

OUTPUT: symbolic expression

Maxima has an optimised representation for multivariate rational expressions. The easiest way to translate those to SR is by first asking Maxima to give the generic representation of the object. That is what RATDISREP does in Maxima.

EXAMPLES:

```

sage: from sage.interfaces.maxima_lib import maxima_lib, mrat_to_sage
sage: var('x y z')
(x, y, z)
sage: c = maxima_lib('(x+y^2+z^9)/x^6+z^8/y').rat()

```

```
sage: c
(_SAGE_VAR_y*_SAGE_VAR_z^9+_SAGE_VAR_x^6*_SAGE_VAR_z^8+_SAGE_VAR_y^3+_SAGE_VAR_x*_SAGE_VAR_y) / (
sage: c.ecl()
<ECL: ( (MRAT SIMP (|$_SAGE_VAR_x| |$_SAGE_VAR_y| |$_SAGE_VAR_z|)
...>
sage: mrat_to_sage(c.ecl())
(x^6*z^8 + y*z^9 + y^3 + x*y) / (x^6*y)
```

`sage.interfaces.maxima_lib.parse_max_string(s)`

Evaluate string in Maxima without *any* further simplification.

INPUT:

•s - string

OUTPUT: ECL object

EXAMPLES:

```
sage: from sage.interfaces.maxima_lib import parse_max_string
sage: parse_max_string('1+1')
<ECL: ( (MPLUS) 1 1)>
```

`sage.interfaces.maxima_lib.pyobject_to_max(obj)`

Convert a (simple) Python object into a Maxima object.

INPUT:

•expr - Python object

OUTPUT: ECL object

---

**Note:** This uses functions defined in `sage.libs.ecl`.

---

EXAMPLES:

```
sage: from sage.interfaces.maxima_lib import pyobject_to_max
sage: pyobject_to_max(4)
<ECL: 4>
sage: pyobject_to_max('z')
<ECL: Z>
sage: var('x')
x
sage: pyobject_to_max(x)
Traceback (most recent call last):
...
TypeError: Unimplemented type for python_to_ecl
```

`sage.interfaces.maxima_lib.reduce_load_MaximaLib()`

Unpickle the (unique) Maxima library interface.

EXAMPLES:

```
sage: from sage.interfaces.maxima_lib import reduce_load_MaximaLib
sage: reduce_load_MaximaLib()
Maxima_lib
```

`sage.interfaces.maxima_lib.sage_rat(x,y)`

Return quotient  $x/y$ .

INPUT:



- x - integer

- y - integer

OUTPUT: rational

EXAMPLES:

```
sage: from sage.interfaces.maxima_lib import sage_rat
sage: sage_rat(1,7)
1/7
```

`sage.interfaces.maxima_lib.sr_to_max(expr)`

Convert a symbolic expression into a Maxima object.

INPUT:

- expr - symbolic expression

OUTPUT: ECL object

EXAMPLES:

```
sage: from sage.interfaces.maxima_lib import sr_to_max
sage: var('x')
x
sage: sr_to_max(x)
<ECL: $X>
sage: sr_to_max(cos(x))
<ECL: ((%COS) $X)>
sage: f = function('f',x)
sage: sr_to_max(f.diff())
<ECL: ((%DERIVATIVE) (($F) $X) $X 1)>
```

TESTS:

We should be able to convert derivatives evaluated at a point, [trac ticket #12796](#):

```
sage: from sage.interfaces.maxima_lib import sr_to_max, max_to_sr
sage: f = function('f')
sage: f_prime = f(x).diff(x)
sage: max_to_sr(sr_to_max(f_prime(x = 1)))
D[0](f)(1)
```

`sage.interfaces.maxima_lib.stdout_to_string(s)`

Evaluate command `s` and catch Maxima stdout (not the result of the command!) into a string.

INPUT:

- s - string; command to evaluate

OUTPUT: string

This is currently used to implement `display2d()`.

EXAMPLES:

```
sage: from sage.interfaces.maxima_lib import stdout_to_string
sage: stdout_to_string('1+1')
''
sage: stdout_to_string('disp(1+1)')
'2\n\n'
```



## INTERFACE TO MUPAD

AUTHOR:

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You must have the optional commercial MuPAD interpreter installed and available as the command code{mupkern} in your PATH in order to use this interface. You do not have to install any optional sage packages.

TESTS:

```
sage: mupad.package("MuPAD-Combinat")
sage: combinat = mupad.combinat
sage: examples = mupad.examples
sage: S = examples.SymmetricFunctions()
sage: S.s[2,1]^2
s[3, 3] + s[4, 2] + s[2, 2, 1, 1] + s[2, 2, 2] + 2 s[3, 2, 1] + s[4, 1, 1] +
s[3, 1, 1, 1]
sage: S.omega( S.s[3] )
s[1, 1, 1]
sage: s = S.s
sage: p = S.p
sage: s(s[2,1] + p[2,1])
s[2, 1] + s[3] - s[1, 1, 1]
sage: s(_)
s[2, 1] + s[3] - s[1, 1, 1]
```

```
sage: combinat.tableaux.list(3) # optional - mupad # note: the order of
```

```

--                                     +---+ --
|                                     | 3 | |
|                                     +---+ +---+ +---+ |
|                                     | 3 | | 2 | | 2 | |
| +---+---+---+ +---+---+ +---+---+ +---+ |
| | 1 | 2 | 3 |, | 1 | 2 |, | 1 | 3 |, | 1 | |
-- +---+---+---+ +---+---+ +---+---+ +---+ --

```

```
sage: three = mupad(3) # optional - mupad
sage: three.combinat.tableaux.list() # optional - mupad
```

```

--                                     +---+ --
|                                     | 3 | |
|                                     +---+ |
|               +---+       +---+   +---+ |
|               | 3 |       | 2 |   | 2 | |
|   +---+ +---+ +---+   +---+ +---+   +---+ |
|   | 1 | 2 | 3 | ,   | 1 | 2 | ,   | 1 | 3 | ,   | 1 | |
--   +---+ +---+ +---+   +---+ +---+   +---+ +---+

```

```
sage: t = _[1] # optional - mupad
sage: t # optional - mupad
```

```

          +---+---+---+
          | 1 | 2 | 3 |
          +---+---+---+
sage: combinat.tableaux.conjugate(t)                                # optional - mupad
          +---+
          | 3 |
          +---+
          | 2 |
          +---+
          | 1 |
          +---+

sage: combinat.ribbonsTableaux.list([2,2],[1,1],2)                # optional - mupad
-- +---+---+ +---+---+ --
| | | 2 | | 2 | |
| + + + , +---+---+ |
| | 1 | | 1 | |
-- +---+---+ +---+---+ --

sage: combinat.tableaux.kAtom([2,1],3)                             # optional - mupad
-- +---+ --
| | 2 | |
| +---+---+ |
| | 1 | 1 | |
-- +---+---+ --

sage: M = S.Macdonald()                                           # optional - mupad
sage: a = M.P[1]^2                                                # optional - mupad
sage: mupad.mapcoeffs(a, 'normal')                                # optional - mupad
          q - t + q t - 1
P[2] + ----- P[1, 1]
          q t - 1

```

```

class sage.interfaces.mupad.Mupad(maxread=1000, script_subdirectory=None, server=None,
                                   server_tmpdir=None, logfile=None)

```

Bases: `sage.interfaces.expect.Expect`

Interface to the MuPAD interpreter.

**completions** (*string*, *strip=False*)

EXAMPLES:

```

sage: mupad.completions('linal') # optional - mupad
['linalg']

```

**console** ()

Spawn a new MuPAD command-line session.

EXAMPLES:

```

sage: mupad.console() #not tested

```

```

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```

**cputime** (*t=None*)

EXAMPLES:

```
sage: t = mupad.cputime() #random, optional - MuPAD
0.11600000000000001
```

**eval** (*code*, *strip=True*, *\*\*kws*)

EXAMPLES:

```
sage: mupad.eval('2+2') # optional - mupad
4
```

**expect** ()

EXAMPLES:

```
sage: a = mupad(1) # optional - mupad
sage: mupad.expect() # optional - mupad
<pexpect.spawn instance at 0x...>
```

**get** (*var*)

Get the value of the variable *var*.

EXAMPLES:

```
sage: mupad.set('a', 4) # optional - mupad
sage: mupad.get('a').strip() # optional - mupad
'4'
```

**set** (*var*, *value*)

Set the variable *var* to the given value.

EXAMPLES:

```
sage: mupad.set('a', 4) # optional - mupad
sage: mupad.get('a').strip() # optional - mupad
'4'
```

**trait\_names** (*verbose=True*, *use\_disk\_cache=True*)

EXAMPLES:

```
sage: names = mupad.trait_names() # optional - mupad
sage: len(names) > 100 # optional - mupad
True
sage: 'combinat' in names # optional - mupad
True
```

**class** sage.interfaces.mupad.**MupadElement** (*parent*, *value*, *is\_name=False*, *name=None*)

Bases: sage.interfaces.expect.ExpectElement

**trait\_names** ()

EXAMPLES:

```
sage: mupad.package('MuPAD-Combinat') # optional - mupad-Combinat
sage: S = mupad.examples.SymmetricFunctions() # optional - mupad-Combinat
sage: 'HallLittlewood' in S.trait_names() # optional - mupad-Combinat
True
```

**class** sage.interfaces.mupad.**MupadFunction** (*parent*, *name*)

Bases: sage.interfaces.expect.ExpectFunction

**trait\_names** ()

EXAMPLES:

```
sage: mupad.linalg.trait_names() # optional - mupad
['addCol',
 'addRow',
 ...
 'wiedemann']
```

```
class sage.interfaces.mupad.MupadFunctionElement(obj, name)
    Bases: sage.interfaces.expect.FunctionElement
```

```
    trait_names()
```

```
    EXAMPLES:
```

```
sage: three = mupad(3) # optional - mupad
```

```
sage: 'list' in three.combinat.tableaux.trait_names() # optional - mupad
True
```

```
sage.interfaces.mupad.mupad_console()
```

```
    Spawn a new MuPAD command-line session.
```

```
    EXAMPLES:
```

```
sage: from sage.interfaces.mupad import mupad_console
```

```
sage: mupad_console() #not tested
```

```

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```

```
sage.interfaces.mupad.reduce_load_mupad()
```

```
    EXAMPLES:
```

```
sage: from sage.interfaces.mupad import reduce_load_mupad
```

```
sage: reduce_load_mupad()
```

```
Mupad
```

## INTERFACE TO MWRANK

`sage.interfaces.mwrank.Mwrank` (*options='', server=None, server\_tmpdir=None*)  
Create and return an mwrank interpreter, with given options.

INPUT:

•options - string; passed when starting mwrank. The format is:

-h	help	prints this info and quits
-q	quiet	turns OFF banner display and prompt
-v n	verbosity	sets verbosity to n (default=1)
-o	PARI/GP output	turns ON extra PARI/GP short output (default is OFF)
-p n	precision	sets precision to n decimals (default=15)
-b n	quartic bound	bound on quartic point search (default=10)
-x n	n aux	number of aux primes used for sieving (default=6)
-l	list	turns ON listing of points (default ON unless v=0)
-s	selmer_only	if set, computes Selmer rank only (default: not set)
-d	skip_2nd_descent	if set, skips the second descent for curves with 2-torsion
-S n	sat_bd	upper bound on saturation primes (default=100, -1 for automatic)

EXAMPLES:

```
sage: M = Mwrank('-v 0 -l')
sage: print M('0 0 1 -1 0')
Curve [0,0,1,-1,0] :      Rank = 1
Generator 1 is [0:-1:1]; height 0.0511114082399688
Regulator = 0.0511114082399688
```

**class** `sage.interfaces.mwrank.Mwrank_class` (*options='', server=None, server\_tmpdir=None*)  
Bases: `sage.interfaces.expect.Expect`

Interface to the Mwrank interpreter.

**console** ()

Start the mwrank console.

EXAMPLE:

```
sage: mwrank.console() # not tested: expects console input
Program mwrank: ...
```

**eval** (*s, \*\*kws*)

Return mwrank's output for the given input.

INPUT:

• `s` (str) - a Sage object which when converted to a string gives valid input to `mwrnk`. The conversion is done by `validate_mwrnk_input()`. Possible formats are:

- a string representing exactly five integers separated by whitespace, for example `'1 2 3 4 5'`
- a string representing exactly five integers separated by commas, preceded by `'['` and followed by `']'` (with arbitrary whitespace), for example `'[1 2 3 4 5]'`
- a list or tuple of exactly 5 integers.

---

**Note:** If a `RuntimeError` exception is raised, then the `mwrnk` interface is restarted and the command is retried once.

---

EXAMPLES:

```
sage: mwrnk.eval('12 3 4 5 6')
'Curve [12,3,4,5,6] :... '
sage: mwrnk.eval('[12, 3, 4, 5, 6]')
'Curve [12,3,4,5,6] :... '
sage: mwrnk.eval([12, 3, 4, 5, 6])
'Curve [12,3,4,5,6] :... '
sage: mwrnk.eval((12, 3, 4, 5, 6))
'Curve [12,3,4,5,6] :... '
```

`sage.interfaces.mwrnk.mwrnk_console()`  
Start the `mwrnk` console.

EXAMPLE:

```
sage: mwrnk_console() # not tested: expects console input
Program mwrnk: ...
```

`sage.interfaces.mwrnk.validate_mwrnk_input(s)`  
Returns a string suitable for `mwrnk` input, or raises an error.

INPUT:

- `s` – one of the following:
  - a list or tuple of 5 integers `[a1,a2,a3,a4,a6]` or `(a1,a2,a3,a4,a6)`
  - a string of the form `'[a1,a2,a3,a4,a6]'` or `'a1 a2 a3 a4 a6'` where `a1, a2, a3, a4, a6` are integers

OUTPUT:

For valid input, a string of the form `'[a1,a2,a3,a4,a6]'`. For invalid input a `ValueError` is raised.

EXAMPLES:

A list or tuple of 5 integers:

```
sage: from sage.interfaces.mwrnk import validate_mwrnk_input
sage: validate_mwrnk_input([1,2,3,4,5])
'[1, 2, 3, 4, 5]'
sage: validate_mwrnk_input((-1,2,-3,4,-55))
'[-1, 2, -3, 4, -55]'
sage: validate_mwrnk_input([1,2,3,4])
Traceback (most recent call last):
...
ValueError: [1, 2, 3, 4] is not valid input to mwrnk (should have 5 entries)
sage: validate_mwrnk_input([1,2,3,4,i])
Traceback (most recent call last):
```



```
...
ValueError: [1, 2, 3, 4, I] is not valid input to mwrnk (entries should be integers)
```

A string of the form '[a1,a2,a3,a4,a6]' with any whitespace and integers ai:

```
sage: validate_mwrnk_input('0 -1 1 -7 6')
'[0,-1,1,-7,6]'
sage: validate_mwrnk_input("[0,-1,1,0,0]\n")
'[0,-1,1,0,0]'
sage: validate_mwrnk_input('0\t-1\t1\t0\t0\n')
'[0,-1,1,0,0]'
sage: validate_mwrnk_input('0 -1 1 -7 ')
Traceback (most recent call last):
...
ValueError: 0 -1 1 -7  is not valid input to mwrnk
```



## INTERFACE TO GNU OCTAVE

GNU Octave is a free software (GPL) MATLAB-like program with numerical routines for integrating, solving systems of equations, special functions, and solving (numerically) differential equations. Please see <http://octave.org/> for more details.

The commands in this section only work if you have the optional “octave” interpreter installed and available in your PATH. It’s not necessary to install any special Sage packages.

EXAMPLES:

```
sage: octave.eval('2+2')      # optional - octave
'ans = 4'
```

```
sage: a = octave(10)         # optional - octave
sage: a**10                   # optional - octave
1e+10
```

LOG: - creation (William Stein) - ? (David Joyner, 2005-12-18) - Examples (David Joyner, 2005-01-03)

### 30.1 Computation of Special Functions

Octave implements computation of the following special functions (see the maxima and gp interfaces for even more special functions):

```
airy
    Airy functions of the first and second kind, and their derivatives.
    airy(0,x) = Ai(x), airy(1,x) = Ai'(x), airy(2,x) = Bi(x), airy(3,x) = Bi'(x)
besselj
    Bessel functions of the first kind.
bessely
    Bessel functions of the second kind.
besseli
    Modified Bessel functions of the first kind.
besselk
    Modified Bessel functions of the second kind.
besselh
    Compute Hankel functions of the first (k = 1) or second (k = 2) kind.
beta
    The Beta function,
        beta(a, b) = gamma(a) * gamma(b) / gamma(a + b).
betainc
    The incomplete Beta function,
erf
    The error function,
```

erfinv

The inverse of the error function.

gamma

The Gamma function,

gammainc

The incomplete gamma function,

For example,

```
sage: octave("airy(3,2)")      # optional - octave
4.10068
sage: octave("beta(2,2)")      # optional - octave
0.166667
sage: octave("betainc(0.2,2,2)") # optional - octave
0.104
sage: octave("besselh(0,2)")   # optional - octave
(0.223891,0.510376)
sage: octave("besselh(0,1)")   # optional - octave
(0.765198,0.088257)
sage: octave("besseli(1,2)")   # optional - octave
1.59064
sage: octave("besselj(1,2)")   # optional - octave
0.576725
sage: octave("besselk(1,2)")   # optional - octave
0.139866
sage: octave("erf(0)")         # optional - octave
0
sage: octave("erf(1)")         # optional - octave
0.842701
sage: octave("erfinv(0.842)")  # optional - octave
0.998315
sage: octave("gamma(1.5)")     # optional - octave
0.886227
sage: octave("gammainc(1.5,1)") # optional - octave
0.77687
```

The Octave interface reads in even very long input (using files) in a robust manner:

```
sage: t = '"%s"'%10^10000      # ten thousand character string.
sage: a = octave.eval(t + ';' ) # optional - octave, < 1/100th of a second
sage: a = octave(t)            # optional - octave
```

Note that actually reading a back out takes forever. This *must* be fixed ASAP - see [http://trac.sagemath.org/sage\\_trac/ticket/940/](http://trac.sagemath.org/sage_trac/ticket/940/).

## 30.2 Tutorial

EXAMPLES:

```
sage: octave('4+10')          # optional - octave
14
sage: octave('date')          # optional - octave; random output
18-Oct-2007
sage: octave('5*10 + 6')      # optional - octave
56
sage: octave('(6+6)/3')       # optional - octave
```

```

4
sage: octave('9')^2          # optional - octave
81
sage: a = octave(10); b = octave(20); c = octave(30)    # optional - octave
sage: avg = (a+b+c)/3      # optional - octave
sage: avg                  # optional - octave
20
sage: parent(avg)          # optional - octave
Octave

sage: my_scalar = octave('3.1415')    # optional - octave
sage: my_scalar                  # optional - octave
3.1415
sage: my_vector1 = octave('[1,5,7]')  # optional - octave
sage: my_vector1                  # optional - octave
1      5      7
sage: my_vector2 = octave('[1;5;7]')  # optional - octave
sage: my_vector2                  # optional - octave
1
5
7
sage: my_vector1 * my_vector2        # optional - octave
75

```

```

class sage.interfaces.octave.Octave(maxread=100, script_subdirectory=None, logfile=None,
                                     server=None, server_tmpdir=None, seed=None)

```

Bases: `sage.interfaces.expect.Expect`

Interface to the Octave interpreter.

EXAMPLES:

```

sage: octave.eval("a = [ 1, 1, 2; 3, 5, 8; 13, 21, 33 ]")    # optional - octave
'a =\n\n 1 1 2\n 3 5 8\n 13 21 33\n\n'
sage: octave.eval("b = [ 1; 3; 13]")    # optional - octave
'b =\n\n 1\n 3\n 13\n\n'
sage: octave.eval("c=a \\ b") # solves linear equation: a*c = b # optional - octave; random out
'c =\n\n 1\n 7.21645e-16\n -7.21645e-16\n\n'
sage: octave.eval("c")    # optional - octave; random output
'c =\n\n 1\n 7.21645e-16\n -7.21645e-16\n\n'

```

**clear** (var)

Clear the variable named var.

EXAMPLES:

```

sage: octave.set('x', '2') # optional - octave
sage: octave.clear('x') # optional - octave
sage: octave.get('x') # optional - octave
"error: `x' undefined near line ... column 1"

```

**console** ()

Spawn a new Octave command-line session.

This requires that the optional octave program be installed and in your PATH, but no optional Sage packages need be installed.

EXAMPLES:

```

sage: octave_console()    # not tested
GNU Octave, version 2.1.73 (i386-apple-darwin8.5.3).

```

```
Copyright (C) 2006 John W. Eaton.  
...  
octave:1> 2+3  
ans = 5  
octave:2> [ctrl-d]
```

Pressing ctrl-d exits the octave console and returns you to Sage. octave, like Sage, remembers its history from one session to another.

### **de\_system\_plot** (*f, ics, trange*)

Plots (using octave's interface to gnuplot) the solution to a  $2 \times 2$  system of differential equations.

INPUT:

- *f* - a pair of strings representing the differential equations; The independent variable must be called *x* and the dependent variable must be called *y*.
- *ics* - a pair  $[x_0, y_0]$  such that  $x(t_0) = x_0$ ,  $y(t_0) = y_0$
- *trange* - a pair  $[t_0, t_1]$

OUTPUT: a gnuplot window appears

EXAMPLES:

```
sage: octave.de_system_plot(['x+y', 'x-y'], [1,-1], [0,2]) # not tested -- does this actually work?
```

This should yield the two plots  $(t, x(t))$ ,  $(t, y(t))$  on the same graph (the  $t$ -axis is the horizontal axis) of the system of ODEs

$$x' = x + y, x(0) = 1; \quad y' = x - y, y(0) = -1, \quad \text{for } 0 < t < 2.$$

### **get** (*var*)

Get the value of the variable *var*.

EXAMPLES:

```
sage: octave.set('x', '2') # optional - octave  
sage: octave.get('x') # optional - octave  
' 2'
```

### **quit** (*verbose=False*)

EXAMPLES:

```
sage: o = Octave()  
sage: o._start() # optional - octave  
sage: o.quit(True) # optional - octave  
Exiting spawned Octave process.
```

### **sage2octave\_matrix\_string** (*A*)

Return an octave matrix from a Sage matrix.

INPUT: A Sage matrix with entries in the rationals or reals.

OUTPUT: A string that evaluates to an Octave matrix.

EXAMPLES:

```
sage: M33 = MatrixSpace(QQ, 3, 3)  
sage: A = M33([1, 2, 3, 4, 5, 6, 7, 8, 0])  
sage: octave.sage2octave_matrix_string(A) # optional - octave  
'[1, 2, 3; 4, 5, 6; 7, 8, 0]'
```

**set** (*var*, *value*)

EXAMPLES:

```
sage: octave.get('x') # optional - octave
```

' 2'

```
set_seed (seed=None)
```

Sets the seed for the random number generator for this octave interpreter.

EXAMPLES:

```
sage: o = Octave() # optional - octave
```

```
sage: o.set_seed(1) # optional - octave
```

1

```
sage: [o.rand() for i in range(5)] # optional - octave
```

```
[ 0.134364,  0.847434,  0.763775,  0.255069,  0.495435]
```

```
solve_linear_system(A, b)
```

Use octave to compute a solution  $x$  to  $A^*x = b$ , as a list.

INPUT:

- $A$  -  $m \times n$  matrix  $A$  with entries in  $QQ$  or  $RR$

- $\mathbf{b}$  - m-vector  $\mathbf{b}$  entries in  $\mathbf{Q}\mathbf{Q}$  or  $\mathbf{R}\mathbf{R}$  (resp)

OUTPUT: An list  $x$  (if it exists) which solves  $M^*x = b$

EXAMPLES:

```
sage: M33 = MatrixSpace(QQ, 3, 3)
```

```
sage: A = M33([1, 2, 3, 4, 5, 6, 7, 8, 0])
```

```
sage: V3 = VectorSpace(QQ, 3)
```

```
sage: b = V3([1, 2, 3])
```

```
sage: octave.solve_linear_system(A,b)      # optional - octave (and output is slightly random)
[-0.333332999999999999, 0.666667000000000000001, -3.523660000000000002e-18]
```

AUTHORS:

•David Joyner and William Stein

```
version()
```

Return the version of Octave.

OUTPUT: string

EXAMPLES:

```
sage: octave.version()      # optional - octave; random output depending on version
'2.1.73'
```

```
class sage.interfaces.octave.OctaveElement (parent, value, is_name=False, name=None)
```

Bases: `sage.interfaces.expect.ExpectElement`

```
sage.interfaces.octave.octave_console()
```

Spawn a new Octave command-line session.

This requires that the optional octave program be installed and in your PATH, but no optional Sage packages need be installed.

**EXAMPLES:**

```
sage: octave_console()          # not tested
GNU Octave, version 2.1.73 (i386-apple-darwin8.5.3).
Copyright (C) 2006 John W. Eaton.
...
octave:1> 2+3
ans = 5
octave:2> [ctrl-d]
```

Pressing ctrl-d exits the octave console and returns you to Sage. octave, like Sage, remembers its history from one session to another.

```
sage.interfaces.octave.octave_version()
```

Return the version of Octave installed.

**EXAMPLES:**

```
sage: octave_version()          # optional - octave; and output is random
'2.9.12'
```

```
sage.interfaces.octave.reduce_load_Octave()
```

**EXAMPLES:**

```
sage: from sage.interfaces.octave import reduce_load_Octave
sage: reduce_load_Octave()
Octave
```



## INTERFACE TO PHC.

PHC computes numerical information about systems of polynomials over the complex numbers.

PHC implements polynomial homotopy methods to exploit structure in order to better approximate all isolated solutions. The package also includes extra tools to handle positive dimensional solution components.

AUTHORS:

- PHC was written by J. Verschelde, R. Cools, and many others (?)
- William Stein and Kelly ?? – first version of interface to PHC
- Marshall Hampton – second version of interface to PHC
- Marshall Hampton and Alex Jokela – third version, path tracking

**class** `sage.interfaces.phc.PHC`

A class to interface with PHCpack, for computing numerical homotopies and root counts.

EXAMPLES:

```
sage: from sage.interfaces.phc import phc
sage: R.<x,y> = PolynomialRing(CDF,2)
sage: testsys = [x^2 + 1, x*y - 1]
sage: phc.mixed_volume(testsys)          # optional -- phc
2
sage: v = phc.blackbox(testsys, R)        # optional -- phc
sage: sols = v.solutions()               # optional -- phc
sage: sols.sort()                       # optional -- phc
sage: sols                              # optional -- phc
[[-1.0000000000000000*I, 1.0000000000000000*I], [1.0000000000000000*I, -1.0000000000000000*I]]
sage: sol_dict = v.solution_dicts()      # optional -- phc
sage: x_sols_from_dict = [d[x] for d in sol_dict] # optional -- phc
sage: x_sols_from_dict.sort(); x_sols_from_dict # optional -- phc
[-1.0000000000000000*I, 1.0000000000000000*I]
sage: residuals = [[test_equation.change_ring(CDF).subs(sol) for test_equation in testsys] for sol in sol_dict]
sage: residuals                          # optional -- phc
[[0, 0], [0, 0]]
```

**blackbox** (*polys, input\_ring, verbose=False*)

Returns as a string the result of running PHC with the given polynomials under blackbox mode (the ‘-b’ option).

INPUT:

- *polys* – a list of multivariate polynomials (elements of a multivariate polynomial ring).
- *input\_ring* – for coercion of the variables into the desired ring.
- *verbose* – print lots of verbose information about what this function does.

OUTPUT:

- a `PHC_Object` object containing the phcpack output string.

EXAMPLES:

```
sage: from sage.interfaces.phc import *
sage: R2.<x,y> = PolynomialRing(QQ,2)
sage: start_sys = [x^6-y^2,y^5-1]
sage: sol = phc.blackbox(start_sys, R2)          # optional -- phc
sage: len(sol.solutions())                      # optional -- phc
30
```

**`mixed_volume`** (*polys*, *verbose=False*)

Computes the mixed volume of the polynomial system given by the input polys.

INPUT:

- *polys* – a list of multivariate polynomials (elements of a multivariate polynomial ring).
- *verbose* – print lots of verbose information about what this function does.

OUTPUT:

- The mixed volume.

EXAMPLES:

```
sage: from sage.interfaces.phc import *
sage: R2.<x,y,z> = PolynomialRing(QQ,3)
sage: test_sys = [(x+y+z)^2-1,x^2-x,y^2-1]
sage: phc.mixed_volume(test_sys)                # optional -- phc
4
```

**`path_track`** (*start\_sys*, *end\_sys*, *input\_ring*, *c\_skew=0.001*, *saved\_start=None*)

This function computes homotopy paths between the solutions of *start\_sys* and *end\_sys*.

INPUT:

- *start\_sys* – a square polynomial system, given as a list of polynomials
- *end\_sys* – same type as *start\_sys*
- *input\_ring* – for coercion of the variables into the desired ring.
- *c\_skew* – optional. the imaginary part of homotopy multiplier; nonzero values are often necessary to avoid intermediate path collisions
- *saved\_start* – optional. A phc output file. If not given, start system solutions are computed via the `phc.blackbox` function.

OUTPUT:

- a list of paths as dictionaries, with the keys variables and t-values on the path.

EXAMPLES:

```
sage: from sage.interfaces.phc import *
sage: R2.<x,y> = PolynomialRing(QQ,2)
sage: start_sys = [x^6-y^2,y^5-1]
sage: sol = phc.blackbox(start_sys, R2)          # optional -- phc
sage: start_save = sol.save_as_start()           # optional -- phc
sage: end_sys = [x^7-2,y^5-x^2]                 # optional -- phc
sage: sol_paths = phc.path_track(start_sys, end_sys, R2, saved_start = start_save) # optional -- phc
sage: len(sol_paths)                            # optional -- phc
30
```

**plot\_paths\_2d**(*start\_sys*, *end\_sys*, *input\_ring*, *c\_skew*=0.001, *endpoints*=True, *saved\_start*=None, *rand\_colors*=False)

This returns a graphics object of solution paths in the complex plane.

INPUT:

- *start\_sys* – a square polynomial system, given as a list of polynomials
- *end\_sys* – same type as *start\_sys*
- *input\_ring* – for coercion of the variables into the desired ring.
- *c\_skew* – optional. the imaginary part of homotopy multiplier; nonzero values are often necessary to avoid intermediate path collisions
- *endpoints* – optional. Whether to draw in the ends of paths as points.
- *saved\_start* – optional. A phc output file. If not given, start system solutions are computed via the `phc.blackbox` function.

OUTPUT:

- lines and points of solution paths

EXAMPLES:

```
sage: from sage.interfaces.phc import *
sage: from sage.structure.sage_object import SageObject
sage: R2.<x,y> = PolynomialRing(QQ,2)
sage: start_sys = [x^5-y^2,y^5-1]
sage: sol = phc.blackbox(start_sys, R2)      # optional -- phc
sage: start_save = sol.save_as_start()      # optional -- phc
sage: end_sys = [x^5-25,y^5-x^2]           # optional -- phc
sage: testing = phc.plot_paths_2d(start_sys, end_sys, R2) # optional -- phc
sage: type(testing)                         # optional -- phc (normally use plot here)
<class 'sage.plot.graphics.Graphics'>
```

**start\_from**(*start\_filename\_or\_string*, *polys*, *input\_ring*, *path\_track\_file*=None, *verbose*=False)

This computes solutions starting from a phcpack solution file.

INPUT:

- *start\_filename\_or\_string* – the filename for a phcpack start system, or the contents of such a file as a string. Variable names must match the inputring variables. The value of the homotopy variable *t* should be 1, not 0.
- *polys* – a list of multivariate polynomials (elements of a multivariate polynomial ring).
- *input\_ring*: for coercion of the variables into the desired ring.
- *path\_track\_file*: whether to save path-tracking information
- *verbose* – print lots of verbose information about what this function does.

OUTPUT:

- A solution in the form of a PHCObject.

EXAMPLES:

```
sage: from sage.interfaces.phc import *
sage: R2.<x,y> = PolynomialRing(QQ,2)
sage: start_sys = [x^6-y^2,y^5-1]
sage: sol = phc.blackbox(start_sys, R2)      # optional -- phc
sage: start_save = sol.save_as_start()      # optional -- phc
sage: end_sys = [x^7-2,y^5-x^2]             # optional -- phc
```

```

sage: sol = phc.start_from(start_save, end_sys, R2) # optional -- phc
sage: len(sol.solutions())                        # optional -- phc
30

```

**class** `sage.interfaces.phc.PHC_Object` (*output\_file\_contents*, *input\_ring*)

A container for data from the PHCpack program - lists of float solutions, etc. Currently the file contents are kept as a string; for really large outputs this would be bad.

INPUT:

- `output_file_contents`: the string output of PHCpack
- `input_ring`: for coercion of the variables into the desired ring.

EXAMPLES:

```

sage: from sage.interfaces.phc import phc
sage: R2.<x,y> = PolynomialRing(QQ,2)
sage: start_sys = [(x-1)^2+(y-1)-1, x^2+y^2-1]
sage: sol = phc.blackbox(start_sys, R2) # optional -- phc
sage: str(sum([x[0] for x in sol.solutions()]).real())[0:3] # optional -- phc
'2.0'

```

**classified\_solution\_dicts** ()

Returns a dictionary of lists of dictionaries of solutions. Its not as crazy as it sounds; the keys are the types of solutions as classified by phcpack: regular vs. singular, complex vs. real

INPUT:

- None

OUTPUT:

- A dictionary of lists of dictionaries of solutions

EXAMPLES:

```

sage: from sage.interfaces.phc import phc
sage: R.<x,y> = PolynomialRing(CC,2)
sage: p_sys = [x^10-y,y^2-1]
sage: sol = phc.blackbox(p_sys,R) # optional -- phc
sage: classifieds = sol.classified_solution_dicts() # optional -- phc
sage: str(sum([q[y] for q in classifieds['real']]))[0:3] # optional -- phc
'2.0'

```

**save\_as\_start** (*start\_filename=None*, *sol\_filter=''*)

Saves a solution as a phcpack start file. The usual output is just as a string, but it can be saved to a file as well. Even if saved to a file, it still returns the output string.

EXAMPLES:

```

sage: from sage.interfaces.phc import phc
sage: R2.<x,y> = PolynomialRing(QQ,2)
sage: start_sys = [x^3-y^2,y^5-1]
sage: sol = phc.blackbox(start_sys, R2) # optional -- phc
sage: start_save = sol.save_as_start() # optional -- phc
sage: end_sys = [x^7-2,y^5-x^2] # optional -- phc
sage: sol = phc.start_from(start_save, end_sys, R2) # optional -- phc
sage: len(sol.solutions()) # optional -- phc
15

```

**solution\_dicts** (*get\_failures=False*)

Returns a list of solutions in dictionary form: variable:value.

INPUT:

- self – for access to self\_out\_file\_contents, the string of raw PHCpack output.
- get\_failures (optional) – a boolean. The default (False) is to not process failed homotopies. These either lie on positive-dimensional components or at infinity.

OUTPUT:

- solution\_dicts: a list of dictionaries. Each dictionary element is of the form variable:value, where the variable is an element of the input\_ring, and the value is in ComplexField.

EXAMPLES:

```
sage: from sage.interfaces.phc import *
sage: R.<x,y,z> = PolynomialRing(QQ,3)
sage: fs = [x^2-1,y^2-x,z^2-y]
sage: sol = phc.blackbox(fs,R)           # optional -- phc
sage: s_list = sol.solution_dicts()      # optional -- phc
sage: s_list.sort()                     # optional -- phc
sage: s_list[0]                          # optional -- phc
{y: 1.000000000000000, z: -1.000000000000000, x: 1.000000000000000}
```

**solutions** (*get\_failures=False*)

Returns a list of solutions in the ComplexField.

Use the variable\_list function to get the order of variables used by PHCpack, which is usually different than the term order of the input\_ring.

INPUT:

- self – for access to self\_out\_file\_contents, the string of raw PHCpack output.
- get\_failures (optional) – a boolean. The default (False) is to not process failed homotopies. These either lie on positive-dimensional components or at infinity.

OUTPUT:

- solutions: a list of lists of ComplexField-valued solutions.

EXAMPLES:

```
sage: from sage.interfaces.phc import *
sage: R2.<x1,x2> = PolynomialRing(QQ,2)
sage: test_sys = [x1^5-x1*x2^2-1, x2^5-x1*x2-1]
sage: sol = phc.blackbox(test_sys, R2)   # optional -- phc
sage: len(sol.solutions())               # optional -- phc
25
```

**variable\_list** ()

Returns the variables, as strings, in the order in which PHCpack has processed them.

EXAMPLES:

```
sage: from sage.interfaces.phc import *
sage: R2.<x1,x2> = PolynomialRing(QQ,2)
sage: test_sys = [x1^5-x1*x2^2-1, x2^5-x1*x2-1]
sage: sol = phc.blackbox(test_sys, R2)   # optional -- phc
sage: sol.variable_list()                # optional -- phc
['x1', 'x2']
```

```
sage.interfaces.phc.get_classified_solution_dicts(output_file_contents, input_ring,
                                                get_failures=True)
```

Returns a dictionary of lists of dictionaries of variable:value (key:value) pairs. Only used internally; see the `classified_solution_dict` function in the `PHC_Object` class definition for details.

INPUT:

- `output_file_contents` – phc solution output as a string
- `input_ring` – a `PolynomialRing` that variable names can be coerced into

OUTPUT:

- a dictionary of lists if dictionaries of solutions, classifies by type

EXAMPLES:

```
sage: from sage.interfaces.phc import *
sage: R2.<x1,x2> = PolynomialRing(QQ,2)
sage: test_sys = [(x1-2)^5-x2, (x2-1)^5-1]
sage: sol = phc.blackbox(test_sys, R2) # optional -- phc
sage: sol_classes = get_classified_solution_dicts(sol.output_file_contents,R2) # optional -- phc
sage: len(sol_classes['real']) # optional -- phc
1
```

```
sage.interfaces.phc.get_solution_dicts(output_file_contents, input_ring,
                                       get_failures=True)
```

Returns a list of dictionaries of variable:value (key:value) pairs. Only used internally; see the `solution_dict` function in the `PHC_Object` class definition for details.

INPUT:

- `output_file_contents` – phc solution output as a string
- `input_ring` – a `PolynomialRing` that variable names can be coerced into

OUTPUT:

a list of dictionaries of solutions

EXAMPLES:

```
sage: from sage.interfaces.phc import *
sage: R2.<x1,x2> = PolynomialRing(QQ,2)
sage: test_sys = [(x1-1)^5-x2, (x2-1)^5-1]
sage: sol = phc.blackbox(test_sys, R2) # optional -- phc
sage: test = get_solution_dicts(sol.output_file_contents,R2) # optional -- phc
sage: str(sum([q[x1].real() for q in test]))[0:4] # optional -- phc
'25.0'
```

```
sage.interfaces.phc.get_variable_list(output_file_contents)
```

Returns the variables, as strings, in the order in which PHCpack has processed them.

EXAMPLES:

```
sage: from sage.interfaces.phc import *
sage: R2.<x1,x2> = PolynomialRing(QQ,2)
sage: test_sys = [(x1-2)^5-x2, (x2-1)^5-1]
sage: sol = phc.blackbox(test_sys, R2) # optional -- phc
sage: get_variable_list(sol.output_file_contents) # optional -- phc
['x1', 'x2']
```

## POV-RAY, THE PERSISTENCE OF VISION RAY TRACER

**class** `sage.interfaces.povray.POVRay`  
POV-Ray The Persistence of Vision Ray Tracer

INPUT:

- `pov_file` – complete path to the .pov file you want to be rendered
- `outfile` – the filename you want to save your result to
- `**kwargs` – additionally keyword arguments you want to pass to POV-Ray

OUTPUT:

Image is written to the file you specified in `outfile`

EXAMPLES:

AUTHOR:

Sage interface written by Yi Qiang (`yqiang_atNOSPAM_gmail.com`)

POV-Ray: <http://www.povray.org>

**usage** ()





## PARALLEL INTERFACE TO THE SAGE INTERPRETER

This is an expect interface to `emph{multiple}` copy of the sage interpreter, which can all run simultaneous calculations. A PSage object does not work as well as the usual Sage object, but does have the great property that when you construct an object in a PSage you get back a prompt immediately. All objects constructed for that PSage print `<<currently executing code>>` until code execution completes, when they print as normal.

note{BUG – currently non-idle PSage subprocesses do not stop when sage exits. I would very much like to fix this but don't know how.}

EXAMPLES:

We illustrate how to factor 3 integers in parallel. First start up 3 parallel Sage interfaces:

```
sage: v = [PSage() for _ in range(3)]
```

Next, request factorization of one random integer in each copy.

```
sage: w = [x('factor(2^%s-1)'% randint(250,310)) for x in v] # long time (5s on sage.math, 2011)
```

Print the status:

```
sage: w # long time, random output (depends on timing)
[3 * 11 * 31^2 * 311 * 11161 * 11471 * 73471 * 715827883 * 2147483647 * 4649919401 * 18158209813151 *
<<currently executing code>>,
9623 * 68492481833 * 23579543011798993222850893929565870383844167873851502677311057483194673]
```

Note that at the point when we printed two of the factorizations had finished but a third one hadn't. A few seconds later all three have finished:

```
sage: w # long time, random output
[3 * 11 * 31^2 * 311 * 11161 * 11471 * 73471 * 715827883 * 2147483647 * 4649919401 * 18158209813151 *
23^2 * 47 * 89 * 178481 * 4103188409 * 199957736328435366769577 * 44667711762797798403039426178361,
9623 * 68492481833 * 23579543011798993222850893929565870383844167873851502677311057483194673]
```

```
class sage.interfaces.p sage.PSage(**kws)
    Bases: sage.interfaces.sage0.Sage

    eval(x, strip=True, **kws)
        x – code strip – ignored

    get(var)
        Get the value of the variable var.

    is_locked()

    set(var, value)
        Set the variable var to the given value.
```

```
class sage.interfaces.psage.PSageElement (parent, value, is_name=False, name=None)  
    Bases: sage.interfaces.sage0.SageElement  
    is_locked()
```

## INTERFACE TO QEPCAD

---

The basic function of QEPCAD is to construct cylindrical algebraic decompositions (CADs) of  $\mathbf{R}^k$ , given a list of polynomials. Using this CAD, it is possible to perform quantifier elimination and formula simplification.

A CAD for a set  $A$  of  $k$ -variate polynomials decomposes  $\mathbf{R}^j$  into disjoint cells, for each  $j$  in  $0 \leq j \leq k$ . The sign of each polynomial in  $A$  is constant in each cell of  $\mathbf{R}^k$ , and for each cell in  $\mathbf{R}^j$  ( $j > 1$ ), the projection of that cell into  $\mathbf{R}^{j-1}$  is a cell of  $\mathbf{R}^{j-1}$ . (This property makes the decomposition ‘cylindrical’.)

Given a formula  $\exists x.P(a,b,x) = 0$  (for a polynomial  $P$ ), and a cylindrical algebraic decomposition for  $P$ , we can eliminate the quantifier (find an equivalent formula in the two variables  $a, b$  without the quantifier  $\exists$ ) as follows. For each cell  $C$  in  $\mathbf{R}^2$ , find the cells of  $\mathbf{R}^3$  which project to  $C$ . (This collection is called the `stack` over  $C$ .) Mark  $C$  as true if some member of the stack has sign = 0; otherwise, mark  $C$  as false. Then, construct a polynomial formula in  $a, b$  which specifies exactly the true cells (this is always possible). The same technique works if the body of the quantifier is any boolean combination of polynomial equalities and inequalities.

Formula simplification is a similar technique. Given a formula which describes a simple set of  $\mathbf{R}^k$  in a complicated way as a boolean combination of polynomial equalities and inequalities, QEPCAD can construct a CAD for the polynomials and recover a simple equivalent formula.

Note that while the following documentation is tutorial in nature, and is written for people who may not be familiar with QEPCAD, it is documentation for the sage interface rather than for QEPCAD. As such, it does not cover several issues that are very important to use QEPCAD efficiently, such as variable ordering, the efficient use of the alternate quantifiers and `_root_` expressions, the `measure-zero-error` command, etc. For more information on QEPCAD, see the online documentation at url{<http://www.cs.usna.edu/~qepcad/B/QEPCAD.html>} and Chris Brown’s tutorial handout and slides from url{<http://www.cs.usna.edu/~wcbrown/research/ISSAC04/Tutorial.html>}. (Several of the examples in this documentation came from these sources.)

The examples below require that the optional qepcad package is installed.

QEPCAD can be run in a fully automatic fashion, or interactively. We first demonstrate the automatic use of QEPCAD.

Since sage has no built-in support for quantifiers, this interface provides `qepcad_formula` which helps construct quantified formulas in the syntax QEPCAD requires.

```
sage: var('a,b,c,d,x,y,z')
(a, b, c, d, x, y, z)
sage: qf = qepcad_formula
```

We start with a simple example. Consider an arbitrarily-selected ellipse:

```
sage: ellipse = 3*x^2 + 2*x*y + y^2 - x + y - 7
```

What is the projection onto the  $x$  axis of this ellipse? First we construct a formula asking this question.

```
sage: F = qf.exists(y, ellipse == 0); F
(E y) [3 x^2 + 2 x y + y^2 - x + y - 7 = 0]
```

Then we run `qepcad` to get the answer:

```
sage: qepcad(F) # optional - qepcad
8 x^2 - 8 x - 29 <= 0
```

How about the projection onto the  $y$  axis?

```
sage: qepcad(qf.exists(x, ellipse == 0)) # optional - qepcad
8 y^2 + 16 y - 85 <= 0
```

QEPCAD deals with more quantifiers than just ‘exists’, of course. Besides the standard ‘forall’, there are also ‘for infinitely many’, ‘for all but finitely many’, ‘for a connected subset’, and ‘for exactly  $k$ ’. The `qepcad()` documentation has examples of all of these; here we will just give one example.

First we construct a circle:

```
sage: circle = x^2 + y^2 - 3
```

For what values  $k$  does a vertical line  $x = k$  intersect the combined figure of the circle and ellipse exactly three times?

```
sage: F = qf.exactly_k(3, y, circle * ellipse == 0); F
(X3 y) [(3 x^2 + 2 x y + y^2 - x + y - 7) (x^2 + y^2 - 3) = 0]
sage: qepcad(F) # not tested (random order)
x^2 - 3 <= 0 /\ 8 x^2 - 8 x - 29 <= 0 /\ 8 x^4 - 26 x^2 - 4 x + 13 >= 0 /\ [ 8 x^4 - 26 x^2 - 4 x + 13 >= 0 /\ 8 x^2 - 8 x - 29 <= 0 /\ x^2 - 3 <= 0 ]
```

Here we see that the solutions are among the eight  $(4 + 2 + 2)$  roots of the three polynomials inside the brackets, but not all of these roots are solutions; the polynomial inequalities outside the brackets are needed to select those roots that are solutions.

QEPCAD also supports an extended formula language, where `_root_ $k$`   $P(\bar{x}, y)$  refers to a particular zero of  $P(\bar{x}, y)$  (viewed as a polynomial in  $y$ ). If there are  $n$  roots, then `_root_1` refers to the least root and `_root_ $n$`  refers to the greatest. Also, `_root_ $-n$`  refers to the least root and `_root_ $-1$`  refers to the greatest.

This extended language is available both on input and output; see the QEPCAD documentation for more information on how to use this syntax on input. We can request output that is intended to be easy to interpret geometrically; then QEPCAD will use the extended language to produce a solution formula without the selection polynomials.

```
sage: qepcad(F, solution='geometric') # not tested (random order)
x = _root_1 8 x^2 - 8 x - 29
\/
8 x^4 - 26 x^2 - 4 x + 13 = 0
\/
x = _root_-1 x^2 - 3
```

We then see that the 6 solutions correspond to the vertical tangent on the left side of the ellipse, the four intersections between the ellipse and the circle, and the vertical tangent on the right side of the circle.

Let us do some basic formula simplification and visualization. We will look at the region which is inside both the ellipse and the circle:

```
sage: F = qf.and_(ellipse < 0, circle < 0); F
[3 x^2 + 2 x y + y^2 - x + y - 7 < 0 /\ x^2 + y^2 - 3 < 0]
sage: qepcad(F) # not tested (random order)
y^2 + 2 x y + y + 3 x^2 - x - 7 < 0 /\ y^2 + x^2 - 3 < 0
```

We get back the same formula we put in. This is not surprising (we started with a pretty simple formula, after all), but it is not very enlightening either. Again, if we ask for a ‘geometric’ output, then we see an output that lets us understand something about the shape of the solution set.

```
sage: qepcad(F, solution='geometric') # not tested (random order)
[
  [
    x = _root_-2 8 x^4 - 26 x^2 - 4 x + 13
    /\
    x = _root_2 8 x^4 - 26 x^2 - 4 x + 13
    /\
    8 x^4 - 26 x^2 - 4 x + 13 < 0
  ]
  /\
  y^2 + 2 x y + y + 3 x^2 - x - 7 < 0
  /\
  y^2 + x^2 - 3 < 0
]
/>\
[
  x > _root_2 8 x^4 - 26 x^2 - 4 x + 13
  /\
  x < _root_-2 8 x^4 - 26 x^2 - 4 x + 13
  /\
  y^2 + x^2 - 3 < 0
]
```

There is another reason to prefer output using `_root_` expressions; not only does it sometimes give added insight into the geometric structure, it also can be more efficient to construct. Consider this formula for the projection of a particular semicircle onto the  $x$  axis:

```
sage: F = qf.exists(y, qf.and_(circle == 0, x + y > 0)); F
(E y) [x^2 + y^2 - 3 = 0 /\ x + y > 0]
sage: qepcad(F) # not tested (random order)
x^2 - 3 <= 0 /\ [ x > 0 /\ 2 x^2 - 3 < 0 ]
```

Here, the formula  $x > 0$  had to be introduced in order to get a solution formula; the original CAD of  $F$  did not include the polynomial  $x$ . To avoid having QEPCAD do the extra work to come up with a solution formula, we can tell it to use the extended language; it is always possible to construct a solution formula in the extended language without introducing new polynomials.

```
sage: qepcad(F, solution='extended') # not tested (random order)
x^2 - 3 <= 0 /\ x > _root_1 2 x^2 - 3
```

Up to this point, all the output we have seen has basically been in the form of strings; there is no support (yet) for parsing these outputs back into sage polynomials (partly because sage does not yet have support for symbolic conjunctions and disjunctions). The function `qepcad()` supports three more output types that give numbers which can be manipulated in sage: any-point, all-points, and cell-points.

These output types give dictionaries mapping variable names to values. With any-point, `qepcad()` either produces a single dictionary specifying a point where the formula is true, or raises an exception if the formula is false everywhere. With all-points, `qepcad()` either produces a list of dictionaries for all points where the formula is true, or raises an exception if the formula is true on infinitely many points. With cell-points, `qepcad()` produces a list of dictionaries with one point for each cell where the formula is true. (This means you will have at least one point in each connected component of the solution, although you will often have many more points than that.)

Let us revisit some of the above examples and get some points to play with. We will start by finding a point on our ellipse.

```
sage: p = qepcad(ellipse == 0, solution='any-point'); p # optional - qepcad
{'x': -1.468501968502953?, 'y': 0.9685019685029527?}
```

(Note that despite the decimal printing and the question marks, these are really exact numbers.)

We can verify that this point is a solution. To do so, we create a copy of ellipse as a polynomial over  $\mathbb{Q}$  (instead of a symbolic expression).

```
sage: pellipse = QQ['x,y'](ellipse)
sage: pellipse(**p) == 0 # optional - qepcad
True
```

For cell-points, let us look at points *not* on the ellipse.

```
sage: pts = qepcad(ellipse != 0, solution='cell-points'); pts # optional - qepcad
[{'x': 4, 'y': 0},
 {'x': 2.468501968502953?, 'y': 1},
 {'x': 2.468501968502953?, 'y': -9},
 {'x': 1/2, 'y': 9},
 {'x': 1/2, 'y': -1},
 {'x': 1/2, 'y': -5},
 {'x': -1.468501968502953?, 'y': 3},
 {'x': -1.468501968502953?, 'y': -1},
 {'x': -3, 'y': 0}]
```

For the points here which are in full-dimensional cells, QEPCAD has the freedom to choose rational sample points, and it does so.

And, of course, all these points really are not on the ellipse.

```
sage: [pellipse(**p) != 0 for p in pts] # optional - qepcad
[True, True, True, True, True, True, True, True, True]
```

Finally, for all-points, let us look again at finding vertical lines that intersect the union of the circle and the ellipse exactly three times.

```
sage: F = qf.exactly_k(3, y, circle * ellipse == 0); F
(X3 y)[(3 x^2 + 2 x y + y^2 - x + y - 7) (x^2 + y^2 - 3) = 0]
sage: pts = qepcad(F, solution='all-points'); pts # optional - qepcad
[{'x': 1.732050807568878?}, {'x': 1.731054913462534?}, {'x': 0.678911384208004?}, {'x': -0.941772737?}]
```

Since  $y$  is bound by the quantifier, the solutions only refer to  $x$ .

We can substitute one of these solutions into the original equation:

```
sage: pt = pts[0] # optional - qepcad
sage: pcombo = QQ['x,y'](circle * ellipse)
sage: intersections = pcombo(y=polygen(AA, 'y'), **pt); intersections # optional - qepcad
y^4 + 4.464101615137755?y^3 + 0.2679491924311227?y^2
```

and verify that it does have three roots:

```
sage: intersections.roots() # optional - qepcad
[(-4.403249005600958?, 1), (-0.06085260953679653?, 1), (0, 2)]
```

Let us check all six solutions.

```
sage: [len(pcombo(y=polygen(AA, 'y'), **p).roots()) for p in pts] # optional - qepcad
[3, 3, 3, 3, 3, 3]
```

We said earlier that we can run QEPCAD either automatically or interactively. Now that we have discussed the automatic modes, let us turn to interactive uses.

If the `qepcad()` function is passed `interact=True`, then instead of returning a result, it returns an object of class `Qepcad` representing a running instance of QEPCAD that you can interact with. For example:

```
sage: qe = qepcad(qf.forall(x, x^2 + b*x + c > 0), interact=True); qe      # optional - qepcad
QEPCAD object in phase 'Before Normalization'
```

This object is a fairly thin wrapper over QEPCAD; most QEPCAD commands are available as methods on the `Qepcad` object. Given a `Qepcad` object `qe`, you can type `qe.[tab]` to see the available QEPCAD commands; to see the documentation for an individual QEPCAD command, for example `d_setting`, you can type `qe.d_setting?`. (In QEPCAD, this command is called `d-setting`. We systematically replace hyphens with underscores for this interface.)

The execution of QEPCAD is divided into four phases. Most commands are not available during all phases. We saw above that QEPCAD starts out in phase 'Before Normalization'. We see that the `d_cell` command is not available in this phase:

```
sage: qe.d_cell()                                # optional - qepcad
Error GETCID: This command is not active here.
```

We will focus here on the fourth (and last) phase, 'Before Solution', because this interface has special support for some operations in this phase. Consult the QEPCAD documentation for information on the other phases.

We can tell QEPCAD to finish off the current phase and move to the next with its `go` command. (There is also the `step` command, which partially completes a phase for phases that have multiple steps, and the `finish` command, which runs QEPCAD to completion.)

```
sage: qe.go()                                    # optional - qepcad
QEPCAD object has moved to phase 'Before Projection (x)'
```

```
sage: qe.go()                                    # optional - qepcad
QEPCAD object has moved to phase 'Before Choice'
```

```
sage: qe.go()                                    # optional - qepcad
QEPCAD object has moved to phase 'Before Solution'
```

Note that the `Qepcad` object returns the new phase whenever the phase changes, as a convenience for interactive use; except that when the new phase is 'EXITED', the solution formula printed by QEPCAD is returned instead.

```
sage: qe.go()                                    # optional - qepcad
4 c - b^2 > 0
```

```
sage: qe                                          # optional - qepcad
QEPCAD object in phase 'EXITED'
```

Let us pick a nice, simple example, return to phase 4, and explore the resulting `qe` object.

```
sage: qe = qepcad(circle == 0, interact=True); qe      # optional - qepcad
QEPCAD object in phase 'Before Normalization'
```

```
sage: qe.go(); qe.go(); qe.go()                    # optional - qepcad
QEPCAD object has moved to phase 'Before Projection (y)'
```

```
QEPCAD object has moved to phase 'Before Choice'
```

```
QEPCAD object has moved to phase 'Before Solution'
```

We said before that QEPCAD creates 'cylindrical algebraic decompositions'; since we have a bivariate polynomial, we get decompositions of  $\mathbf{R}^0$ ,  $\mathbf{R}^1$ , and  $\mathbf{R}^2$ . In this case, where our example is a circle of radius  $\sqrt{3}$  centered on the origin, these decompositions are as follows:

The decomposition of  $\mathbf{R}^0$  is trivial (of course). The decomposition of  $\mathbf{R}^1$  has five cells:  $x < -\sqrt{3}$ ,  $x = -\sqrt{3}$ ,  $-\sqrt{3} < x < \sqrt{3}$ ,  $x = \sqrt{3}$ , and  $x > \sqrt{3}$ . These five cells comprise the `stack` over the single cell in the trivial

decomposition of  $\mathbf{R}^0$ .

These five cells give rise to five stacks in  $\mathbf{R}^2$ . The first and fifth stack have just one cell apiece. The second and fourth stacks have three cells:  $y < 0$ ,  $y = 0$ , and  $y > 0$ . The third stack has five cells: below the circle, the lower semicircle, the interior of the circle, the upper semicircle, and above the circle.

QEPCAD (and this QEPCAD interface) number the cells in a stack starting with 1. Each cell has an `index`, which is a tuple of integers describing the path to the cell in the tree of all cells. For example, the cell ‘below the circle’ has index (3,1) (the first cell in the stack over the third cell of  $\mathbf{R}^1$ ) and the interior of the circle has index (3,3).

We can view these cells with the QEPCAD command `d_cell`. For instance, let us look at the cell for the upper semicircle:

```
sage: qe.d_cell(3, 4)                                     # optional - qepcad
----- Information about the cell (3,4) -----
Level                               : 2
Dimension                           : 1
Number of children                   : 0
Truth value                         : T      by trial evaluation.
Degrees after substitution          : Not known yet or No polynomial.
Multiplicities                      : ((1,1))
Signs of Projection Factors
Level 1 : (-)
Level 2 : (0)
----- Sample point -----
The sample point is in a PRIMITIVE representation.

alpha = the unique root of x^2 - 3 between 0 and 4
      = 1.7320508076-

Coordinate 1 = 0
            = 0.0000000000
Coordinate 2 = alpha
            = 1.7320508076-
-----
```

We see that, the level of this cell is 2, meaning that it is part of the decomposition of  $\mathbf{R}^2$ . The dimension is 1, meaning that the cell is homeomorphic to a line (rather than a plane or a point). The sample point gives the coordinates of one point in the cell, both symbolically and numerically.

For programmatic access to cells, we have defined a sage wrapper class `QepcadCell`. These cells can be created with the `cell()` method; for example:

```
sage: c = qe.cell(3, 4); c                               # optional - qepcad
QEPCAD cell (3, 4)
```

A `QepcadCell` has accessor methods for the important state held within a cell. For instance:

```
sage: c.level()                                           # optional - qepcad
2
sage: c.index()                                           # optional - qepcad
(3, 4)
sage: qe.cell(3).number_of_children()                   # optional - qepcad
5
sage: len(qe.cell(3))                                     # optional - qepcad
5
```

One particularly useful thing we can get from a cell is its sample point, as sage algebraic real numbers.



```

sage: c.sample_point()                                # optional - qepcad
(0, 1.732050807568878?)
sage: c.sample_point_dict()                           # optional - qepcad
{'x': 0, 'y': 1.732050807568878?}

```

We have seen that we can get cells using the `cell()` method. There are several QEPCAD commands that print lists of cells; we can also get cells using the `make_cells()` method, passing it the output of one of these commands.

```

sage: qe.make_cells(qe.d_true_cells())                 # optional - qepcad
[QEPCAD cell (4, 2), QEPCAD cell (3, 4), QEPCAD cell (3, 2),
QEPCAD cell (2, 2)]

```

Also, the cells in the stack over a given cell can be accessed using array subscripting or iteration. (Remember that cells in a stack are numbered starting with one; we preserve this convention in the array-subscripting syntax.)

```

sage: c = qe.cell(3)                                  # optional - qepcad
sage: c[1]                                             # optional - qepcad
QEPCAD cell (3, 1)
sage: [c2 for c2 in c]                                # optional - qepcad
[QEPCAD cell (3, 1), QEPCAD cell (3, 2), QEPCAD cell (3, 3),
QEPCAD cell (3, 4), QEPCAD cell (3, 5)]

```

We can do one more thing with a cell: we can set its truth value. Once the truth values of the cells have been set, we can get QEPCAD to produce a formula which is true in exactly the cells we have selected. This is useful if QEPCAD's quantifier language is insufficient to express your problem.

For example, consider again our combined figure of the circle and the ellipse. Suppose you want to find all vertical lines that intersect the circle twice, and also intersect the ellipse twice. The vertical lines that intersect the circle twice can be found by simplifying:

```

sage: F = qf.exactly_k(2, y, circle == 0); F
(X2 y) [x^2 + y^2 - 3 = 0]

```

and the vertical lines that intersect the ellipse twice are expressed by:

```

sage: G = qf.exactly_k(2, y, ellipse == 0); G
(X2 y) [3 x^2 + 2 x y + y^2 - x + y - 7 = 0]

```

and the lines that intersect both figures would be:

```

sage: qf.and_(F, G)
Traceback (most recent call last):
...
ValueError: QEPCAD formulas must be in prenex (quantifiers outermost) form

```

...except that QEPCAD does not support formulas like this; in QEPCAD input, all logical connectives must be inside all quantifiers.

Instead, we can get QEPCAD to construct a CAD for our combined figure and set the truth values ourselves. (The exact formula we use doesn't matter, since we're going to replace the truth values in the cells; we just need to use a formula that uses both polynomials.)

```

sage: qe = qepcad(qf.and_(ellipse == 0, circle == 0), interact=True) # optional - qepcad
sage: qe.go(); qe.go(); qe.go()                                     # optional - qepcad
QEPCAD object has moved to phase 'Before Projection (y)'
QEPCAD object has moved to phase 'Before Choice'
QEPCAD object has moved to phase 'Before Solution'

```

Now we want to find all cells  $c$  in the decomposition of  $\mathbf{R}^1$  such that the stack over  $c$  contains exactly two cells on the ellipse, and also contains exactly two cells on the circle.

Our input polynomials are ‘level-2 projection factors’, we see:

```
sage: qe.d_proj_factors()                                # optional - qepcad
P_1,1 = fac(J_1,1) = fac(dis(A_2,1))
      = 8 x^2 - 8 x - 29
P_1,2 = fac(J_1,2) = fac(dis(A_2,2))
      = x^2 - 3
P_1,3 = fac(J_1,3) = fac(res(A_2,1|A_2,2))
      = 8 x^4 - 26 x^2 - 4 x + 13
A_2,1 = input
      = y^2 + 2 x y + y + 3 x^2 - x - 7
A_2,2 = input
      = y^2 + x^2 - 3
```

so we can test whether a cell is on the ellipse by checking that the sign of the corresponding projection factor is 0 in our cell. For instance, the cell (12,2) is on the ellipse:

```
sage: qe.cell(12,2).signs()[1][0]                      # optional - qepcad
0
```

So we can update the truth values as desired like this:

```
sage: for c in qe.cell():                               # optional - qepcad
.....:     count_ellipse = 0
.....:     count_circle = 0
.....:     for c2 in c:
.....:         count_ellipse += (c2.signs()[1][0] == 0)
.....:         count_circle += (c2.signs()[1][1] == 0)
.....:     c.set_truth(count_ellipse == 2 and count_circle == 2)
```

and then we can get our desired solution formula. (The ‘G’ stands for ‘geometric’, and gives solutions using the same rules as `solution='geometric'` described above.)

```
sage: qe.solution_extension('G')                        # not tested (random order)
8 x^2 - 8 x - 29 < 0
/\
x^2 - 3 < 0
```

## TESTS:

Check the qepcad configuration file:

```
sage: open('%s/default.qepcadrc'%SAGE_LOCAL).readlines()[-1]
'SINGULAR .../local/bin\n'
```

Tests related to the not tested examples (nondeterministic order of atoms):

```
sage: from sage.interfaces.qepcad import _qepcad_atoms
sage: var('a,b,c,d,x,y,z')
(a, b, c, d, x, y, z)
sage: qf = qepcad_formula
sage: ellipse = 3*x^2 + 2*x*y + y^2 - x + y - 7
sage: circle = x^2 + y^2 - 3

sage: F = qf.exactly_k(3, y, circle * ellipse == 0)
sage: _qepcad_atoms(qepcad(F))                          # optional - qepcad
```

```

{'8 x^2 - 8 x - 29 <= 0',
 '8 x^2 - 8 x - 29 = 0',
 '8 x^4 - 26 x^2 - 4 x + 13 = 0',
 '8 x^4 - 26 x^2 - 4 x + 13 >= 0',
 'x^2 - 3 <= 0',
 'x^2 - 3 = 0'}
sage: _gepcad_atoms(gepcad(F, solution='geometric'))           # optional - gepcad
{'8 x^4 - 26 x^2 - 4 x + 13 = 0',
 'x = _root_-1 x^2 - 3',
 'x = _root_1 8 x^2 - 8 x - 29'}

sage: F = qf.and_(ellipse < 0, circle < 0)
sage: _gepcad_atoms(gepcad(F))                                 # optional - gepcad
{'y^2 + 2 x y + y + 3 x^2 - x - 7 < 0', 'y^2 + x^2 - 3 < 0'}
sage: _gepcad_atoms(gepcad(F, solution='geometric'))         # optional - gepcad
{'8 x^4 - 26 x^2 - 4 x + 13 < 0',
 'x < _root_-2 8 x^4 - 26 x^2 - 4 x + 13',
 'x = _root_-2 8 x^4 - 26 x^2 - 4 x + 13',
 'x = _root_2 8 x^4 - 26 x^2 - 4 x + 13',
 'x > _root_2 8 x^4 - 26 x^2 - 4 x + 13',
 'y^2 + 2 x y + y + 3 x^2 - x - 7 < 0',
 'y^2 + x^2 - 3 < 0'}

sage: F = qf.exists(y, qf.and_(circle == 0, x + y > 0))
sage: _gepcad_atoms(gepcad(F))                                 # optional - gepcad
{'2 x^2 - 3 < 0', 'x > 0', 'x^2 - 3 <= 0'}
sage: _gepcad_atoms(gepcad(F, solution='extended'))          # optional - gepcad
{'x > _root_1 2 x^2 - 3', 'x^2 - 3 <= 0'}

sage: qe = gepcad(qf.and_(ellipse == 0, circle == 0), interact=True) # optional - gepcad
sage: qe.go(); qe.go(); qe.go()                                   # optional - gepcad
QEPCAD object has moved to phase 'Before Projection (y)'
QEPCAD object has moved to phase 'Before Choice'
QEPCAD object has moved to phase 'Before Solution'
sage: for c in qe.cell():                                         # optional - gepcad
....:     count_ellipse = 0
....:     count_circle = 0
....:     for c2 in c:
....:         count_ellipse += (c2.signs()[1][0] == 0)
....:         count_circle += (c2.signs()[1][1] == 0)
....:     c.set_truth(count_ellipse == 2 and count_circle == 2)
sage: _gepcad_atoms(qe.solution_extension('G'))                 # optional - gepcad
{'8 x^2 - 8 x - 29 < 0', 'x^2 - 3 < 0'}

```

**AUTHORS:**

- Carl Witty (2008-03): initial version
- Thierry Monteil (2015-07) repackaging + noncommutative doctests.

**class** sage.interfaces.gepcad.**Qepcad** (formula, vars=None, logfile=None, verbose=False, mem-cells=None, server=None)

The wrapper for QEPCAD.

**answer** ()

For a QEPCAD instance which is finished, return the simplified quantifier-free formula that it printed just before exiting.

EXAMPLES:

```

sage: qe = qepcad(x^3 - x == 0, interact=True) # optional - qepcad
sage: qe.finish() # not tested (random order)
x - 1 <= 0 /\ x + 1 >= 0 /\ [ x = 0 \/ x - 1 = 0 \/ x + 1 = 0 ]
sage: qe.answer() # not tested (random order)
x - 1 <= 0 /\ x + 1 >= 0 /\ [ x = 0 \/ x - 1 = 0 \/ x + 1 = 0 ]

```

#### TESTS:

Tests related to the not tested examples (nondeterministic order of atoms):

```

sage: from sage.interfaces.qepcad import _qepcad_atoms
sage: qe = qepcad(x^3 - x == 0, interact=True) # optional - qepcad
sage: qe.finish() # random, optional - qepcad
x - 1 <= 0 /\ x + 1 >= 0 /\ [ x = 0 \/ x - 1 = 0 \/ x + 1 = 0 ]
sage: _qepcad_atoms(qe.answer()) # optional - qepcad
{'x + 1 = 0', 'x + 1 >= 0', 'x - 1 <= 0', 'x - 1 = 0', 'x = 0'}

```

#### **assume** (*assume*)

The following documentation is from `qepcad.help`.

Add an assumption to the problem. These will not be included in the solution formula.

For example, with input (E x)[ a x<sup>2</sup> + b x + c = 0], if we issue the command

```
assume [ a /= 0 ]
```

we will get the solution formula  $b^2 - 4ac \geq 0$ . Without the assumption we'd get something like  $[a = 0 / b \neq 0] / [a \neq 0 / 4ac - b^2 \leq 0] / [a = 0 / b = 0 / c = 0]$ .

#### EXAMPLES:

```

sage: var('a,b,c,x')
(a, b, c, x)
sage: qf = qepcad_formula
sage: qe = qepcad(qf.exists(x, a*x^2 + b*x + c == 0), interact=True) # optional - qepcad
sage: qe.assume(a != 0) # optional - qepcad
sage: qe.finish() # optional - qepcad
4 a c - b^2 <= 0

```

#### **cell** (*\*index*)

Given a cell index, returns a `QepcadCell` wrapper for that cell. Uses a cache for efficiency.

#### EXAMPLES:

```

sage: qe = qepcad(x + 3 == 42, interact=True) # optional - qepcad
sage: qe.go(); qe.go(); qe.go() # optional - qepcad
QEPCAD object has moved to phase 'At the end of projection phase'
QEPCAD object has moved to phase 'Before Choice'
QEPCAD object has moved to phase 'Before Solution'
sage: qe.cell(2) # optional - qepcad
QEPCAD cell (2)
sage: qe.cell(2) is qe.cell(2) # optional - qepcad
True

```

#### **final\_stats** ()

For a QEPCAD instance which is finished, return the statistics that it printed just before exiting.

#### EXAMPLES:

```

sage: qe = qepcad(x == 0, interact=True) # optional - qepcad
sage: qe.finish() # optional - qepcad
x = 0

```

```
sage: qe.final_stats() # random, optional - qepcad
```

```
-----
0 Garbage collections, 0 Cells and 0 Arrays reclaimed, in 0 milliseconds.
492840 Cells in AVAIL, 500000 Cells in SPACE.
System time: 8 milliseconds.
System time after the initialization: 4 milliseconds.
-----
```

### **make\_cells** (*text*)

Given the result of some QEPCAD command that returns cells (such as `d_cell()`, `d_witness_list()`, etc.), return a list of cell objects.

#### EXAMPLES:

```
sage: var('x,y')
```

```
(x, y)
```

```
sage: qe = qepcad(x^2 + y^2 == 1, interact=True) # optional - qepcad
```

```
sage: qe.go(); qe.go(); qe.go() # optional - qepcad
```

```
QEPCAD object has moved to phase 'Before Projection (y)'
```

```
QEPCAD object has moved to phase 'Before Choice'
```

```
QEPCAD object has moved to phase 'Before Solution'
```

```
sage: qe.make_cells(qe.d_false_cells()) # optional - qepcad
```

```
[QEPCAD cell (5, 1), QEPCAD cell (4, 3), QEPCAD cell (4, 1), QEPCAD cell (3, 5), QEPCAD cell
```

### **phase** ()

Return the current phase of the QEPCAD program.

#### EXAMPLES:

```
sage: qe = qepcad(x > 2/3, interact=True) # optional - qepcad
```

```
sage: qe.phase() # optional - qepcad
```

```
'Before Normalization'
```

```
sage: qe.go() # optional - qepcad
```

```
QEPCAD object has moved to phase 'At the end of projection phase'
```

```
sage: qe.phase() # optional - qepcad
```

```
'At the end of projection phase'
```

```
sage: qe.go() # optional - qepcad
```

```
QEPCAD object has moved to phase 'Before Choice'
```

```
sage: qe.phase() # optional - qepcad
```

```
'Before Choice'
```

```
sage: qe.go() # optional - qepcad
```

```
QEPCAD object has moved to phase 'Before Solution'
```

```
sage: qe.phase() # optional - qepcad
```

```
'Before Solution'
```

```
sage: qe.go() # optional - qepcad
```

```
3 x - 2 > 0
```

```
sage: qe.phase() # optional - qepcad
```

```
'EXITED'
```

### **set\_truth\_value** (*index*, *nv*)

Given a cell index (or a cell) and an integer, set the truth value of the cell to that integer.

Valid integers are 0 (false), 1 (true), and 2 (undetermined).

#### EXAMPLES:

```
sage: qe = qepcad(x == 1, interact=True) # optional - qepcad
```

```
sage: qe.go(); qe.go(); qe.go() # optional - qepcad
```

```
QEPCAD object has moved to phase 'At the end of projection phase'
```

```
QEPCAD object has moved to phase 'Before Choice'
```

QEPCAD object has moved to phase 'Before Solution'  
**sage:** `qe.set_truth_value(1, 1) # optional - qepcad`

### **solution\_extension** (*kind*)

The following documentation is modified from `qepcad.help`:

solution-extension x

Use an alternative solution formula construction method. The parameter x is allowed to be T,E, or G. If x is T, then a formula in the usual language of Tarski formulas is produced. If x is E, a formula in the language of Extended Tarski formulas is produced. If x is G, then a geometry-based formula is produced.

#### EXAMPLES:

```
sage: var('x,y')
(x, y)
sage: qf = qepcad_formula
sage: qe = qepcad(qf.and_(x^2 + y^2 - 3 == 0, x + y > 0), interact=True) # optional - qepcad
sage: qe.go(); qe.go(); qe.go() # optional - qepcad
QEPCAD object has moved to phase 'Before Projection (y)'
QEPCAD object has moved to phase 'Before Choice'
QEPCAD object has moved to phase 'Before Solution'
sage: qe.solution_extension('E') # not tested (random order)
x > _root_1 2 x^2 - 3 /\ y^2 + x^2 - 3 = 0 /\ [ 2 x^2 - 3 > 0 \/ y = _root_-1 y^2 + x^2 - 3
sage: qe.solution_extension('G') # not tested (random order)
[
  [
    2 x^2 - 3 < 0
    \/
    x = _root_-1 2 x^2 - 3
  ]
  /\
  y = _root_-1 y^2 + x^2 - 3
]
\/
[
  x^2 - 3 <= 0
  /\
  x > _root_-1 2 x^2 - 3
  /\
  y^2 + x^2 - 3 = 0
]
sage: qe.solution_extension('T') # not tested (random order)
y + x > 0 /\ y^2 + x^2 - 3 = 0
```

#### TESTS:

Tests related to the not tested examples (nondeterministic order of atoms):

```
sage: from sage.interfaces.qepcad import _qepcad_atoms
sage: var('x,y')
(x, y)
sage: qf = qepcad_formula
sage: qe = qepcad(qf.and_(x^2 + y^2 - 3 == 0, x + y > 0), interact=True) # optional - qepcad
sage: qe.go(); qe.go(); qe.go() # optional - qepcad
QEPCAD object has moved to phase 'Before Projection (y)'
QEPCAD object has moved to phase 'Before Choice'
QEPCAD object has moved to phase 'Before Solution'
sage: _qepcad_atoms(qe.solution_extension('E')) # optional - qepcad
{'2 x^2 - 3 > 0',
```

```

'x > _root_-1 2 x^2 - 3',
'y = _root_-1 y^2 + x^2 - 3',
'y^2 + x^2 - 3 = 0'}
sage: _qepcad_atoms(qe.solution_extension('G'))      # optional - qepcad
{'2 x^2 - 3 < 0',
'x = _root_-1 2 x^2 - 3',
'x > _root_-1 2 x^2 - 3',
'x^2 - 3 <= 0',
'y = _root_-1 y^2 + x^2 - 3',
'y^2 + x^2 - 3 = 0'}
sage: _qepcad_atoms(qe.solution_extension('T'))      # optional - qepcad
{'y + x > 0', 'y^2 + x^2 - 3 = 0'}

```

### trait\_names()

Return a list of the QEPCAD commands which are available as extra methods on a `Qepcad` object.

#### EXAMPLES:

```

sage: qe = qepcad(x^2 < 0, interact=True) # optional - qepcad
sage: len(qe.trait_names()) # random, optional - qepcad
97
sage: 'd_cell' in qe.trait_names() # optional - qepcad
True

```

**class** sage.interfaces.qepcad.**QepcadCell**(parent, lines)

A wrapper for a QEPCAD cell.

### index()

Give the index of a QEPCAD cell.

#### EXAMPLES:

```

sage: var('x,y')
(x, y)
sage: qe = qepcad(x^2 + y^2 == 1, interact=True) # optional - qepcad
sage: qe.go(); qe.go(); qe.go() # optional - qepcad
QEPCAD object has moved to phase 'Before Projection (y)'
QEPCAD object has moved to phase 'Before Choice'
QEPCAD object has moved to phase 'Before Solution'
sage: qe.cell().index() # optional - qepcad
()
sage: qe.cell(1).index() # optional - qepcad
(1,)
sage: qe.cell(2, 2).index() # optional - qepcad
(2, 2)

```

### level()

Return the level of a QEPCAD cell.

#### EXAMPLES:

```

sage: var('x,y')
(x, y)
sage: qe = qepcad(x^2 + y^2 == 1, interact=True) # optional - qepcad
sage: qe.go(); qe.go(); qe.go() # optional - qepcad
QEPCAD object has moved to phase 'Before Projection (y)'
QEPCAD object has moved to phase 'Before Choice'
QEPCAD object has moved to phase 'Before Solution'
sage: qe.cell().level() # optional - qepcad
0

```

```

sage: qe.cell(1).level() # optional - qepcad
1
sage: qe.cell(2, 2).level() # optional - qepcad
2

```

**number\_of\_children()**

Return the number of elements in the stack over a QEPCAD cell. (This is always an odd number, if the stack has been constructed.)

**EXAMPLES:**

```

sage: var('x,y')
(x, y)
sage: qe = qepcad(x^2 + y^2 == 1, interact=True) # optional - qepcad
sage: qe.go(); qe.go(); qe.go() # optional - qepcad
QEPCAD object has moved to phase 'Before Projection (y)'
QEPCAD object has moved to phase 'Before Choice'
QEPCAD object has moved to phase 'Before Solution'
sage: qe.cell().number_of_children() # optional - qepcad
5
sage: [c.number_of_children() for c in qe.cell()] # optional - qepcad
[1, 3, 5, 3, 1]

```

**sample\_point()**

Return the coordinates of a point in the cell, as a tuple of sage algebraic reals.

**EXAMPLES:**

```

sage: qe = qepcad(x^2 - x - 1 == 0, interact=True) # optional - qepcad
sage: qe.go(); qe.go(); qe.go() # optional - qepcad
QEPCAD object has moved to phase 'At the end of projection phase'
QEPCAD object has moved to phase 'Before Choice'
QEPCAD object has moved to phase 'Before Solution'
sage: v1 = qe.cell(2).sample_point()[0]; v1 # optional - qepcad
-0.618033988749895?
sage: v2 = qe.cell(4).sample_point()[0]; v2 # optional - qepcad
1.618033988749895?
sage: v1 + v2 == 1 # optional - qepcad
True

```

**sample\_point\_dict()**

Return the coordinates of a point in the cell, as a dictionary mapping variable names (as strings) to sage algebraic reals.

**EXAMPLES:**

```

sage: qe = qepcad(x^2 - x - 1 == 0, interact=True) # optional - qepcad
sage: qe.go(); qe.go(); qe.go() # optional - qepcad
QEPCAD object has moved to phase 'At the end of projection phase'
QEPCAD object has moved to phase 'Before Choice'
QEPCAD object has moved to phase 'Before Solution'
sage: qe.cell(4).sample_point_dict() # optional - qepcad
{'x': 1.618033988749895?}

```

**set\_truth(v)**

Set the truth value of this cell, as used by QEPCAD for solution formula construction.

The argument *v* should be either a boolean or None (which will set the truth value to 'undetermined').



## EXAMPLES:

```

sage: var('x,y')
(x, y)
sage: qe = qepcad(x^2 + y^2 == 1, interact=True) # optional - qepcad
sage: qe.go(); qe.go(); qe.go() # optional - qepcad
QEPCAD object has moved to phase 'Before Projection (y)'
QEPCAD object has moved to phase 'Before Choice'
QEPCAD object has moved to phase 'Before Solution'
sage: qe.solution_extension('T') # optional - qepcad
y^2 + x^2 - 1 = 0
sage: qe.cell(3, 3).set_truth(True) # optional - qepcad
sage: qe.solution_extension('T') # optional - qepcad
y^2 + x^2 - 1 <= 0

```

**signs()**

Return the sign vector of a QEPCAD cell.

This is a list of lists. The outer list contains one element for each level of the cell; the inner list contains one element for each projection factor at that level. These elements are either -1, 0, or 1.

## EXAMPLES:

```

sage: var('x,y')
(x, y)
sage: qe = qepcad(x^2 + y^2 == 1, interact=True) # optional - qepcad
sage: qe.go(); qe.go(); qe.go() # optional - qepcad
QEPCAD object has moved to phase 'Before Projection (y)'
QEPCAD object has moved to phase 'Before Choice'
QEPCAD object has moved to phase 'Before Solution'
sage: from sage.interfaces.qepcad import QepcadCell
sage: all_cells = flatten(qe.cell(), ltypes=QepcadCell, max_level=1) # optional - qepcad
sage: [(c, c.signs()[1][0]) for c in all_cells] # optional - qepcad
[(QEPCAD cell (1, 1), 1), (QEPCAD cell (2, 1), 1), (QEPCAD cell (2, 2), 0), (QEPCAD cell (2,

```

**class** sage.interfaces.qepcad.**QepcadFunction**(parent, name)

Bases: sage.interfaces.expect.ExpectFunction

A wrapper for a QEPCAD command.

**class** sage.interfaces.qepcad.**Qepcad\_expect**(memcells=None, maxread=100000, log-file=None, server=None)

Bases: sage.interfaces.expect.Expect

The low-level wrapper for QEPCAD.

sage.interfaces.qepcad.**qepcad**(formula, assume=None, interact=False, solution=None, vars=None, \*\*kwargs)

Quantifier elimination and formula simplification using QEPCAD B.

If `assume` is specified, then the given formula is 'assumed', which is taken into account during final solution formula construction.

If `interact=True` is given, then a `Qepcad` object is returned which can be interacted with either at the command line or programmatically.

The type of solution returned can be adjusted with `solution`. The options are 'geometric', which tries to construct a solution formula with geometric meaning; 'extended', which gives a solution formula in an extended language that may be more efficient to construct; 'any-point', which returns any point where the formula is true; 'all-points', which returns a list of all points where the formula is true (or raises an exception if there are infinitely many); and 'cell-points', which returns one point in each cell where the formula is true.

All other keyword arguments are passed through to the `Qepcad` constructor.

For much more documentation and many more examples, see the module docstring for this module (type `sage.interfaces.qepcad?` to read this docstring from the sage command line).

The examples below require that the optional qepcad package is installed.

#### EXAMPLES:

```
sage: qf = qepcad_formula
```

```
sage: var('a,b,c,d,x,y,z,long_with_underscore_314159')
(a, b, c, d, x, y, z, long_with_underscore_314159)
```

```
sage: K.<q,r> = QQ[]
```

```
sage: qepcad('(E x)[a x + b > 0]', vars='(a,b,x)')          # not tested (random order)
a /= 0 /\ b > 0
```

```
sage: qepcad(a > b)                                          # optional - qepcad
b - a < 0
```

```
sage: qepcad(qf.exists(x, a*x^2 + b*x + c == 0))          # not tested (random order)
4 a c - b^2 <= 0 /\ [ c = 0 /\ a /= 0 /\ 4 a c - b^2 < 0 ]
```

```
sage: qepcad(qf.exists(x, a*x^2 + b*x + c == 0), assume=(a != 0)) # optional - qepcad
4 a c - b^2 <= 0
```

For which values of  $a, b, c$  does  $ax^2 + bx + c$  have 2 real zeroes?

```
sage: exact2 = qepcad(qf.exactly_k(2, x, a*x^2 + b*x + c == 0)); exact2 # not tested (random order)
a /= 0 /\ 4 a c - b^2 < 0
```

one real zero?

```
sage: exact1 = qepcad(qf.exactly_k(1, x, a*x^2 + b*x + c == 0)); exact1 # not tested (random order)
[ a > 0 /\ 4 a c - b^2 = 0 ] /\ [ a < 0 /\ 4 a c - b^2 = 0 ] /\ [ a = 0 /\ 4 a c - b^2 < 0 ]
```

No real zeroes?

```
sage: exact0 = qepcad(qf.forall(x, a*x^2 + b*x + c != 0)); exact0 # not tested (random order)
4 a c - b^2 >= 0 /\ c /= 0 /\ [ b = 0 /\ 4 a c - b^2 > 0 ]
```

$3^{75}$  real zeroes?

```
sage: qepcad(qf.exactly_k(3^75, x, a*x^2 + b*x + c == 0)) # optional - qepcad
FALSE
```

We can check that the results don't overlap:

```
sage: qepcad(r'[[%s] /\ [%s]]' % (exact0, exact1), vars='a,b,c') # not tested (random order)
FALSE
```

```
sage: qepcad(r'[[%s] /\ [%s]]' % (exact0, exact2), vars='a,b,c') # not tested (random order)
FALSE
```

```
sage: qepcad(r'[[%s] /\ [%s]]' % (exact1, exact2), vars='a,b,c') # not tested (random order)
FALSE
```

and that the union of the results is as expected:

```
sage: qepcad(r'[[%s] \/ [%s] \/ [%s]]' % (exact0, exact1, exact2), vars=(a,b,c)) # not tested (random order)
b /= 0 /\ a /= 0 /\ c /= 0
```

So we have finitely many zeroes if  $a, b$ , or  $c$  is nonzero; which means we should have infinitely many zeroes if they are all zero.

```
sage: qepcad(qf.infinitely_many(x, a*x^2 + b*x + c == 0))           # not tested (random order)
a = 0 /\ b = 0 /\ c = 0
```

The polynomial is nonzero almost everywhere iff it is not identically zero.

```
sage: qepcad(qf.all_but_finitely_many(x, a*x^2 + b*x + c != 0))    # not tested (random order)
b /= 0 /\ a /= 0 /\ c /= 0
```

The non-zeroes are continuous iff there are no zeroes or if the polynomial is zero.

```
sage: qepcad(qf.connected_subset(x, a*x^2 + b*x + c != 0))        # not tested (random order)
4 a c - b^2 >= 0 /\ [ a = 0 /\ 4 a c - b^2 > 0 ]
```

The zeroes are continuous iff there are no or one zeroes, or if the polynomial is zero:

```
sage: qepcad(qf.connected_subset(x, a*x^2 + b*x + c == 0))        # not tested (random order)
a = 0 /\ 4 a c - b^2 >= 0
sage: qepcad(r'[[%s] /\ [%s] /\ [a = 0 /\ b = 0 /\ c = 0]]' % (exact0, exact1), vars='a,b,c')
a = 0 /\ 4 a c - b^2 >= 0
```

Since polynomials are continuous and  $y > 0$  is an open set, they are positive infinitely often iff they are positive at least once.

```
sage: qepcad(qf.infinitely_many(x, a*x^2 + b*x + c > 0))          # not tested (random order)
c > 0 /\ a > 0 /\ 4 a c - b^2 < 0
sage: qepcad(qf.exists(x, a*x^2 + b*x + c > 0))                   # not tested (random order)
c > 0 /\ a > 0 /\ 4 a c - b^2 < 0
```

However, since  $y \geq 0$  is not open, the equivalence does not hold if you replace ‘positive’ with ‘nonnegative’. (We assume  $a \neq 0$  to get simpler formulas.)

```
sage: qepcad(qf.infinitely_many(x, a*x^2 + b*x + c >= 0), assume=(a != 0)) # not tested (random order)
a > 0 /\ 4 a c - b^2 < 0
sage: qepcad(qf.exists(x, a*x^2 + b*x + c >= 0), assume=(a != 0))        # not tested (random order)
a > 0 /\ 4 a c - b^2 <= 0
```

## TESTS:

We verify that long variable names work. (Note that QEPCAD does not support underscores, so they are stripped from the formula.)

```
sage: qepcad(qf.exists(a, a*long_with_underscore_314159 == 1))      # optional - qepcad
longwithunderscore314159 /= 0
```

Tests related to the not tested examples (nondeterministic order of atoms):

```
sage: from sage.interfaces.qepcad import _qepcad_atoms
sage: var('a,b,c,d,x,y,z,long_with_underscore_314159')
(a, b, c, d, x, y, z, long_with_underscore_314159)
sage: K.<q,r> = QQ[]
```

```
sage: _qepcad_atoms(qepcad('(E x)[a x + b > 0]', vars='(a,b,x)'))    # optional - qepcad
{'a /= 0', 'b > 0'}
```

```
sage: _qepcad_atoms(qepcad(qf.exists(x, a*x^2 + b*x + c == 0)))      # optional - qepcad
{'4 a c - b^2 < 0', '4 a c - b^2 <= 0', 'a /= 0', 'c = 0'}
```

```
sage: exact2 = qepcad(qf.exactly_k(2, x, a*x^2 + b*x + c == 0)); _qepcad_atoms(exact2) # optional - qepcad
{'4 a c - b^2 < 0', 'a /= 0'}
```

```
sage: exact1 = qepcad(qf.exactly_k(1, x, a*x^2 + b*x + c == 0)); _qepcad_atoms(exact1) # optional - qepcad
```

```

{'4 a c - b^2 < 0', '4 a c - b^2 = 0', 'a < 0', 'a = 0', 'a > 0'}

sage: exact0 = qepcad(qf.forall(x, a*x^2 + b*x + c != 0)); _qepcad_atoms(exact0) # optional
{'4 a c - b^2 > 0', '4 a c - b^2 >= 0', 'b = 0', 'c != 0'}

sage: qepcad(r'[[%s] /\ [%s]]' % (exact0, exact1), vars='a,b,c') # optional
FALSE
sage: qepcad(r'[[%s] /\ [%s]]' % (exact0, exact2), vars='a,b,c') # optional
FALSE
sage: qepcad(r'[[%s] /\ [%s]]' % (exact1, exact2), vars='a,b,c') # optional
FALSE

sage: _qepcad_atoms(qepcad(r'[[%s] \/ [%s] \/ [%s]]' % (exact0, exact1, exact2), vars=(a,b,c)))
{'a != 0', 'b != 0', 'c != 0'}

sage: _qepcad_atoms(qepcad(qf.infinitely_many(x, a*x^2 + b*x + c == 0))) # optional
{'a = 0', 'b = 0', 'c = 0'}

sage: _qepcad_atoms(qepcad(qf.all_but_finitely_many(x, a*x^2 + b*x + c != 0))) # optional
{'a != 0', 'b != 0', 'c != 0'}

sage: _qepcad_atoms(qepcad(qf.connected_subset(x, a*x^2 + b*x + c != 0))) # optional
{'4 a c - b^2 > 0', '4 a c - b^2 >= 0', 'a = 0'}

sage: _qepcad_atoms(qepcad(qf.connected_subset(x, a*x^2 + b*x + c == 0))) # optional
{'4 a c - b^2 >= 0', 'a = 0'}

sage: _qepcad_atoms(qepcad(r'[[%s] \/ [%s] \/ [a = 0 /\ b = 0 /\ c = 0]]' % (exact0, exact1), vars=(a,b,c)))
{'4 a c - b^2 >= 0', 'a = 0'}

sage: _qepcad_atoms(qepcad(qf.infinitely_many(x, a*x^2 + b*x + c > 0))) # optional
{'4 a c - b^2 < 0', 'a > 0', 'c > 0'}

sage: _qepcad_atoms(qepcad(qf.exists(x, a*x^2 + b*x + c > 0))) # optional
{'4 a c - b^2 < 0', 'a > 0', 'c > 0'}

sage: _qepcad_atoms(qepcad(qf.infinitely_many(x, a*x^2 + b*x + c >= 0), assume=(a != 0))) # optional
{'4 a c - b^2 < 0', 'a > 0'}

sage: _qepcad_atoms(qepcad(qf.exists(x, a*x^2 + b*x + c >= 0), assume=(a != 0))) # optional
{'4 a c - b^2 <= 0', 'a > 0'}

```

sage.interfaces.qepcad.**qepcad\_banner()**

Return the QEPCAD startup banner.

EXAMPLES:

```

sage: from sage.interfaces.qepcad import qepcad_banner
sage: qepcad_banner() # optional - qepcad
=====
Quantifier Elimination
      in
Elementary Algebra and Geometry
      by
Partial Cylindrical Algebraic Decomposition
...
      by
      Hoon Hong
      (hhong@math.ncsu.edu)

```

With contributions by: Christopher W. Brown, George E. Collins, Mark J. Encarnacion, Jeremy R. Johnson, Werner Krandick, Richard Liska, Scott McCallum, Nicolas Robidoux, and Stanly Steinberg

=====

```
sage.interfaces.qepcad.qepcad_console(memcells=None)
Run QEPCAD directly. To exit early, press Control-C.
```

#### EXAMPLES:

```
sage: qepcad_console() # not tested
...
Enter an informal description between '[' and ']':
```

```
class sage.interfaces.qepcad.qepcad_formula_factory
Contains routines to help construct formulas in QEPCAD syntax.
```

#### **A**(*v, formula*)

Given a variable (or list of variables) and a formula, returns the universal quantification of the formula over the variables.

This method is available both as `forall()` and `A()` (the QEPCAD name for this quantifier).

The input formula may be a `qformula` as returned by the methods of `qepcad_formula`, a symbolic equality or inequality, or a polynomial  $p$  (meaning  $p = 0$ ).

#### EXAMPLE:

```
sage: var('a,b')
(a, b)
sage: qf = qepcad_formula
sage: qf.forall(a, a^2 + b > b^2 + a)
(A a) [a^2 + b > b^2 + a]
sage: qf.forall((a, b), a^2 + b^2 > 0)
(A a) (A b) [a^2 + b^2 > 0]
sage: qf.A(b, b^2 != a)
(A b) [b^2 /= a]
```

#### **C**(*v, formula, allow\_multi=False*)

Given a variable and a formula, returns a new formula which is true iff the set of values for the variable at which the original formula was true is connected (including cases where this set is empty or is a single point).

This method is available both as `connected_subset()` and `C()` (the QEPCAD name for this quantifier).

The input formula may be a `qformula` as returned by the methods of `qepcad_formula`, a symbolic equality or inequality, or a polynomial  $p$  (meaning  $p = 0$ ).

#### EXAMPLE:

```
sage: var('a,b')
(a, b)
sage: qf = qepcad_formula
sage: qf.connected_subset(a, a^2 + b > b^2 + a)
(C a) [a^2 + b > b^2 + a]
sage: qf.C(b, b^2 != a)
(C b) [b^2 /= a]
```

**E**(*v, formula*)

Given a variable (or list of variables) and a formula, returns the existential quantification of the formula over the variables.

This method is available both as `exists()` and `E()` (the QEPCAD name for this quantifier).

The input formula may be a `qformula` as returned by the methods of `qepcad_formula`, a symbolic equality or inequality, or a polynomial  $p$  (meaning  $p = 0$ ).

EXAMPLE:

```
sage: var('a,b')
(a, b)
sage: qf = qepcad_formula
sage: qf.exists(a, a^2 + b > b^2 + a)
(E a) [a^2 + b > b^2 + a]
sage: qf.exists((a, b), a^2 + b^2 < 0)
(E a) (E b) [a^2 + b^2 < 0]
sage: qf.E(b, b^2 == a)
(E b) [b^2 = a]
```

**F**(*v, formula*)

Given a variable and a formula, returns a new formula which is true iff the original formula was true for infinitely many values of the variable.

This method is available both as `infinitely_many()` and `F()` (the QEPCAD name for this quantifier).

The input formula may be a `qformula` as returned by the methods of `qepcad_formula`, a symbolic equality or inequality, or a polynomial  $p$  (meaning  $p = 0$ ).

EXAMPLE:

```
sage: var('a,b')
(a, b)
sage: qf = qepcad_formula
sage: qf.infinitely_many(a, a^2 + b > b^2 + a)
(F a) [a^2 + b > b^2 + a]
sage: qf.F(b, b^2 != a)
(F b) [b^2 /= a]
```

**G**(*v, formula*)

Given a variable and a formula, returns a new formula which is true iff the original formula was true for all but finitely many values of the variable.

This method is available both as `all_but_finitely_many()` and `G()` (the QEPCAD name for this quantifier).

The input formula may be a `qformula` as returned by the methods of `qepcad_formula`, a symbolic equality or inequality, or a polynomial  $p$  (meaning  $p = 0$ ).

EXAMPLE:

```
sage: var('a,b')
(a, b)
sage: qf = qepcad_formula
sage: qf.all_but_finitely_many(a, a^2 + b > b^2 + a)
(G a) [a^2 + b > b^2 + a]
sage: qf.G(b, b^2 != a)
(G b) [b^2 /= a]
```

**X**(*k*, *v*, *formula*, *allow\_multi=False*)

Given a nonnegative integer *k*, a variable, and a formula, returns a new formula which is true iff the original formula is true for exactly *k* values of the variable.

This method is available both as `exactly_k()` and `X()` (the QEPCAD name for this quantifier).

(Note that QEPCAD does not support  $k = 0$  with this syntax, so if  $k = 0$  is requested we implement it with `forall()` and `not_()`.)

The input formula may be a `qformula` as returned by the methods of `qepcad_formula`, a symbolic equality or inequality, or a polynomial *p* (meaning  $p = 0$ ).

EXAMPLE:

```
sage: var('a,b')
(a, b)
sage: qf = qepcad_formula
sage: qf.exactly_k(1, x, x^2 + a*x + b == 0)
(X1 x) [a x + x^2 + b = 0]
sage: qf.exactly_k(0, b, a*b == 1)
(A b) [~a b = 1]
```

**all\_but\_finitely\_many**(*v*, *formula*)

Given a variable and a formula, returns a new formula which is true iff the original formula was true for all but finitely many values of the variable.

This method is available both as `all_but_finitely_many()` and `G()` (the QEPCAD name for this quantifier).

The input formula may be a `qformula` as returned by the methods of `qepcad_formula`, a symbolic equality or inequality, or a polynomial *p* (meaning  $p = 0$ ).

EXAMPLE:

```
sage: var('a,b')
(a, b)
sage: qf = qepcad_formula
sage: qf.all_but_finitely_many(a, a^2 + b > b^2 + a)
(G a) [a^2 + b > b^2 + a]
sage: qf.G(b, b^2 != a)
(G b) [b^2 /= a]
```

**and\_**(*\*formulas*)

Return the conjunction of its input formulas.

(This method would be named ‘and’ if that were not a Python keyword.)

Each input formula may be a `qformula` as returned by the methods of `qepcad_formula`, a symbolic equality or inequality, or a polynomial *p* (meaning  $p = 0$ ).

EXAMPLES:

```
sage: var('a,b,c,x')
(a, b, c, x)
sage: qf = qepcad_formula
sage: qf.and_(a*b, a*c, b*c != 0)
[a b = 0 /\ a c = 0 /\ b c /= 0]
sage: qf.and_(a*x^2 == 3, qf.or_(a > b, b > c))
[a x^2 = 3 /\ [a > b \/ b > c]]
```

**atomic**(*lhs*, *op*='=', *rhs*=0)

Construct a QEPCAD formula from the given inputs.

INPUT:

- `lhs` – a polynomial, or a symbolic equality or inequality
- `op` – a relational operator, default `'='`
- `rhs` – a polynomial, default 0

If `lhs` is a symbolic equality or inequality, then `op` and `rhs` are ignored.

This method works by printing the given polynomials, so we do not care what ring they are in (as long as they print with integral or rational coefficients).

EXAMPLES:

```
sage: qf = qepcad_formula
sage: var('a,b,c')
(a, b, c)
sage: K.<x,y> = QQ[]
sage: def test_qf(qf):
....:     return qf, qf.vars
sage: test_qf(qf.atomic(a^2 + 17))
(a^2 + 17 = 0, frozenset({'a'}))
sage: test_qf(qf.atomic(a*b*c <= c^3))
(a b c <= c^3, frozenset({'a', 'b', 'c'}))
sage: test_qf(qf.atomic(x+y^2, '!=', a+b))
(y^2 + x /= a + b, frozenset({'a', 'b', 'x', 'y'}))
sage: test_qf(qf.atomic(x, operator.lt))
(x < 0, frozenset({'x'}))
```

**connected\_subset** (*v, formula, allow\_multi=False*)

Given a variable and a formula, returns a new formula which is true iff the set of values for the variable at which the original formula was true is connected (including cases where this set is empty or is a single point).

This method is available both as `connected_subset()` and `C()` (the QEPCAD name for this quantifier).

The input formula may be a `qformula` as returned by the methods of `qepcad_formula`, a symbolic equality or inequality, or a polynomial  $p$  (meaning  $p = 0$ ).

EXAMPLE:

```
sage: var('a,b')
(a, b)
sage: qf = qepcad_formula
sage: qf.connected_subset(a, a^2 + b > b^2 + a)
(C a) [a^2 + b > b^2 + a]
sage: qf.C(b, b^2 != a)
(C b) [b^2 /= a]
```

**exactly\_k** (*k, v, formula, allow\_multi=False*)

Given a nonnegative integer  $k$ , a variable, and a formula, returns a new formula which is true iff the original formula is true for exactly  $k$  values of the variable.

This method is available both as `exactly_k()` and `X()` (the QEPCAD name for this quantifier).

(Note that QEPCAD does not support  $k = 0$  with this syntax, so if  $k = 0$  is requested we implement it with `forall()` and `not_()`.)

The input formula may be a `qformula` as returned by the methods of `qepcad_formula`, a symbolic equality or inequality, or a polynomial  $p$  (meaning  $p = 0$ ).



EXAMPLE:

```
sage: var('a,b')
(a, b)
sage: qf = qepcad_formula
sage: qf.exactly_k(1, x, x^2 + a*x + b == 0)
(X1 x) [a x + x^2 + b = 0]
sage: qf.exactly_k(0, b, a*b == 1)
(A b) [~a b = 1]
```

**exists** (*v*, *formula*)

Given a variable (or list of variables) and a formula, returns the existential quantification of the formula over the variables.

This method is available both as `exists()` and `E()` (the QEPCAD name for this quantifier).

The input formula may be a `qformula` as returned by the methods of `qepcad_formula`, a symbolic equality or inequality, or a polynomial  $p$  (meaning  $p = 0$ ).

EXAMPLE:

```
sage: var('a,b')
(a, b)
sage: qf = qepcad_formula
sage: qf.exists(a, a^2 + b > b^2 + a)
(E a) [a^2 + b > b^2 + a]
sage: qf.exists((a, b), a^2 + b^2 < 0)
(E a) (E b) [a^2 + b^2 < 0]
sage: qf.E(b, b^2 == a)
(E b) [b^2 = a]
```

**forall** (*v*, *formula*)

Given a variable (or list of variables) and a formula, returns the universal quantification of the formula over the variables.

This method is available both as `forall()` and `A()` (the QEPCAD name for this quantifier).

The input formula may be a `qformula` as returned by the methods of `qepcad_formula`, a symbolic equality or inequality, or a polynomial  $p$  (meaning  $p = 0$ ).

EXAMPLE:

```
sage: var('a,b')
(a, b)
sage: qf = qepcad_formula
sage: qf.forall(a, a^2 + b > b^2 + a)
(A a) [a^2 + b > b^2 + a]
sage: qf.forall((a, b), a^2 + b^2 > 0)
(A a) (A b) [a^2 + b^2 > 0]
sage: qf.A(b, b^2 != a)
(A b) [b^2 /= a]
```

**formula** (*formula*)

Constructs a QEPCAD formula from the given input.

INPUTS:

- *formula* – a polynomial, a symbolic equality or inequality, or a list of polynomials, equalities, or inequalities

A polynomial  $p$  is interpreted as the equation  $p = 0$ . A list is interpreted as the conjunction ('and') of the elements.

EXAMPLES:

```
sage: var('a,b,c,x')
(a, b, c, x)
sage: qf = qepcad_formula
sage: qf.formula(a*x + b)
a x + b = 0
sage: qf.formula((a*x^2 + b*x + c, a != 0))
[a x^2 + b x + c = 0 /\ a != 0]
```

**iff** (*f1*, *f2*)

Return the equivalence of its input formulas (that is, given formulas  $P$  and  $Q$ , returns ' $P$  iff  $Q$ ').

The input formulas may be a `qformula` as returned by the methods of `qepcad_formula`, a symbolic equality or inequality, or a polynomial  $p$  (meaning  $p = 0$ ).

EXAMPLES:

```
sage: var('a,b')
(a, b)
sage: qf = qepcad_formula
sage: qf.iff(a, b)
[a = 0 <==> b = 0]
sage: qf.iff(a^2 < b, b^2 < a)
[a^2 < b <==> b^2 < a]
```

**implies** (*f1*, *f2*)

Return the implication of its input formulas (that is, given formulas  $P$  and  $Q$ , returns ' $P$  implies  $Q$ ').

The input formulas may be a `qformula` as returned by the methods of `qepcad_formula`, a symbolic equality or inequality, or a polynomial  $p$  (meaning  $p = 0$ ).

EXAMPLES:

```
sage: var('a,b')
(a, b)
sage: qf = qepcad_formula
sage: qf.implies(a, b)
[a = 0 ==> b = 0]
sage: qf.implies(a^2 < b, b^2 < a)
[a^2 < b ==> b^2 < a]
```

**infinitely\_many** (*v*, *formula*)

Given a variable and a formula, returns a new formula which is true iff the original formula was true for infinitely many values of the variable.

This method is available both as `infinitely_many()` and `F()` (the QEPCAD name for this quantifier).

The input formula may be a `qformula` as returned by the methods of `qepcad_formula`, a symbolic equality or inequality, or a polynomial  $p$  (meaning  $p = 0$ ).

EXAMPLE:

```
sage: var('a,b')
(a, b)
sage: qf = qepcad_formula
sage: qf.infinitely_many(a, a^2 + b > b^2 + a)
(F a) [a^2 + b > b^2 + a]
sage: qf.F(b, b^2 != a)
(F b) [b^2 != a]
```

**not\_**(*formula*)

Return the negation of its input formula.

(This method would be named 'not' if that were not a Python keyword.)

The input formula may be a `qformula` as returned by the methods of `qepcad_formula`, a symbolic equality or inequality, or a polynomial  $p$  (meaning  $p = 0$ ).

EXAMPLES:

```

sage: var('a,b')
(a, b)
sage: qf = qepcad_formula
sage: qf.not_(a > b)
[~a > b]
sage: qf.not_(a^2 + b^2)
[~a^2 + b^2 = 0]
sage: qf.not_(qf.and_(a > 0, b < 0))
[~[a > 0 /\ b < 0]]

```

**or\_**(*\*formulas*)

Return the disjunction of its input formulas.

(This method would be named 'or' if that were not a Python keyword.)

Each input formula may be a `qformula` as returned by the methods of `qepcad_formula`, a symbolic equality or inequality, or a polynomial  $p$  (meaning  $p = 0$ ).

EXAMPLES:

```

sage: var('a,b,c,x')
(a, b, c, x)
sage: qf = qepcad_formula
sage: qf.or_(a*b, a*c, b*c != 0)
[a b = 0 \/ a c = 0 \/ b c != 0]
sage: qf.or_(a*x^2 == 3, qf.and_(a > b, b > c))
[a x^2 = 3 \/ [a > b /\ b > c]]

```

**quantifier**(*kind, v, formula, allow\_multi=True*)

A helper method for building quantified QEPCAD formulas; not expected to be called directly.

Takes the quantifier kind (the string label of this quantifier), a variable or list of variables, and a formula, and returns the quantified formula.

EXAMPLES:

```

sage: var('a,b')
(a, b)
sage: qf = qepcad_formula
sage: qf.quantifier('NOT_A_REAL_QEPCAD_QUANTIFIER', a, a*b==0)
(NOT_A_REAL_QEPCAD_QUANTIFIER a) [a b = 0]
sage: qf.quantifier('FOO', (a, b), a*b)
(FOO a) (FOO b) [a b = 0]

```

`sage.interfaces.qepcad.qepcad_version()`

Return a string containing the current QEPCAD version number.

EXAMPLES:

```

sage: qepcad_version() # random, optional - qepcad
'Version B 1.69, 16 Mar 2012'

```

TESTS:

```
sage: qepcad_version() # optional - qepcad  
      'Version B ..., ...'
```

**class** `sage.interfaces.qepcad.qformula` (*formula*, *vars*, *qvars*=[])

A `qformula` holds a string describing a formula in QEPCAD's syntax, and a set of variables used.

## INTERFACE TO BILL HART'S QUADRATIC SIEVE

`sage.interfaces.qsieve.data_to_list(out, n, time)`

Convert output of Hart's sieve and `n` to a list and time.

INPUT:

- `out` – snapshot of text output of Hart's QuadraticSieve program
- `n` – the integer being factored

OUTPUT:

- `list` – proper factors found so far
- `str` – time information

`sage.interfaces.qsieve.qsieve(n, block=True, time=False, verbose=False)`

Run Hart's quadratic sieve and return the distinct proper factors of the integer `n` that it finds.

CONDITIONS:

The conditions for the quadratic sieve to work are as follows:

- No small factors
- Not a perfect power
- Not prime

If any of these fails, the sieve will also.

INPUT:

- `n` – an integer with at least 40 digits
- `block` – (default: `True`) if `True`, you must wait until the sieve computation is complete before using Sage further. If `False`, Sage will run while the sieve computation runs in parallel. If `q` is the returned object, use `q.quit()` to terminate a running factorization.
- `time` – (default: `False`) if `True`, time the command using the UNIX “time” command (which you might have to install).
- `verbose` – (default: `False`) if `True`, print out verbose logging information about what happened during the Sieve run (for non-blocking Sieve, verbose information is always available via the `log()` method.)

OUTPUT:

- `list` – a list of the distinct proper factors of `n` found
- `str` – the time in cpu seconds that the computation took, as given by the command line `time` command. (If `time` is `False`, this is always an empty string.)

EXAMPLES:

```
sage: k = 19; n = next_prime(10^k)*next_prime(10^(k+1))
sage: factor(n) # (currently) uses PARI
1000000000000000000051 * 1000000000000000000039
sage: v, t = qsieve(n, time=True) # uses qsieve; optional - time
sage: v # optional - time
[1000000000000000000051, 1000000000000000000039]
sage: t # random; optional - time
'0.36 real          0.19 user          0.00 sys'
```

`sage.interfaces.qsieve.qsieve_block(n, time, verbose=False)`

Compute the factorization of  $n$  using Hart's quadratic Sieve blocking until complete.

**class** `sage.interfaces.qsieve.qsieve_nonblock(n, time)`

A non-blocking version of Hart's quadratic sieve.

The sieve starts running when you create the object, but you can still use Sage in parallel.

EXAMPLES:

```
sage: k = 19; n = next_prime(10^k)*next_prime(10^(k+1))
sage: q = qsieve(n, block=False, time=True) # optional - time
sage: q # random output; optional - time
Proper factors so far: []
sage: q # random output; optional - time
([1000000000000000000051, 1000000000000000000039], '0.21')
sage: q.list() # random output; optional - time
[1000000000000000000051, 1000000000000000000039]
sage: q.time() # random output; optional - time
'0.21'

sage: q = qsieve(next_prime(10^20)*next_prime(10^21), block=False)
sage: q # random output
Proper factors so far: [1000000000000000000039, 1000000000000000000117]
sage: q # random output
[1000000000000000000039, 1000000000000000000117]
```

**cputime()**

Return the time in seconds (as a string) that it took to factor  $n$ , or return “?” if the factorization has not completed or the time is unknown.

**done()**

Return True if the sieve process has completed.

**list()**

Return a list of the factors found so far, as Sage integers.

**log()**

Return all output of running the sieve so far.

**n()**

Return the integer that is being factored.

**pid()**

Return the PIN id of the QuadraticSieve process (actually of the time process that spawns the sieve process).

**quit()**

Terminate the QuadraticSieve process, in case you want to give up on computing this factorization.

EXAMPLES:

```
sage: n = next_prime(2^310)*next_prime(2^300)
sage: qs = qsieve(n, block=False)
sage: qs
Proper factors so far: []
sage: qs.quit()
sage: qs
Factorization was terminated early.
```

**time()**

Return the time in seconds (as a string) that it took to factor  $n$ , or return “?” if the factorization has not completed or the time is unknown.

```
sage.interfaces.qsieve.tmpdir()
```





## INTERFACE TO R

The following examples try to follow “An Introduction to R” which can be found at <http://cran.r-project.org/doc/manuals/R-intro.html>.

### EXAMPLES:

Simple manipulations; numbers and vectors

The simplest data structure in R is the numeric vector which consists of an ordered collection of numbers. To create a vector named  $x$  using the R interface in Sage, you pass the R interpreter object a list or tuple of numbers:

```
sage: x = r([10.4, 5.6, 3.1, 6.4, 21.7]); x
[1] 10.4  5.6  3.1  6.4 21.7
```

You can invert elements of a vector  $x$  in R by using the invert operator or by doing  $1/x$ :

```
sage: ~x
[1] 0.09615385 0.17857143 0.32258065 0.15625000 0.04608295
sage: 1/x
[1] 0.09615385 0.17857143 0.32258065 0.15625000 0.04608295
```

The following assignment creates a vector  $y$  with 11 entries which consists of two copies of  $x$  with a 0 in between:

```
sage: y = r([x, 0, x]); y
[1] 10.4  5.6  3.1  6.4 21.7  0.0 10.4  5.6  3.1  6.4 21.7
```

### Vector Arithmetic

The following command generates a new vector  $v$  of length 11 constructed by adding together (element by element)  $2x$  repeated 2.2 times,  $y$  repeated just once, and 1 repeated 11 times:

```
sage: v = 2*x+y+1; v
[1] 32.2 17.8 10.3 20.2 66.1 21.8 22.6 12.8 16.9 50.8 43.5
```

One can compute the sum of the elements of an R vector in the following two ways:

```
sage: sum(x)
[1] 47.2
sage: x.sum()
[1] 47.2
```

One can calculate the sample variance of a list of numbers:

```
sage: ((x-x.mean())^2/(x.length()-1)).sum()
[1] 53.853
sage: x.var()
[1] 53.853
```

```
sage: x.sort()
[1] 3.1  5.6  6.4 10.4 21.7
sage: x.min()
[1] 3.1
sage: x.max()
[1] 21.7
sage: x
[1] 10.4  5.6  3.1  6.4 21.7

sage: r(-17).sqrt()
[1] NaN
sage: r('-17+0i').sqrt()
[1] 0+4.123106i
```

Generating an arithmetic sequence:

```
sage: r('1:10')
[1] 1  2  3  4  5  6  7  8  9 10
```

Because `from` is a keyword in Python, it can't be used as a keyword argument. Instead, `from_` can be passed, and R will recognize it as the correct thing:

```
sage: r.seq(length=10, from_=-1, by=.2)
[1] -1.0 -0.8 -0.6 -0.4 -0.2  0.0  0.2  0.4  0.6  0.8

sage: x = r([10.4, 5.6, 3.1, 6.4, 21.7]);
sage: x.rep(2)
[1] 10.4  5.6  3.1  6.4 21.7 10.4  5.6  3.1  6.4 21.7
sage: x.rep(times=2)
[1] 10.4  5.6  3.1  6.4 21.7 10.4  5.6  3.1  6.4 21.7
sage: x.rep(each=2)
[1] 10.4 10.4  5.6  5.6  3.1  3.1  6.4  6.4 21.7 21.7
```

Missing Values:

```
sage: na = r('NA')
sage: z = r([1, 2, 3, na])
sage: z
[1] 1  2  3 NA
sage: ind = r.is_na(z)
sage: ind
[1] FALSE FALSE FALSE  TRUE
sage: zero = r(0)
sage: zero / zero
[1] NaN
sage: inf = r('Inf')
sage: inf-inf
[1] NaN
sage: r.is_na(inf)
[1] FALSE
sage: r.is_na(inf-inf)
[1] TRUE
sage: r.is_na(zero/zero)
[1] TRUE
sage: r.is_na(na)
[1] TRUE
sage: r.is_nan(inf-inf)
[1] TRUE
```

```
sage: r.is_nan(zero/zero)
[1] TRUE
sage: r.is_nan(na)
[1] FALSE
```

#### Character Vectors:

```
sage: labs = r.paste('c("X","Y")', '1:10', sep='"); labs
[1] "X1" "Y2" "X3" "Y4" "X5" "Y6" "X7" "Y8" "X9" "Y10"
```

#### Index vectors; selecting and modifying subsets of a data set:

```
sage: na = r('NA')
sage: x = r([10.4,5.6,3.1,6.4,21.7,na]); x
[1] 10.4 5.6 3.1 6.4 21.7 NA
sage: x['!is.na(self)']
[1] 10.4 5.6 3.1 6.4 21.7

sage: x = r([10.4,5.6,3.1,6.4,21.7,na]); x
[1] 10.4 5.6 3.1 6.4 21.7 NA
sage: (x+1)['(!is.na(self)) & self>0']
[1] 11.4 6.6 4.1 7.4 22.7
sage: x = r([10.4,-2,3.1,-0.5,21.7,na]); x
[1] 10.4 -2.0 3.1 -0.5 21.7 NA
sage: (x+1)['(!is.na(self)) & self>0']
[1] 11.4 4.1 0.5 22.7
```

#### Distributions:

```
sage: r.options(width="60");
$width
[1] 100

sage: rr = r.dnorm(r.seq(-3,3,0.1))
sage: rr
[1] 0.004431848 0.005952532 0.007915452 0.010420935
[5] 0.013582969 0.017528300 0.022394530 0.028327038
[9] 0.035474593 0.043983596 0.053990967 0.065615815
[13] 0.078950158 0.094049077 0.110920835 0.129517596
[17] 0.149727466 0.171368592 0.194186055 0.217852177
[21] 0.241970725 0.266085250 0.289691553 0.312253933
[25] 0.333224603 0.352065327 0.368270140 0.381387815
[29] 0.391042694 0.396952547 0.398942280 0.396952547
[33] 0.391042694 0.381387815 0.368270140 0.352065327
[37] 0.333224603 0.312253933 0.289691553 0.266085250
[41] 0.241970725 0.217852177 0.194186055 0.171368592
[45] 0.149727466 0.129517596 0.110920835 0.094049077
[49] 0.078950158 0.065615815 0.053990967 0.043983596
[53] 0.035474593 0.028327038 0.022394530 0.017528300
[57] 0.013582969 0.010420935 0.007915452 0.005952532
[61] 0.004431848
```

#### Convert R Data Structures to Python/Sage:

```
sage: rr = r.dnorm(r.seq(-3,3,0.1))
sage: sum(rr._sage_())
9.9772125168981...
```

Or you get a dictionary to be able to access all the information:

```
sage: rs = r.summary(r.c(1,4,3,4,3,2,5,1))
sage: rs
  Min. 1st Qu.  Median    Mean 3rd Qu.    Max.
1.000  1.750   3.000   2.875   4.000   5.000
sage: d = rs._sage_()
sage: d['DATA']
[1, 1.75, 3, 2.875, 4, 5]
sage: d['_Names']
['Min.', '1st Qu.', 'Median', 'Mean', '3rd Qu.', 'Max.']
sage: d['_r_class']
['summaryDefault', 'table']
```

It is also possible to access the plotting capabilities of R through Sage. For more information see the documentation of `r.plot()` or `r.png()`.

#### AUTHORS:

- Mike Hansen (2007-11-01)
- William Stein (2008-04-19)
- Harald Schilly (2008-03-20)
- Mike Hansen (2008-04-19)

**class** `sage.interfaces.r.HelpExpression`

Bases: `str`

Used to improve printing of output of `r.help`.

**class** `sage.interfaces.r.R` (*maxread=100000, script\_subdirectory=None, server\_tmpdir=None, log-file=None, server=None, init\_list\_length=1024, seed=None*)

Bases: `sage.interfaces.expect.Expect`

An interface to the R interpreter.

R is a comprehensive collection of methods for statistics, modelling, bioinformatics, data analysis and much more. For more details, see <http://www.r-project.org/about.html>

Resources:

- <http://r-project.org/> provides more information about R.
- <http://rseek.org/> R's own search engine.

#### EXAMPLES:

```
sage: r.summary(r.c(1,2,3,111,2,3,2,3,2,5,4))
Min. 1st Qu.  Median    Mean 3rd Qu.    Max.
1.00    2.00    3.00   12.55    3.50   111.00
```

#### TESTS:

```
sage: r == loads(dumps(r))
True
```

**available\_packages()**

Returns a list of all available R package names.

This list is not necessarily sorted.

OUTPUT: list of strings

**Note:** This requires an internet connection. The CRAN server that is checked is defined at the top of `sage/interfaces/r.py`.

EXAMPLES:

```
sage: ap = r.available_packages()      # optional - internet
sage: len(ap) > 20                     # optional - internet
True
```

**call** (*function\_name*, \**args*, \*\**kwds*)

This is an alias for `function_call()`.

EXAMPLES:

```
sage: r.call('length', [1,2,3])
[1] 3
```

**chdir** (*dir*)

Changes the working directory to *dir*

INPUT:

- *dir* – the directory to change to.

EXAMPLES:

```
sage: import tempfile
sage: tmpdir = tempfile.mkdtemp()
sage: r.chdir(tmpdir)
```

Check that `tmpdir` and `r.getwd()` refer to the same directory. We need to use `realpath()` in case `$TMPDIR` (by default `/tmp`) is a symbolic link (see [trac ticket #10264](#)).

```
sage: os.path.realpath(tmpdir) == sageobj(r.getwd()) # known bug (:trac:'9970')
True
```

**completions** (*s*)

Return all command names that complete the command starting with the string *s*. This is like typing `s[Ctrl-T]` in the R interpreter.

INPUT:

- *s* – string

OUTPUT: list – a list of strings

EXAMPLES:

```
sage: dummy = r.trait_names(use_disk_cache=False) #clean doctest
sage: r.completions('tes')
['testInheritedMethods', 'testPlatformEquivalence', 'testVirtual']
```

**console** ()

Runs the R console as a separate new R process.

EXAMPLES:

```
sage: r.console() # not tested
R version 2.6.1 (2007-11-26)
Copyright (C) 2007 The R Foundation for Statistical Computing
ISBN 3-900051-07-0
...
```

**convert\_r\_list** (*l*)

Converts an R list to a Python list.

EXAMPLES:

```
sage: s = 'c(".GlobalEnv", "package:stats", "package:graphics", "package:grDevices", \n"pack
sage: r.convert_r_list(s)
['.GlobalEnv',
 'package:stats',
 'package:graphics',
 'package:grDevices',
 'package:utils',
 'package:datasets',
 'package:methods',
 'Autoloads',
 'package:base']
```

**eval** (*code*, *globals*=None, *locals*=None, *synchronize*=True, \**args*, \*\**kws*)

Evaluates a command inside the R interpreter and returns the output as a string.

EXAMPLES:

```
sage: r.eval('1+1')
'[1] 2'
```

**function\_call** (*function*, *args*=None, *kws*=None)

Return the result of calling an R function, with given args and keyword args.

OUTPUT: RElement – an object in R

EXAMPLES:

```
sage: r.function_call('length', args=[ [1,2,3] ])
[1] 3
```

**get** (*var*)

Returns the string representation of the variable var.

INPUT:

- var – a string

OUTPUT: string

EXAMPLES:

```
sage: r.set('a', 2)
sage: r.get('a')
'[1] 2'
```

**help** (*command*)

Returns help string for a given command.

INPUT: - command – a string

OUTPUT: HelpExpression – a subclass of string whose `__repr__` method is `__str__`, so it prints nicely

EXAMPLES:

```
sage: r.help('c')
c                                     package:base                R Documentation
...
.. note::
```

This is similar to typing `r.command?`.

### **install\_packages** (*package\_name*)

Install an R package into Sage's R installation.

EXAMPLES:

```
sage: r.install_packages('aaMI')           # not tested
...
R is free software and comes with ABSOLUTELY NO WARRANTY.
You are welcome to redistribute it under certain conditions.
Type 'license()' or 'licence()' for distribution details.
...
Please restart Sage in order to use 'aaMI'.
```

### **library** (*library\_name*)

Load the library `library_name` into the R interpreter.

This function raises an `ImportError` if the given library is not known.

INPUT:

- `library_name` – string

EXAMPLES:

```
sage: r.library('grid')
sage: 'grid' in r.eval('(.packages())')
True
sage: r.library('foobar')
Traceback (most recent call last):
...
ImportError: ...
```

### **na** ()

Returns the NA in R.

OUTPUT: RElement – an element of R

EXAMPLES:

```
sage: r.na()
[1] NA
```

### **plot** (*\*args, \*\*kws*)

The R plot function. Type `r.help('plot')` for much more extensive documentation about this function. See also below for a brief introduction to more plotting with R.

If one simply wants to view an R graphic, using this function is sufficient (because it calls `dev.off()` to turn off the device).

However, if one wants to save the graphic to a specific file, it should be used as in the example below to write the output.

EXAMPLES:

This example saves a plot to the standard R output, usually a filename like `Rplot001.png` - from the command line, in the current directory, and in the cell directory in the notebook:

```
sage: d=r.setwd('%s'%SAGE_TMP)           # for doctesting only; ignore if you are trying this;
sage: r.plot("1:10")                     # optional -- rgraphics
```

```

null device
      1

```

To save to a specific file name, one should use `png()` to set the output device to that file. If this is done in the notebook, it must be done in the same cell as the plot itself:

```

sage: filename = tmp_filename() + '.png'
sage: r.png(filename="%s"%filename) # Note the double quotes in single quotes!; optional -- rgraphics
NULL
sage: x = r([1,2,3])
sage: y = r([4,5,6])
sage: r.plot(x,y) # optional -- rgraphics
null device
      1
sage: import os; os.unlink(filename) # For doctesting, we remove the file; optional -- rgraphics

```

Please note that for more extensive use of R's plotting capabilities (such as the lattices package), it is advisable to either use an interactive plotting device or to use the notebook. The following examples are not tested, because they differ depending on operating system:

```

sage: r.X11() # not tested - opens interactive device on systems with X11 support
sage: r.quartz() # not tested - opens interactive device on OSX
sage: r.hist("rnorm(100)") # not tested - makes a plot
sage: r.library("lattice") # not tested - loads R lattice plotting package
sage: r.histogram(x = "~ wt | cyl", data="mtcars") # not tested - makes a lattice plot
sage: r.dev_off() # not tested, turns off the interactive viewer

```

In the notebook, one can use `r.png()` to open the device, but would need to use the following since R lattice graphics do not automatically print away from the command line:

```

sage: filename = tmp_filename() + '.png' # Not needed in notebook, used for doctesting
sage: r.png(filename="%s"%filename) # filename not needed in notebook, used for doctesting
NULL
sage: r.library("lattice")
sage: r("print(histogram(~wt | cyl, data=mtcars))") # plot should appear; optional -- rgraphics
sage: import os; os.unlink(filename) # We remove the file for doctesting, not needed in notebook

```

### `png(*args, **kws)`

Creates an R PNG device.

This should primarily be used to save an R graphic to a custom file. Note that when using this in the notebook, one must plot in the same cell that one creates the device. See `r.plot()` documentation for more information about plotting via R in Sage.

These examples won't work on the many platforms where R still gets built without graphics support.

#### EXAMPLES:

```

sage: filename = tmp_filename() + '.png'
sage: r.png(filename="%s"%filename) # optional -- rgraphics
NULL
sage: x = r([1,2,3])
sage: y = r([4,5,6])
sage: r.plot(x,y) # This saves to filename, but is not viewable from command line; optional -- rgraphics
null device
      1
sage: import os; os.unlink(filename) # We remove the file for doctesting; optional -- rgraphics

```

We want to make sure that we actually can view R graphics, which happens differently on different platforms:



```

sage: s = r.eval('capabilities("png")') # Should be on Linux and Solaris
sage: t = r.eval('capabilities("aqua")') # Should be on all supported Mac versions
sage: "TRUE" in s+t                      # optional -- rgraphics
True

```

**read**(filename)

Read filename into the R interpreter by calling R's source function on a read-only file connection.

## EXAMPLES:

```

sage: filename = tmp_filename()
sage: f = open(filename, 'w')
sage: f.write('a <- 2+2\n')
sage: f.close()
sage: r.read(filename)
sage: r.get('a')
'[1] 4'

```

**require**(library\_name)

Load the library library\_name into the R interpreter.

This function raises an ImportError if the given library is not known.

## INPUT:

- library\_name – string

## EXAMPLES:

```

sage: r.library('grid')
sage: 'grid' in r.eval('(.packages())')
True
sage: r.library('foobar')
Traceback (most recent call last):
...
ImportError: ...

```

**set**(var, value)

Set the variable var in R to what the string value evaluates to in R.

## INPUT:

- var – a string
- value – a string

## EXAMPLES:

```

sage: r.set('a', '2 + 3')
sage: r.get('a')
'[1] 5'

```

**set\_seed**(seed=None)

Sets the seed for R interpreter. The seed should be an integer.

## EXAMPLES:

```

sage: r = R()
sage: r.set_seed(1)
1
sage: r.sample("1:10", 5)
[1] 3 4 5 7 2

```

**source** (*s*)

Display the R source (if possible) about the function named *s*.

INPUT:

- *s* – a string representing the function whose source code you want to see

OUTPUT: string – source code

EXAMPLES:

```
sage: print r.source("c")
function (... , recursive = FALSE) .Primitive("c")
```

**trait\_names** (*verbose=True, use\_disk\_cache=True*)

Return list of all R functions.

INPUT:

- *verbose* – bool (default: True); if True, display debugging information
- *use\_disk\_cache* – bool (default: True); if True, use the disk cache of trait names to save time.

OUTPUT: list – list of string

EXAMPLES:

```
sage: t = r.trait_names(verbose=False)
sage: len(t) > 200
True
```

**version** ()

Return the version of R currently running.

OUTPUT: tuple of ints; string

EXAMPLES:

```
sage: r.version() # not tested
((3, 0, 1), 'R version 3.0.1 (2013-05-16)')
sage: rint, rstr = r.version()
sage: rint[0] >= 3
True
sage: rstr.startswith('R version')
True
```

**class** `sage.interfaces.r.RElement` (*parent, value, is\_name=False, name=None*)

Bases: `sage.interfaces.expect.ExpectElement`

**dot\_product** (*other*)

Implements the notation `self . other`.

INPUT:

- *self, other* – R elements

OUTPUT: R element

EXAMPLES:

```
sage: c = r.c(1,2,3,4)
sage: c.dot_product(c.t())
[,1] [,2] [,3] [,4]
[1,]  1   2   3   4
[2,]  2   4   6   8
[3,]  3   6   9  12
```

```
[4,]      4      8     12     16
```

```
sage: v = r([3,-1,8])
sage: v.dot_product(v)
      [ ,1]
[1,]    74
```

**stat\_model(x)**

The tilde regression operator in R.

EXAMPLES:

```
sage: x = r([1,2,3,4,5])
sage: y = r([3,5,7,9,11])
sage: a = r.lm( y.tilde(x) ) # lm( y ~ x )
sage: d = a._sage_()
sage: d['DATA']['coefficients']['DATA'][1]
2
```

**tilde(x)**

The tilde regression operator in R.

EXAMPLES:

```
sage: x = r([1,2,3,4,5])
sage: y = r([3,5,7,9,11])
sage: a = r.lm( y.tilde(x) ) # lm( y ~ x )
sage: d = a._sage_()
sage: d['DATA']['coefficients']['DATA'][1]
2
```

**trait\_names()**

Return a list of all methods of this object.

---

**Note:** Currently returns all R commands.

---

EXAMPLES:

```
sage: a = r([1,2,3])
sage: t = a.trait_names()
sage: len(t) > 200
True
```

**class** sage.interfaces.r.**RFunction**(parent, name, r\_name=None)

Bases: sage.interfaces.expect.ExpectFunction

A Function in the R interface.

INPUT:

- parent – the R interface
- name – the name of the function for Python
- r\_name – the name of the function in R itself (which can have dots in it)

EXAMPLES:

```
sage: length = r.length
sage: type(length)
<class 'sage.interfaces.r.RFunction'>
```

```
sage: loads(dumps(length))
length
```

```
class sage.interfaces.r.RFunctionElement(obj, name)
    Bases: sage.interfaces.expect.FunctionElement
```

```
sage.interfaces.r.is_RElement(x)
    Return True if x is an element in an R interface.
```

INPUT:

- x – object

OUTPUT: bool

EXAMPLES:

```
sage: from sage.interfaces.r import is_RElement
sage: is_RElement(2)
False
sage: is_RElement(r(2))
True
```

```
sage.interfaces.r.r_console()
    Spawn a new R command-line session.
```

EXAMPLES:

```
sage: r_console() # not tested
      R version 2.6.1 (2007-11-26)
      Copyright (C) 2007 The R Foundation for Statistical Computing
      ISBN 3-900051-07-0
      ...
```

```
sage.interfaces.r.r_version()
    Return the R version.
```

EXAMPLES:

```
sage: r_version() # not tested
((3, 0, 1), 'R version 3.0.1 (2013-05-16)')
sage: rint, rstr = r_version()
sage: rint[0] >= 3
True
sage: rstr.startswith('R version')
True
```

```
sage.interfaces.r.reduce_load_R()
    Used for reconstructing a copy of the R interpreter from a pickle.
```

EXAMPLES:

```
sage: from sage.interfaces.r import reduce_load_R
sage: reduce_load_R()
R Interpreter
```

## INTERFACE TO SEVERAL RUBIK'S CUBE SOLVERS.

The first is by Michael Reid, and tries to find an optimal solution given the cube's state, and may take a long time. See [http://www.math.ucf.edu/~reid/Rubik/optimal\\_solver.html](http://www.math.ucf.edu/~reid/Rubik/optimal_solver.html)

The second is by Eric Dietz, and uses a standard (?) algorithm to solve the cube one level at a time. It is extremely fast, but often returns a far from optimal solution. See <http://wrongway.org/?rubiksource>

The third is by Dik Winter and implements Kociemba's algorithm which finds reasonable solutions relatively quickly, and if it is kept running will eventually find the optimal solution.

**AUTHOR:** – Optimal was written by Michael Reid <[reid@math.ucf.edu](mailto:reid@math.ucf.edu)> (2004) – Cubex was written by Eric Dietz <[root@wrongway.org](mailto:root@wrongway.org)> (2003) – Kociemba was written by Dik T. Winter <[dik.winter@cw.nl](mailto:dik.winter@cw.nl)> (1993) – Initial interface by Robert Bradshaw (2007-08)

```
class sage.interfaces.rubik.CubexSolver
```

```
    format_cube (facets)
```

```
    solve (facets)
```

```
    EXAMPLES: sage: from sage.interfaces.rubik import * sage: C = RubiksCube("R U") sage: Cubex-
    Solver().solve(C.facets()) 'R U' sage: C = RubiksCube("R U F L B D") sage: sol = Cubex-
    Solver().solve(C.facets()); sol "U' L' L' U L U' L U D L L D' L' D L' D' L D L' U' L D' L' U
    L' B' U' L' U B L D L D' U' L' U L B L B' L' U L U' L' F' L' F L' F L F' L' D' L' D D L D' B
    L B' L B' L B F' L F F B' L F' B D' D' L D B' B' L' D' B U' U' L' B' D' F' F' L D F" sage:
    RubiksCube(sol) == C True sage: C = RubiksCube("R2 F") sage: CubexSolver().solve(C.facets())
    "R' R' F" sage: C = RubiksCube().scramble() sage: sol = CubexSolver().solve(C.facets()) sage: C
    == RubiksCube(sol) True
```

```
class sage.interfaces.rubik.DikSolver
```

```
    format_cube (facets)
```

```
    solve (facets, timeout=10, extra_time=2)
```

```
    EXAMPLES: sage: from sage.interfaces.rubik import * sage: C = RubiksCube().move("R U") sage:
    DikSolver().solve(C.facets()) 'R U' sage: C = RubiksCube().move("R U F L B D") sage: Dik-
    Solver().solve(C.facets()) 'R U F L B D' sage: C = RubiksCube().move("R2 F") sage: Dik-
    Solver().solve(C.facets()) "R2 F"
```

```
class sage.interfaces.rubik.OptimalSolver (verbose=False, wait=True)
    Interface to Michael Reid's optimal Rubik's Cube solver.
```

```
    format_cube (facets)
```

```
    ready ()
```

**solve** (*facets*)

The initial startup and precomputation are substantial...

TODO: Let it keep searching once it found a solution?

**EXAMPLES:** sage: from sage.interfaces.rubik import \* sage: solver = DikSolver() sage: solver = OptimalSolver() # long time (28s on sage.math, 2012) Initializing tables... Done. sage: C = RubiksCube("R U") sage: solver.solve(C.facets()) 'R U' sage: C = RubiksCube("R U F L B D") sage: solver.solve(C.facets()) 'R U F L B D' sage: C = RubiksCube("R2 D2") sage: solver.solve(C.facets()) 'R2 D2'

**start** ()

**stop** ()

**class** sage.interfaces.rubik.**SingNot** (*s*)

This class is to resolve difference between various Singmaster notation. Case is ignored, and the second and third letters may be swapped.

**EXAMPLE:** sage: from sage.interfaces.rubik import SingNot sage: SingNot("acb") == SingNot("ACB") True  
sage: SingNot("acb") == SingNot("bca") False

## INTERFACE TO SAGE

This is an expect interface to *another* copy of the Sage interpreter.

```
class sage.interfaces.sage0.Sage (logfile=None, preparse=True, python=False, init_code=None,
                                server=None, server_tmpdir=None, remote_cleaner=True,
                                **kwds)
```

Bases: `sage.interfaces.expect.Expect`

Expect interface to the Sage interpreter itself.

INPUT:

- `server` - (optional); if specified runs Sage on a remote machine with address. You must have ssh keys setup so you can login to the remote machine by typing “ssh remote\_machine” and no password, call `_install_hints_ssh()` for hints on how to do that.

The version of Sage should be the same as on the local machine, since pickling is used to move data between the two Sage process.

EXAMPLES: We create an interface to a copy of Sage. This copy of Sage runs as an external process with its own memory space, etc.

```
sage: s = Sage()
```

Create the element 2 in our new copy of Sage, and cube it.

```
sage: a = s(2)
```

```
sage: a^3
```

```
8
```

Create a vector space of dimension 4, and compute its generators:

```
sage: V = s('QQ^4')
```

```
sage: V.gens()
```

```
((1, 0, 0, 0), (0, 1, 0, 0), (0, 0, 1, 0), (0, 0, 0, 1))
```

Note that `V` is not a vector space, it's a wrapper around an object (which happens to be a vector space), in another running instance of Sage.

```
sage: type(V)
```

```
<class 'sage.interfaces.sage0.SageElement'>
```

```
sage: V.parent()
```

```
Sage
```

```
sage: g = V.0; g
```

```
(1, 0, 0, 0)
```

```
sage: g.parent()
```

```
Sage
```

We can still get the actual parent by using the name attribute of `g`, which is the variable name of the object in the child process.

```
sage: s('%s.parent()' % g.name())
Vector space of dimension 4 over Rational Field
```

Note that the memory space is completely different.

```
sage: x = 10
sage: s('x = 5')
5
sage: x
10
sage: s('x')
5
```

We can have the child interpreter itself make another child Sage process, so now three copies of Sage are running:

```
sage: s3 = s('Sage()')
sage: a = s3(10)
sage: a
10
```

This  $a = 10$  is in a subprocess of a subprocesses of your original Sage.

```
sage: _ = s.eval('%s.eval("x=8")' % s3.name())
sage: s3('"x"')
8
sage: s('x')
5
sage: x
10
```

The double quotes are needed because the call to `s3` first evaluates its arguments using the `s` interpreter, so the call to `s3` is passed `s('"x"')`, which is the string `"x"` in the `s` interpreter.

**clear** (*var*)

Clear the variable named `var`.

Note that the exact format of the `NameError` for a cleared variable is slightly platform dependent, see trac #10539.

EXAMPLES:

```
sage: sage0.set('x', '2')
sage: sage0.get('x')
'2'
sage: sage0.clear('x')
sage: 'NameError' in sage0.get('x')
True
```

**console** ()

Spawn a new Sage command-line session.

EXAMPLES:

```
sage: sage0.console() #not tested
```

```
-----
| SageMath Version ..., Release Date: ...                               |
| Type notebook() for the GUI, and license() for information.           |
-----
```



...

**cputime** (*t=None*)

Return cputime since this Sage subprocess was started.

EXAMPLES:

```
sage: sage0.cputime()      # random output
1.3530439999999999
sage: sage0('factor(2^157-1)')
852133201 * 60726444167 * 1654058017289 * 2134387368610417
sage: sage0.cputime()      # random output
1.6462939999999999
```

**eval** (*line, strip=True, \*\*kws*)Send the code *x* to a second instance of the Sage interpreter and return the output as a string.

This allows you to run two completely independent copies of Sage at the same time in a unified way.

INPUT:

- line* - input line of code
- strip* - ignored

EXAMPLES:

```
sage: sage0.eval('2+2')
'4'
```

**get** (*var*)Get the value of the variable *var*.

EXAMPLES:

```
sage: sage0.set('x', '2')
sage: sage0.get('x')
'2'
```

**new** (*x*)

EXAMPLES:

```
sage: sage0.new(2)
2
sage: _.parent()
Sage
```

**preparse** (*x*)Returns the preparsed version of the string *s*.

EXAMPLES:

```
sage: sage0.preparse('2+2')
'Integer(2)+Integer(2)'
```

**set** (*var, value*)Set the variable *var* to the given value.

EXAMPLES:

```
sage: sage0.set('x', '2')
sage: sage0.get('x')
'2'
```

**trait\_names()**

EXAMPLES:

```
sage: t = sage0.trait_names()
sage: len(t) > 100
True
sage: 'gcd' in t
True
```

**version()**

EXAMPLES:

```
sage: sage0.version()
'SageMath Version ..., Release Date: ...'
sage: sage0.version() == version()
True
```

**class** `sage.interfaces.sage0.SageElement` (*parent, value, is\_name=False, name=None*)

Bases: `sage.interfaces.expect.ExpectElement`

**class** `sage.interfaces.sage0.SageFunction` (*obj, name*)

Bases: `sage.interfaces.expect.FunctionElement`

`sage.interfaces.sage0.reduce_load_Sage()`

EXAMPLES:

```
sage: from sage.interfaces.sage0 import reduce_load_Sage
sage: reduce_load_Sage()
Sage
```

`sage.interfaces.sage0.reduce_load_element(s)`

EXAMPLES:

```
sage: from sage.interfaces.sage0 import reduce_load_element
sage: s = dumps(1/2)
sage: half = reduce_load_element(s); half
1/2
sage: half.parent()
Sage
```

`sage.interfaces.sage0.sage0_console()`

Spawn a new Sage command-line session.

EXAMPLES:

```
sage: sage0_console() #not tested
```

```
-----
| SageMath Version ..., Release Date: ...                               |
| Type notebook() for the GUI, and license() for information.           |
|-----|
...

```

`sage.interfaces.sage0.sage0_version()`

EXAMPLES:

```
sage: from sage.interfaces.sage0 import sage0_version
sage: sage0_version() == version()
True
```

## INTERFACE TO SCILAB

Scilab is a scientific software package for numerical computations providing a powerful open computing environment for engineering and scientific applications. Scilab includes hundreds of mathematical functions with the possibility to add interactively programs from various languages (C, C++, Fortran...). It has sophisticated data structures (including lists, polynomials, rational functions, linear systems...), an interpreter and a high level programming language.

The commands in this section only work if you have the “scilab” interpreter installed and available in your PATH. It’s not necessary to install any special Sage packages.

### EXAMPLES:

```
sage: scilab.eval('2+2')           # optional - scilab
'ans  =\n \n      4.'
```

```
sage: scilab('2+2')               # optional - scilab
4.
```

```
sage: a = scilab(10)              # optional - scilab
sage: a**10                       # optional - scilab
1.000D+10
```

Tutorial based the MATLAB interface tutorial:

### EXAMPLES:

```
sage: scilab('4+10')              # optional - scilab
14.
```

```
sage: scilab('date')              # optional - scilab; random output
15-Feb-2010
```

```
sage: scilab('5*10 + 6')           # optional - scilab
56.
```

```
sage: scilab('(6+6)/3')            # optional - scilab
4.
```

```
sage: scilab('9')^2               # optional - scilab
81.
```

```
sage: a = scilab(10); b = scilab(20); c = scilab(30)    # optional - scilab
sage: avg = (a+b+c)/3              # optional - scilab
sage: avg                          # optional - scilab
20.
```

```
sage: parent(avg)                 # optional - scilab
Scilab
```

```
sage: my_scalar = scilab('3.1415') # optional - scilab
sage: my_scalar                    # optional - scilab
3.1415
```

```
sage: my_vector1 = scilab('[1,5,7]') # optional - scilab
sage: my_vector1                  # optional - scilab
1.      5.      7.
```

```
sage: my_vector2 = scilab('[1;5;7]') # optional - scilab
```

```
sage: my_vector2                                # optional - scilab
1.
5.
7.
sage: my_vector1 * my_vector2                    # optional - scilab
75.

sage: row_vector1 = scilab('[1 2 3]')            # optional - scilab
sage: row_vector2 = scilab('[3 2 1]')            # optional - scilab
sage: matrix_from_row_vec = scilab('[%s; %s]'%(row_vector1.name(), row_vector2.name())) # optional - scilab
sage: matrix_from_row_vec                        # optional - scilab
1.      2.      3.
3.      2.      1.

sage: column_vector1 = scilab('[1;3]')           # optional - scilab
sage: column_vector2 = scilab('[2;8]')           # optional - scilab
sage: matrix_from_col_vec = scilab('[%s %s]%(column_vector1.name(), column_vector2.name())') # optional - scilab
sage: matrix_from_col_vec                        # optional - scilab
1.      2.
3.      8.

sage: my_matrix = scilab('[8, 12, 19; 7, 3, 2; 12, 4, 23; 8, 1, 1]') # optional - scilab
sage: my_matrix                                  # optional - scilab
8.      12.     19.
7.      3.      2.
12.     4.      23.
8.      1.      1.

sage: combined_matrix = scilab('[%s, %s]%(my_matrix.name(), my_matrix.name())')
sage: combined_matrix                            # optional - scilab
8.      12.     19.     8.      12.     19.
7.      3.      2.      7.      3.      2.
12.     4.      23.     12.     4.      23.
8.      1.      1.      8.      1.      1.

sage: tm = scilab('0.5:2:10')                    # optional - scilab
sage: tm                                          # optional - scilab
0.5     2.5     4.5     6.5     8.5

sage: my_vector1 = scilab('[1,5,7]')             # optional - scilab
sage: my_vector1(1)                             # optional - scilab
1.
sage: my_vector1(2)                             # optional - scilab
5.
sage: my_vector1(3)                             # optional - scilab
7.
```

Matrix indexing works as follows:

```
sage: my_matrix = scilab('[8, 12, 19; 7, 3, 2; 12, 4, 23; 8, 1, 1]') # optional - scilab
sage: my_matrix(3,2)                                                # optional - scilab
4.
```

One can also use square brackets:

```
sage: my_matrix[3,2]                                                # optional - scilab
4.
```

Setting using parenthesis cannot work (because of how the Python language works). Use square brackets or the set function:

```
sage: my_matrix = scilab('[8, 12, 19; 7, 3, 2; 12, 4, 23; 8, 1, 1]') # optional - scilab
sage: my_matrix.set(2,3, 1999) # optional - scilab
sage: my_matrix # optional - scilab
      8.      12.      19.
      7.       3.     1999.
     12.       4.      23.
      8.       1.       1.
sage: my_matrix[2,3] = -126 # optional - scilab
sage: my_matrix # optional - scilab
      8.      12.      19.
      7.       3.     - 126.
     12.       4.      23.
      8.       1.       1.
```

#### TESTS:

```
sage: M = scilab(x) # optional - scilab
Traceback (most recent call last):
...
TypeError: _interface_init_() takes exactly one argument (0 given)
sage: M = scilab(matrix(3,range(9))); M # optional - scilab
      0.      1.      2.
      3.      4.      5.
      6.      7.      8.
sage: M(10) # optional - scilab
Traceback (most recent call last):
...
TypeError: Error executing code in Scilab
...
Invalid index.
sage: M[10] # optional - scilab
Traceback (most recent call last):
...
TypeError: Error executing code in Scilab
...
Invalid index.
sage: M(4,2) # optional - scilab
Traceback (most recent call last):
...
TypeError: Error executing code in Scilab
...
Invalid index.
sage: M[2,4] # optional - scilab
Traceback (most recent call last):
...
TypeError: Error executing code in Scilab
...
Invalid index.
sage: M(9) = x # optional - scilab
Traceback (most recent call last):
...
SyntaxError: can't assign to function call (... , line 1)
```

#### AUTHORS:

– Ronan Paixao (2008-11-26), based on the MATLAB tutorial by William Stein (2006-10-11)

```
class sage.interfaces.scilab.Scilab(maxread=100, script_subdirectory=None, logfile=None,
                                   server=None, server_tmpdir=None, seed=None)
```

Bases: `sage.interfaces.expect.Expect`

Interface to the Scilab interpreter.

EXAMPLES:

```
sage: a = scilab('[ 1, 1, 2; 3, 5, 8; 13, 21, 33 ]')      # optional - scilab
sage: b = scilab('[ 1; 3; 13]')                        # optional - scilab
sage: c = a * b                                         # optional - scilab
sage: print c                                           # optional - scilab
30.
122.
505.
```

**console()**

Starts Scilab console.

EXAMPLES:

```
sage: scilab.console()                                # optional - scilab; not tested
```

**eval**(command, \*args, \*\*kws)

Evaluates commands.

EXAMPLES:

```
sage: scilab.eval("5")                                # optional - scilab
'ans =
```

5.' sage: scilab.eval("d=44") # optional - scilab 'd =

44.'

**get**(var)

Get the value of the variable var.

EXAMPLES:

```
sage: scilab.eval('b=124;')                            # optional - scilab
''
sage: scilab.get('b')                                  # optional - scilab
'
```

124.'

**sage2scilab\_matrix\_string**(A)

Return a Scilab matrix from a Sage matrix.

**INPUT:** A Sage matrix with entries in the rationals or reals.

**OUTPUT:** A string that evaluates to an Scilab matrix.

EXAMPLES:

```
sage: M33 = MatrixSpace(QQ, 3, 3)                      # optional - scilab
sage: A = M33([1, 2, 3, 4, 5, 6, 7, 8, 0])              # optional - scilab
sage: scilab.sage2scilab_matrix_string(A)              # optional - scilab
'[1, 2, 3; 4, 5, 6; 7, 8, 0]'
```

**set** (*var*, *value*)

Set the variable *var* to the given value.

EXAMPLES:

```
sage: scilab.set('a', 123)           # optional - scilab
sage: scilab.get('a')              # optional - scilab
'
123.'
```

**set\_seed** (*seed=None*)

Sets the seed for gp interpreter. The seed should be an integer.

EXAMPLES:

```
sage: from sage.interfaces.scilab import Scilab # optional - scilab
sage: s = Scilab() # optional - scilab
sage: s.set_seed(1) # optional - scilab
1
sage: [s.rand() for i in range(5)] # optional - scilab
[
    0.6040239,
    0.0079647,
    0.6643966,
    0.9832111,
    0.5321420]
```

**version** ()

Returns the version of the Scilab software used.

EXAMPLES:

```
sage: scilab.version()              # optional - scilab
'scilab-...'
```

**whos** (*name=None*, *typ=None*)

Returns information about current objects. Arguments: *nam*: first characters of selected names *typ*: name of selected Scilab variable type

EXAMPLES:

```
sage: scilab.whos("core")           # optional - scilab
'Name                               Type                               Size                               Bytes...'
sage: scilab.whos(typ='function')   # optional - scilab
'Name                               Type                               Size                               Bytes...'
```

**class** `sage.interfaces.scilab.ScilabElement` (*parent*, *value*, *is\_name=False*, *name=None*)

Bases: `sage.interfaces.expect.ExpectElement`

**set** (*i*, *j*, *x*)

Set the variable *var* to the given value.

EXAMPLES:

```
sage: scilab.set('c', 125)           # optional - scilab
sage: scilab.get('c')              # optional - scilab
,
```

125.'

`sage.interfaces.scilab.scilab_console()`

This requires that the optional Scilab program be installed and in your PATH, but no optional Sage packages need to be installed.

EXAMPLES:

```
sage: from sage.interfaces.scilab import scilab_console # optional - scilab
sage: scilab_console()                                # optional - scilab; not tested
```

---

scilab-5.0.3

Consortium Scilab (DIGITEO)  
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Copyright (c) 1989-2007 (ENPC)

---

Startup execution:

loading initial environment

-->2+3

ans =

5.

-->quit

Typing quit exits the Scilab console and returns you to Sage. Scilab, like Sage, remembers its history from one session to another.

`sage.interfaces.scilab.scilab_version()`

Return the version of Scilab installed.

EXAMPLES:

```
sage: from sage.interfaces.scilab import scilab_version # optional - scilab
sage: scilab_version()    # optional - scilab
'scilab-...'
```



## INTERFACE TO SINGULAR

### AUTHORS:

- David Joyner and William Stein (2005): first version
- Martin Albrecht (2006-03-05): code so `singular.[tab]` and `x = singular(...)`, `x.[tab]` includes all singular commands.
- Martin Albrecht (2006-03-06): This patch adds the equality symbol to singular. Also fix a problem in which `""` as prompt means comparison will break all further communication with Singular.
- Martin Albrecht (2006-03-13): added `current_ring()` and `current_ring_name()`
- William Stein (2006-04-10): Fixed problems with ideal constructor
- Martin Albrecht (2006-05-18): added `sage_poly`.
- Simon King (2010-11-23): Reduce the overhead caused by waiting for the Singular prompt by doing garbage collection differently.
- Simon King (2011-06-06): Make conversion from Singular to Sage more flexible.
- Simon King (2015): Extend pickling capabilities.

## 40.1 Introduction

This interface is extremely flexible, since it's exactly like typing into the Singular interpreter, and anything that works there should work here.

The Singular interface will only work if Singular is installed on your computer; this should be the case, since Singular is included with Sage. The interface offers three pieces of functionality:

1. `singular_console()` - A function that dumps you into an interactive command-line Singular session.
2. `singular(expr, type='def')` - Creation of a Singular object. This provides a Pythonic interface to Singular. For example, if `f=singular(10)`, then `f.factorize()` returns the factorization of 10 computed using Singular.
3. `singular.eval(expr)` - Evaluation of arbitrary Singular expressions, with the result returned as a string.

Of course, there are polynomial rings and ideals in Sage as well (often based on a C-library interface to Singular). One can convert an object in the Singular interpreter interface to Sage by the method `sage()`.

## 40.2 Tutorial

EXAMPLES: First we illustrate multivariate polynomial factorization:

```

sage: R1 = singular.ring(0, '(x,y)', 'dp')
sage: R1
// characteristic : 0
// number of vars : 2
//      block   1 : ordering dp
//              : names   x y
//      block   2 : ordering C
sage: f = singular('9x16 - 18x13y2 - 9x12y3 + 9x10y4 - 18x11y2 + 36x8y4 + 18x7y5 - 18x5y6 + 9x6y4 - 18x4y5 + 9x3y6 - 9x2y7 + 9x1y8 - 9')
sage: f
9*x^16-18*x^13*y^2-9*x^12*y^3+9*x^10*y^4-18*x^11*y^2+36*x^8*y^4+18*x^7*y^5-18*x^5*y^6+9*x^6*y^4-18*x^4*y^5+9*x^3*y^6-9*x^2*y^7+9*x*y^8-9
sage: f.parent()
Singular

```

```

sage: F = f.factorize(); F
[1]:
  _[1]=9
  _[2]=x^6-2*x^3*y^2-x^2*y^3+y^4
  _[3]=-x^5+y^2
[2]:
  1,1,2

```

```

sage: F[1]
9,
x^6-2*x^3*y^2-x^2*y^3+y^4,
-x^5+y^2
sage: F[1][2]
x^6-2*x^3*y^2-x^2*y^3+y^4

```

We can convert  $f$  and each exponent back to Sage objects as well.

```

sage: g = f.sage(); g
9*x^16 - 18*x^13*y^2 - 9*x^12*y^3 + 9*x^10*y^4 - 18*x^11*y^2 + 36*x^8*y^4 + 18*x^7*y^5 - 18*x^5*y^6 + 9*x^6*y^4 - 18*x^4*y^5 + 9*x^3*y^6 - 9*x^2*y^7 + 9*x*y^8 - 9
sage: F[1][2].sage()
x^6 - 2*x^3*y^2 - x^2*y^3 + y^4
sage: g.parent()
Multivariate Polynomial Ring in x, y over Rational Field

```

This example illustrates polynomial GCD's:

```

sage: R2 = singular.ring(0, '(x,y,z)', 'lp')
sage: a = singular.new('3x2*(x+y)')
sage: b = singular.new('9x*(y2-x2)')
sage: g = a.gcd(b)
sage: g
x^2+x*y

```

This example illustrates computation of a Groebner basis:

```

sage: R3 = singular.ring(0, '(a,b,c,d)', 'lp')
sage: I = singular.ideal(['a + b + c + d', 'a*b + a*d + b*c + c*d', 'a*b*c + a*b*d + a*c*d + b*c*d', 'a^2*d^6-c^2*d^2-d^4+1,
sage: I2 = I.groebner()
sage: I2
c^2*d^6-c^2*d^2-d^4+1,
c^3*d^2+c^2*d^3-c-d,
b*d^4-b+d^5-d,
b*c-b*d^5+c^2*d^4+c*d-d^6-d^2,
b^2+2*b*d+d^2,

```

$a+b+c+d$

The following example is the same as the one in the Singular - Gap interface documentation:

```
sage: R = singular.ring(0, '(x0,x1,x2)', 'lp')
sage: I1 = singular.ideal(['x0*x1*x2 -x0^2*x2', 'x0^2*x1*x2-x0*x1^2*x2-x0*x1*x2^2', 'x0*x1-x0*x2-x1*x2'])
sage: I2 = I1.groebner()
sage: I2
x1^2*x2^2,
x0*x2^3-x1^2*x2^2+x1*x2^3,
x0*x1-x0*x2-x1*x2,
x0^2*x2-x0*x2^2-x1*x2^2
sage: I2.sage()
Ideal (x1^2*x2^2, x0*x2^3 - x1^2*x2^2 + x1*x2^3, x0*x1 - x0*x2 - x1*x2, x0^2*x2 - x0*x2^2 - x1*x2^2)
```

This example illustrates moving a polynomial from one ring to another. It also illustrates calling a method of an object with an argument.

```
sage: R = singular.ring(0, '(x,y,z)', 'dp')
sage: f = singular('x3+y3+(x-y)*x2y2+z2')
sage: f
x^3*y^2-x^2*y^3+x^3+y^3+z^2
sage: R1 = singular.ring(0, '(x,y,z)', 'ds')
sage: f = R.fetch(f)
sage: f
z^2+x^3+y^3+x^3*y^2-x^2*y^3
```

We can calculate the Milnor number of  $f$ :

```
sage: _=singular.LIB('sing.lib')      # assign to _ to suppress printing
sage: f.milnor()
4
```

The Jacobian applied twice yields the Hessian matrix of  $f$ , with which we can compute.

```
sage: H = f.jacob().jacob()
sage: H
6*x+6*x*y^2-2*y^3, 6*x^2*y-6*x*y^2, 0,
6*x^2*y-6*x*y^2, 6*y+2*x^3-6*x^2*y, 0,
0, 0, 2
sage: H.sage()
[6*x + 6*x*y^2 - 2*y^3      6*x^2*y - 6*x*y^2      0]
[      6*x^2*y - 6*x*y^2  6*y + 2*x^3 - 6*x^2*y      0]
[      0      0      0      2]
sage: H.det()      # This is a polynomial in Singular
72*x*y+24*x^4-72*x^3*y+72*x*y^3-24*y^4-48*x^4*y^2+64*x^3*y^3-48*x^2*y^4
sage: H.det().sage()      # This is the corresponding polynomial in Sage
72*x*y + 24*x^4 - 72*x^3*y + 72*x*y^3 - 24*y^4 - 48*x^4*y^2 + 64*x^3*y^3 - 48*x^2*y^4
```

The 1x1 and 2x2 minors:

```
sage: H.minor(1)
2,
6*y+2*x^3-6*x^2*y,
6*x^2*y-6*x*y^2,
6*x^2*y-6*x*y^2,
6*x+6*x*y^2-2*y^3
sage: H.minor(2)
```

```
12*y+4*x^3-12*x^2*y,
12*x^2*y-12*x*y^2,
12*x^2*y-12*x*y^2,
12*x+12*x*y^2-4*y^3,
-36*x*y-12*x^4+36*x^3*y-36*x*y^3+12*y^4+24*x^4*y^2-32*x^3*y^3+24*x^2*y^4
```

```
sage: _=singular.eval('option(redSB)')
sage: H.minor(1).groebner()
1
```

## 40.3 Computing the Genus

We compute the projective genus of ideals that define curves over  $\mathbb{Q}$ . It is *very important* to load the `normal.lib` library before calling the `genus` command, or you'll get an error message.

EXAMPLE:

```
sage: singular.lib('normal.lib')
sage: R = singular.ring(0, '(x,y)', 'dp')
sage: i2 = singular.ideal('y^9 - x^2*(x-1)^9 + x')
sage: i2.genus()
40
```

Note that the genus can be much smaller than the degree:

```
sage: i = singular.ideal('y^9 - x^2*(x-1)^9')
sage: i.genus()
0
```

## 40.4 An Important Concept

AUTHORS:

- Neal Harris

The following illustrates an important concept: how Sage interacts with the data being used and returned by Singular. Let's compute a Groebner basis for some ideal, using Singular through Sage.

```
sage: singular.lib('poly.lib')
sage: singular.ring(32003, '(a,b,c,d,e,f)', 'lp')
// characteristic : 32003
// number of vars : 6
//      block   1 : ordering lp
//                  : names      a b c d e f
//      block   2 : ordering C
sage: I = singular.ideal('cyclic(6)')
sage: g = singular('groebner(I)')
Traceback (most recent call last):
...
TypeError: Singular error:
...
```

We restart everything and try again, but correctly.

```

sage: singular.quit()
sage: singular.lib('poly.lib'); R = singular.ring(32003, '(a,b,c,d,e,f)', 'lp')
sage: I = singular.ideal('cyclic(6)')
sage: I.groebner()
f^48-2554*f^42-15674*f^36+12326*f^30-12326*f^18+15674*f^12+2554*f^6-1,
...

```

It's important to understand why the first attempt at computing a basis failed. The line where we gave singular the input 'groebner(I)' was useless because Singular has no idea what 'I' is! Although 'I' is an object that we computed with calls to Singular functions, it actually lives in Sage. As a consequence, the name 'I' means nothing to Singular. When we called `I.groebner()`, Sage was able to call the `groebner` function on 'I' in Singular, since 'I' actually means something to Sage.

## 40.5 Long Input

The Singular interface reads in even very long input (using files) in a robust manner, as long as you are creating a new object.

```

sage: t = '%"s"%10^15000 # 15 thousand character string (note that normal Singular input must be a
sage: a = singular.eval(t)
sage: a = singular(t)

```

TESTS:

We test an automatic coercion:

```

sage: a = 3*singular('2'); a
6
sage: type(a)
<class 'sage.interfaces.singular.SingularElement'>
sage: a = singular('2')*3; a
6
sage: type(a)
<class 'sage.interfaces.singular.SingularElement'>

```

Create a ring over GF(9) to check that `gftables` has been installed, see [trac ticket #11645](#):

```

sage: singular.eval("ring testgf9 = (9,x), (a,b,c,d,e,f), (M((1,2,3,0)), wp(2,3), lp);")
''

```

```

class sage.interfaces.singular.Singular(maxread=1000, script_subdirectory=None, log-
                                         file=None, server=None, server_tmpdir=None,
                                         seed=None)

```

Bases: `sage.interfaces.expect.Expect`

Interface to the Singular interpreter.

EXAMPLES: A Groebner basis example.

```

sage: R = singular.ring(0, '(x0,x1,x2)', 'lp')
sage: I = singular.ideal([ 'x0*x1*x2 -x0^2*x2', 'x0^2*x1*x2-x0*x1^2*x2-x0*x1*x2^2', 'x0*x1-x0*x2
sage: I.groebner()
x1^2*x2^2,
x0*x2^3-x1^2*x2^2+x1*x2^3,
x0*x1-x0*x2-x1*x2,
x0^2*x2-x0*x2^2-x1*x2^2

```

AUTHORS:

•David Joyner and William Stein

**LIB** (*lib*, *reload=False*)

Load the Singular library named *lib*.

Note that if the library was already loaded during this session it is not reloaded unless the optional *reload* argument is *True* (the default is *False*).

EXAMPLES:

```
sage: singular.lib('sing.lib')
sage: singular.lib('sing.lib', reload=True)
```

**clear** (*var*)

Clear the variable named *var*.

EXAMPLES:

```
sage: singular.set('int', 'x', '2')
sage: singular.get('x')
'2'
sage: singular.clear('x')
```

“Clearing the variable” means to allow to free the memory that it uses in the Singular sub-process. However, the actual deletion of the variable is only committed when the next element in the Singular interface is created:

```
sage: singular.get('x')
'2'
sage: a = singular(3)
sage: singular.get('x')
'`x`'
```

**console** ()

EXAMPLES:

```
sage: singular_console() #not tested
SINGULAR / Development
A Computer Algebra System for Polynomial Computations / version 3-0-4
0<
by: G.-M. Greuel, G. Pfister, H. Schoenemann \ Nov 2007
FB Mathematik der Universitaet, D-67653 Kaiserslautern
```

**cputime** (*t=None*)

Returns the amount of CPU time that the Singular session has used. If *t* is not *None*, then it returns the difference between the current CPU time and *t*.

EXAMPLES:

```
sage: t = singular.cputime()
sage: R = singular.ring(0, '(x0,x1,x2)', 'lp')
sage: I = singular.ideal([ 'x0*x1*x2 -x0^2*x2', 'x0^2*x1*x2-x0*x1^2*x2-x0*x1*x2^2', 'x0*x1-x0^2*x1^2' ])
sage: gb = I.groebner()
sage: singular.cputime(t) #random
0.02
```

**current\_ring** ()

Returns the current ring of the running Singular session.

EXAMPLES:

```

sage: r = PolynomialRing(GF(127), 3, 'xyz', order='invlex')
sage: r._singular_()
// characteristic : 127
// number of vars : 3
//      block   1 : ordering rp
//              : names   x y z
//      block   2 : ordering C
sage: singular.current_ring()
// characteristic : 127
// number of vars : 3
//      block   1 : ordering rp
//              : names   x y z
//      block   2 : ordering C

```

### **current\_ring\_name()**

Returns the Singular name of the currently active ring in Singular.

OUTPUT: currently active ring's name

EXAMPLES:

```

sage: r = PolynomialRing(GF(127), 3, 'xyz')
sage: r._singular_().name() == singular.current_ring_name()
True

```

### **eval(x, allow\_semicolon=True, strip=True, \*\*kws)**

Send the code x to the Singular interpreter and return the output as a string.

INPUT:

- x - string (of code)
- allow\_semicolon - default: False; if False then raise a TypeError if the input line contains a semicolon.
- strip - ignored

EXAMPLES:

```

sage: singular.eval('2 > 1')
'1'
sage: singular.eval('2 + 2')
'4'

```

if the verbosity level is > 1 comments are also printed and not only returned.

```

sage: r = singular.ring(0, '(x,y,z)', 'dp')
sage: i = singular.ideal(['x^2', 'y^2', 'z^2'])
sage: s = i.std()
sage: singular.eval('hilb(%s)'%(s.name()))
'// 1 t^0\n// -3 t^2\n// 3 t^4\n// -1 t^6\n\n// 1 t^0\n// 3 t^1\n// 3 t^2\n// 1 t^3\n// dimension (affine) = 0\n// degree (affine) = 8'

```

```

sage: set_verbosity(1)
sage: o = singular.eval('hilb(%s)'%(s.name()))
//      1 t^0
//      -3 t^2
//      3 t^4
//      -1 t^6
//      1 t^0
//      3 t^1

```

```
//      3 t^2
//      1 t^3
// dimension (affine) = 0
// degree (affine) = 8
```

This is mainly useful if this method is called implicitly. Because then intermediate results, debugging outputs and printed statements are printed

```
sage: o = s.hilb()
//      1 t^0
//      -3 t^2
//      3 t^4
//      -1 t^6
//      1 t^0
//      3 t^1
//      3 t^2
//      1 t^3
// dimension (affine) = 0
// degree (affine) = 8
// ** right side is not a datum, assignment ignored
```

rather than ignored

```
sage: set_verbose(0)
sage: o = s.hilb()
```

**get** (*var*)

Get string representation of variable named *var*.

EXAMPLES:

```
sage: singular.set('int', 'x', '2')
sage: singular.get('x')
'2'
```

**ideal** (*\*gens*)

Return the ideal generated by *gens*.

INPUT:

- *gens* - list or tuple of Singular objects (or objects that can be made into Singular objects via evaluation)

OUTPUT: the Singular ideal generated by the given list of *gens*

EXAMPLES: A Groebner basis example done in a different way.

```
sage: _ = singular.eval("ring R=0, (x0,x1,x2),lp")
sage: i1 = singular.ideal([ 'x0*x1*x2 -x0^2*x2', 'x0^2*x1*x2-x0*x1^2*x2-x0*x1*x2^2', 'x0*x1-
sage: i1
-x0^2*x2+x0*x1*x2,
x0^2*x1*x2-x0*x1^2*x2-x0*x1*x2^2,
x0*x1-x0*x2-x1*x2

sage: i2 = singular.ideal('groebner(%s);'%i1.name())
sage: i2
x1^2*x2^2,
x0*x2^3-x1^2*x2^2+x1*x2^3,
x0*x1-x0*x2-x1*x2,
x0^2*x2-x0*x2^2-x1*x2^2
```



**lib** (*lib*, *reload=False*)

Load the Singular library named *lib*.

Note that if the library was already loaded during this session it is not reloaded unless the optional *reload* argument is *True* (the default is *False*).

EXAMPLES:

```
sage: singular.lib('sing.lib')
sage: singular.lib('sing.lib', reload=True)
```

**list** (*x*)

Creates a list in Singular from a Sage list *x*.

EXAMPLES:

```
sage: singular.list([1,2])
[1]:
  1
[2]:
  2
```

**load** (*lib*, *reload=False*)

Load the Singular library named *lib*.

Note that if the library was already loaded during this session it is not reloaded unless the optional *reload* argument is *True* (the default is *False*).

EXAMPLES:

```
sage: singular.lib('sing.lib')
sage: singular.lib('sing.lib', reload=True)
```

**matrix** (*nrows*, *ncols*, *entries=None*)

EXAMPLES:

```
sage: singular.lib("matrix")
sage: R = singular.ring(0, '(x,y,z)', 'dp')
sage: A = singular.matrix(3,2,'1,2,3,4,5,6')
sage: A
1, 2,
3, 4,
5, 6
sage: A.gauss_col()
2, -1,
1, 0,
0, 1
```

AUTHORS:

•Martin Albrecht (2006-01-14)

**option** (*cmd=None*, *val=None*)

Access to Singular's options as follows:

Syntax: `option()` Returns a string of all defined options.

Syntax: `option( 'option_name' )` Sets an option. Note to disable an option, use the prefix `no`.

Syntax: `option( 'get' )` Returns an intvec of the state of all options.

Syntax: `option( 'set', intvec_expression )` Restores the state of all options from an intvec (produced by `option('get')`).

EXAMPLES:

```
sage: singular.option()
//options: redefine loadLib usage prompt
sage: singular.option('get')
0,
10321
sage: old_options = _
sage: singular.option('noredefine')
sage: singular.option()
//options: loadLib usage prompt
sage: singular.option('set', old_options)
sage: singular.option('get')
0,
10321
```

**ring** (*char=0, vars='(x)', order='lp', check=True*)  
Create a Singular ring and makes it the current ring.

INPUT:

- *char* - characteristic of the base ring (see examples below), which must be either 0, prime (!), or one of several special codes (see examples below).
- *vars* - a tuple or string that defines the variable names
- *order* - string - the monomial order (default: 'lp')
- *check* - if True, check primality of the characteristic if it is an integer.

OUTPUT: a Singular ring

---

**Note:** This function is *not* identical to calling the Singular `ring` function. In particular, it also attempts to “kill” the variable names, so they can actually be used without getting errors, and it sets printing of elements for this range to short (i.e., with \*’s and carets).

---

EXAMPLES: We first declare  $\mathbb{Q}[x, y, z]$  with degree reverse lexicographic ordering.

```
sage: R = singular.ring(0, '(x,y,z)', 'dp')
sage: R
// characteristic : 0
// number of vars : 3
//      block   1 : ordering dp
//              : names   x y z
//      block   2 : ordering C

sage: R1 = singular.ring(32003, '(x,y,z)', 'dp')
sage: R2 = singular.ring(32003, '(a,b,c,d)', 'lp')
```

This is a ring in variables named  $x(1)$  through  $x(10)$  over the finite field of order 7:

```
sage: R3 = singular.ring(7, '(x(1..10))', 'ds')
```

This is a polynomial ring over the transcendental extension  $\mathbb{Q}(a)$  of  $\mathbb{Q}$ :

```
sage: R4 = singular.ring('(0,a)', '(mu,nu)', 'lp')
```

This is a ring over the field of single-precision floats:

```
sage: R5 = singular.ring('real', '(a,b)', 'lp')
```

This is over 50-digit floats:

```
sage: R6 = singular.ring(' (real,50)', ' (a,b)', 'lp')
sage: R7 = singular.ring(' (complex,50,i)', ' (a,b)', 'lp')
```

To use a ring that you’ve defined, use the `set_ring()` method on the ring. This sets the ring to be the “current ring”. For example,

```
sage: R = singular.ring(7, ' (a,b)', 'ds')
sage: S = singular.ring('real', ' (a,b)', 'lp')
sage: singular.new('10*a')
1.000e+01*a
sage: R.set_ring()
sage: singular.new('10*a')
3*a
```

**set** (*type, name, value*)

Set the variable with given name to the given value.

REMARK:

If a variable in the Singular interface was previously marked for deletion, the actual deletion is done here, before the new variable is created in Singular.

EXAMPLES:

```
sage: singular.set('int', 'x', '2')
sage: singular.get('x')
'2'
```

We test that an unused variable is only actually deleted if this method is called:

```
sage: a = singular(3)
sage: n = a.name()
sage: del a
sage: singular.eval(n)
'3'
sage: singular.set('int', 'y', '5')
sage: singular.eval('defined(%s)'%n)
'0'
```

**set\_ring** (*R*)

Sets the current Singular ring to *R*.

EXAMPLES:

```
sage: R = singular.ring(7, ' (a,b)', 'ds')
sage: S = singular.ring('real', ' (a,b)', 'lp')
sage: singular.current_ring()
// characteristic : 0 (real)
// number of vars : 2
//      block   1 : ordering lp
//                : names   a b
//      block   2 : ordering C
sage: singular.set_ring(R)
sage: singular.current_ring()
// characteristic : 7
// number of vars : 2
//      block   1 : ordering ds
//                : names   a b
//      block   2 : ordering C
```

**set\_seed** (*seed=None*)

Sets the seed for singular interpreter. The seed should be an integer at least 1 and not more than 30 bits. See [http://www.singular.uni-kl.de/Manual/html/sing\\_19.htm#SEC26](http://www.singular.uni-kl.de/Manual/html/sing_19.htm#SEC26) and [http://www.singular.uni-kl.de/Manual/html/sing\\_283.htm#SEC323](http://www.singular.uni-kl.de/Manual/html/sing_283.htm#SEC323)

EXAMPLES:

```
sage: s = Singular()
sage: s.set_seed(1)
1
sage: [s.random(1,10) for i in range(5)]
[8, 10, 4, 9, 1]
```

**setring**(*R*)

Sets the current Singular ring to *R*.

EXAMPLES:

```
sage: R = singular.ring(7, '(a,b)', 'ds')
sage: S = singular.ring('real', '(a,b)', 'lp')
sage: singular.current_ring()
// characteristic : 0 (real)
// number of vars : 2
//      block 1 : ordering lp
//      : names a b
//      block 2 : ordering C
sage: singular.set_ring(R)
sage: singular.current_ring()
// characteristic : 7
// number of vars : 2
//      block 1 : ordering ds
//      : names a b
//      block 2 : ordering C
```

**string**(*x*)

Creates a Singular string from a Sage string. Note that the Sage string has to be “double-quoted”.

EXAMPLES:

```
sage: singular.string(' "Sage" ')
Sage
```

**trait\_names**()

Return a list of all Singular commands.

EXAMPLES:

```
sage: singular.trait_names()
['exteriorPower',
 ...
 'stdfglm']
```

**version**()

EXAMPLES:

**class** sage.interfaces.singular.**SingularElement** (*parent, type, value, is\_name=False*)  
Bases: sage.interfaces.expect.ExpectElement

EXAMPLES:

```
sage: a = singular(2)
sage: loads(dumps(a))
2
```

**attrib** (*name*, *value=None*)

Get and set attributes for self.

INPUT:

- name - string to choose the attribute
- value - boolean value or None for reading, (default:None)

VALUES: isSB - the standard basis property is set by all commands computing a standard basis like groebner, std, stdhilb etc.; used by lift, dim, degree, mult, hilb, vdim, kbase isHomog - the weight vector for homogeneous or quasihomogeneous ideals/modules isCI - complete intersection property isCM - Cohen-Macaulay property rank - set the rank of a module (see nrows) withSB - value of type ideal, resp. module, is std withHilb - value of type intvec is hilb(,1) (see hilb) withRes - value of type list is a free resolution withDim - value of type int is the dimension (see dim) withMult - value of type int is the multiplicity (see mult)

EXAMPLE:

```
sage: P.<x,y,z> = PolynomialRing(QQ)
sage: I = Ideal([z^2, y*z, y^2, x*z, x*y, x^2])
sage: Ibar = I._singular_()
sage: Ibar.attrib('isSB')
0
sage: singular.eval('vdim(%s)'%Ibar.name()) # sage7 name is random
// ** sage7 is no standard basis
4
sage: Ibar.attrib('isSB',1)
sage: singular.eval('vdim(%s)'%Ibar.name())
'4'
```

**is\_string** ()

Tell whether this element is a string.

EXAMPLES:

```
sage: singular('"abc"').is_string()
True
sage: singular('1').is_string()
False
```

**sage\_flattened\_str\_list** ()

EXAMPLES:

```
sage: R=singular.ring(0,'(x,y)', 'dp')
sage: RL = R.ringlist()
sage: RL.sage_flattened_str_list()
['0', 'x', 'y', 'dp', '1,1', 'C', '0', '_[1]=0']
```

**sage\_global\_ring** ()

Return the current basering in Singular as a polynomial ring or quotient ring.

EXAMPLE:

```
sage: singular.eval('ring r1 = (9,x), (a,b,c,d,e,f), (M((1,2,3,0)),wp(2,3),lp)')
,,
sage: R = singular('r1').sage_global_ring()
sage: R
Multivariate Polynomial Ring in a, b, c, d, e, f over Finite Field in x of size 3^2
sage: R.term_order()
Block term order with blocks:
(Matrix term order with matrix
```

```

[1 2]
[3 0],
  Weighted degree reverse lexicographic term order with weights (2, 3),
  Lexicographic term order of length 2)

sage: singular.eval('ring r2 = (0,x), (a,b,c), dp')
''
sage: singular('r2').sage_global_ring()
Multivariate Polynomial Ring in a, b, c over Fraction Field of Univariate Polynomial Ring in
sage: singular.eval('ring r3 = (3,z), (a,b,c), dp')
''
sage: singular.eval('minpoly = 1+z+z2+z3+z4')
''
sage: singular('r3').sage_global_ring()
Multivariate Polynomial Ring in a, b, c over Finite Field in z of size 3^4

```

Real and complex fields in both Singular and Sage are defined with a precision. The precision in Singular is given in terms of digits, but in Sage it is given in terms of bits. So, the digit precision is internally converted to a reasonable bit precision:

```

sage: singular.eval('ring r4 = (real,20), (a,b,c), dp')
''
sage: singular('r4').sage_global_ring()
Multivariate Polynomial Ring in a, b, c over Real Field with 70 bits of precision

```

The case of complex coefficients is not fully supported, yet, since the generator of a complex field in Sage is always called “I”:

```

sage: singular.eval('ring r5 = (complex,15,j), (a,b,c), dp')
''
sage: R = singular('r5').sage_global_ring(); R
Multivariate Polynomial Ring in a, b, c over Complex Field with 54 bits of precision
sage: R.base_ring() ('j')
Traceback (most recent call last):
...
NameError: name 'j' is not defined
sage: R.base_ring() ('I')
1.000000000000000*I

```

In our last example, the base ring is a quotient ring:

```

sage: singular.eval('ring r6 = (9,a), (x,y,z), lp')
''
sage: Q = singular('std(ideal(x^2,x+y^2+z^3))', type='qring')
sage: Q.sage_global_ring()
Quotient of Multivariate Polynomial Ring in x, y, z over Finite Field in a of size 3^2 by th

```

AUTHOR:

- Simon King (2011-06-06)

**sage\_matrix** (*R*, *sparse=True*)  
Returns Sage matrix for self

INPUT:

- R* - (default: None); an optional ring, over which the resulting matrix is going to be defined. By default, the output of `sage_global_ring()` is used.

- `sparse` - (default: `True`); determines whether the resulting matrix is sparse or not.

## EXAMPLES:

```
sage: R = singular.ring(0, '(x,y,z)', 'dp')
sage: A = singular.matrix(2,2)
sage: A.sage_matrix(ZZ)
[0 0]
[0 0]
sage: A.sage_matrix(RDF)
[0.0 0.0]
[0.0 0.0]
```

**sage\_poly** (*R=None, kcache=None*)

Returns a Sage polynomial in the ring *r* matching the provided poly which is a singular polynomial.

## INPUT:

- R* - (default: `None`); an optional polynomial ring. If it is provided, then you have to make sure that it matches the current singular ring as, e.g., returned by `singular.current_ring()`. By default, the output of `sage_global_ring()` is used.
- kcache* - (default: `None`); an optional dictionary for faster finite field lookups, this is mainly useful for finite extension fields

## OUTPUT: MPolynomial

## EXAMPLES:

```
sage: R = PolynomialRing(GF(2^8, 'a'), 2, 'xy')
sage: f=R('a^20*x^2*y+a^10+x')
sage: f._singular_().sage_poly(R)==f
True
sage: R = PolynomialRing(GF(2^8, 'a'), 1, 'x')
sage: f=R('a^20*x^3+x^2+a^10')
sage: f._singular_().sage_poly(R)==f
True

sage: P.<x,y> = PolynomialRing(QQ, 2)
sage: f = x*y**3 - 1/9 * x + 1; f
x*y^3 - 1/9*x + 1
sage: singular(f)
x*y^3-1/9*x+1
sage: P(singular(f))
x*y^3 - 1/9*x + 1
```

## TESTS:

```
sage: singular.eval('ring r = (3,z), (a,b,c), dp')
''
sage: singular.eval('minpoly = 1+z+z2+z3+z4')
''
sage: p = singular('z^4*a^3+z^2*a*b*c')
sage: p.sage_poly()
(-z^3 - z^2 - z - 1)*a^3 + (z^2)*a*b*c
sage: singular('z^4')
(-z3-z2-z-1)
```

## AUTHORS:

- Martin Albrecht (2006-05-18)

- Simon King (2011-06-06): Deal with Singular's short polynomial representation, automatic construction of a polynomial ring, if it is not explicitly given.

---

**Note:** For very simple polynomials `eval(SingularElement.sage_polystring())` is faster than `SingularElement.sage_poly(R)`, maybe we should detect the crossover point (in dependence of the string length) and choose an appropriate conversion strategy

---

#### **sage\_polystring()**

If this Singular element is a polynomial, return a string representation of this polynomial that is suitable for evaluation in Python. Thus `*` is used for multiplication and `**` for exponentiation. This function is primarily used internally.

The `short=0` option *must* be set for the parent ring or this function will not work as expected. This option is set by default for rings created using `singular.ring` or set using `ring_name.set_ring()`.

EXAMPLES:

```
sage: R = singular.ring(0, '(x,y)')
sage: f = singular('x^3 + 3*y^11 + 5')
sage: f
x^3+3*y^11+5
sage: f.sage_polystring()
'x**3+3*y**11+5'
```

#### **sage\_structured\_str\_list()**

If self is a Singular list of lists of Singular elements, returns corresponding Sage list of lists of strings.

EXAMPLES:

```
sage: R=singular.ring(0, '(x,y)', 'dp')
sage: RL=R.ringlist()
sage: RL
[1]:
  0
[2]:
  [1]:
    x
  [2]:
    y
[3]:
  [1]:
    [1]:
      dp
    [2]:
      1,1
  [2]:
    [1]:
      C
    [2]:
      0
[4]:
  _[1]=0
sage: RL.sage_structured_str_list()
['0', ['x', 'y'], [['dp', '1,\n1 '], ['C', '0 ']], '0']
```

#### **set\_ring()**

Sets the current ring in Singular to be self.

EXAMPLES:



```

sage: R = singular.ring(7, '(a,b)', 'ds')
sage: S = singular.ring('real', '(a,b)', 'lp')
sage: singular.current_ring()
// characteristic : 0 (real)
// number of vars : 2
//      block 1 : ordering lp
//      : names a b
//      block 2 : ordering C
sage: R.set_ring()
sage: singular.current_ring()
// characteristic : 7
// number of vars : 2
//      block 1 : ordering ds
//      : names a b
//      block 2 : ordering C

```

**trait\_names()**

Returns the possible tab-completions for self. In this case, we just return all the tab completions for the Singular object.

EXAMPLES:

```

sage: R = singular.ring(0, '(x,y)', 'dp')
sage: R.trait_names()
['exteriorPower',
 ...
 'stdfglm']

```

**type()**

Returns the internal type of this element.

EXAMPLES:

```

sage: R = PolynomialRing(GF(2^8, 'a'), 2, 'x')
sage: R._singular_.type()
'ring'
sage: fs = singular('x0^2', 'poly')
sage: fs.type()
'poly'

```

**exception sage.interfaces.singular.SingularError**

Bases: `exceptions.RuntimeError`

Raised if Singular printed an error message

**class sage.interfaces.singular.SingularFunction** (*parent, name*)

Bases: `sage.interfaces.expect.ExpectFunction`

**class sage.interfaces.singular.SingularFunctionElement** (*obj, name*)

Bases: `sage.interfaces.expect.FunctionElement`

**class sage.interfaces.singular.SingularGBDefaultContext** (*singular=None*)

Within this context all Singular Groebner basis calculations are reduced automatically.

AUTHORS:

- Martin Albrecht
- Simon King

**class** `sage.interfaces.singular.SingularGBLogPrettyPrinter` (*verbosity=1*)

A device which prints Singular Groebner basis computation logs more verbatim.

**flush()**

EXAMPLE:

```
sage: from sage.interfaces.singular import SingularGBLogPrettyPrinter
sage: s3 = SingularGBLogPrettyPrinter(verbosity=3)
sage: s3.flush()
```

**write(s)**

EXAMPLE:

```
sage: from sage.interfaces.singular import SingularGBLogPrettyPrinter
sage: s3 = SingularGBLogPrettyPrinter(verbosity=3)
sage: s3.write("(S:1337)")
Performing complete reduction of 1337 elements.
sage: s3.write("M[389,12]")
Parallel reduction of 389 elements with 12 non-zero output elements.
```

`sage.interfaces.singular.generate_docstring_dictionary()`

Generate global dictionaries which hold the docstrings for Singular functions.

EXAMPLE:

```
sage: from sage.interfaces.singular import generate_docstring_dictionary
sage: generate_docstring_dictionary()
```

`sage.interfaces.singular.get_docstring(name)`

Return the docstring for the function name.

INPUT:

- name - a Singular function name

EXAMPLE:

```
sage: from sage.interfaces.singular import get_docstring
sage: 'groebner' in get_docstring('groebner')
True
sage: 'standard.lib' in get_docstring('groebner')
True
```

`sage.interfaces.singular.is_SingularElement(x)`

Returns True if x is of type SingularElement.

EXAMPLES:

```
sage: from sage.interfaces.singular import is_SingularElement
sage: is_SingularElement(singular(2))
True
sage: is_SingularElement(2)
False
```

`sage.interfaces.singular.reduce_load()`

This is for backwards compatibility only.

To be precise, it only serves at unpickling the invalid singular elements that are stored in the pickle jar.

EXAMPLES:

```
sage: from sage.interfaces.singular import reduce_load
sage: reduce_load()
```

```
doctest:...: DeprecationWarning: This function is only used to unpickle invalid objects
See http://trac.sagemath.org/18848 for details.
(invalid object -- defined in terms of closed session)
```

By [trac ticket #18848](#), pickling actually often works:

```
sage: loads(dumps(singular.ring()))
// characteristic : 0
// number of vars : 1
//      block    1 : ordering lp
//              : names      x
//      block    2 : ordering C
```

`sage.interfaces.singular.reduce_load_Singular()`

EXAMPLES:

```
sage: from sage.interfaces.singular import reduce_load_Singular
sage: reduce_load_Singular()
Singular
```

`sage.interfaces.singular.singular_console()`

Spawn a new Singular command-line session.

EXAMPLES:

```
sage: singular_console() #not tested
SINGULAR / Development
A Computer Algebra System for Polynomial Computations / version 3-0-4
0<
by: G.-M. Greuel, G. Pfister, H. Schoenemann \ Nov 2007
FB Mathematik der Universitaet, D-67653 Kaiserslautern
```

`sage.interfaces.singular.singular_gb_standard_options(func)`

Decorator to force a reduced Singular groebner basis.

TESTS:

```
sage: P.<a,b,c,d,e> = PolynomialRing(GF(127))
sage: J = sage.rings.ideal.Cyclic(P).homogenize()
sage: from sage.misc.sageinspect import sage_getsource
sage: "basis" in sage_getsource(J.interreduced_basis) #indirect doctest
True
```

The following tests against a bug that was fixed in [trac ticket #11298](#):

```
sage: from sage.misc.sageinspect import sage_getsourcelines, sage_getargspec
sage: P.<x,y> = QQ[]
sage: I = P*[x,y]
sage: sage_getargspec(I.interreduced_basis)
ArgSpec(args=['self'], varargs=None, keywords=None, defaults=None)
sage: sage_getsourcelines(I.interreduced_basis)
([' @singular_gb_standard_options\n',
  ' @libsingular_gb_standard_options\n',
  ' def interreduced_basis(self):\n', '
...
  ' return self.basis.reduced()\n'], ...)
```

---

**Note:** This decorator is used automatically internally so the user does not need to use it manually.

---

`sage.interfaces.singular.singular_version()`

Returns the version of Singular being used.

EXAMPLES:

## THE TACHYON RAY TRACER

### AUTHOR:

- John E. Stone

**class** `sage.interfaces.tachyon.TachyonRT`

Bases: `sage.structure.sage_object.SageObject`

The Tachyon Ray Tracer

`tachyon_rt(model, outfile='sage.png', verbose=1, block=True, extra_opts='')`

### INPUT:

- `model` - a string that describes a 3d model in the Tachyon modeling format. Type `tachyon_rt.help()` for a description of this format.
- `outfile` - (default: 'sage.png') output filename; the extension of the filename determines the type. Supported types include:
  - `tga` - 24-bit (uncompressed)
  - `bmp` - 24-bit Windows BMP (uncompressed)
  - `ppm` - 24-bit PPM (uncompressed)
  - `rgb` - 24-bit SGI RGB (uncompressed)
  - `png` - 24-bit PNG (compressed, lossless)
- `verbose` - integer; (default: 1)
  - 0 - silent
  - 1 - some output
  - 2 - very verbose output
- `block` - bool (default: True); if False, run the rendering command in the background.
- `extra_opts` - passed directly to tachyon command line. Use `tachyon_rt.usage()` to see some of the possibilities.

### OUTPUT:

- Some text may be displayed onscreen.
- The file `outfile` is created.

### EXAMPLES:

**\_\_call\_\_** (*model*, *outfile*='sage.png', *verbose*=1, *extra\_opts*='')

This executes the tachyon program, given a scene file input.

INPUT:

- *model* – string. The tachyon model.
- *outfile* – string, default 'sage.png'. The filename to save the model to.
- *verbose* – 0, 1, (default) or 2. The verbosity level.
- *extra\_opts* – string (default: empty string). Extra options that will be appended to the tachyon commandline.

EXAMPLES:

```
sage: from sage.interfaces.tachyon import TachyonRT
sage: tgen = Tachyon()
sage: tgen.texture('t1')
sage: tgen.sphere((0,0,0),1,'t1')
sage: tgen.str()[30:40]
'resolution'
sage: t = TachyonRT()
sage: import os
sage: t(tgen.str(), outfile=os.devnull)
tachyon ...
Tachyon Parallel/Multiprocessor Ray Tracer...
```

TESTS:

```
sage: from sage.env import SAGE_EXTCODE
sage: filename = os.path.join(SAGE_EXTCODE, 'doctest', 'invalid', 'syntax_error.tachyon')
sage: syntax_error = open(filename, 'r').read()
sage: t(syntax_error, outfile=os.devnull)
Traceback (most recent call last):
...
RuntimeError: Tachyon Parallel/Multiprocessor Ray Tracer...
...
Parser failed due to an input file syntax error.
Aborting render.
```

**help** (*use\_pager*=True)

Prints (pages) the help file written by John Stone describing scene files for Tachyon. The output is paged unless *use\_pager*=False.

TESTS:

```
sage: from sage.interfaces.tachyon import TachyonRT
sage: t = TachyonRT()
sage: t.help(use_pager=False)
This help, which was written by John Stone, describes ...
```

**usage** (*use\_pager*=True)

Returns the basic description of using the Tachyon raytracer (simply what is returned by running tachyon with no input). The output is paged unless *use\_pager*=False.

TESTS:

```
sage: from sage.interfaces.tachyon import TachyonRT
sage: t = TachyonRT()
sage: t.usage(use_pager=False)
Tachyon Parallel/Multiprocessor Ray Tracer    Version...
```

## INTERFACE TO TIDES

This module contains tools to write the .c files needed for TIDES [TI] .

Tides is an integration engine based on the Taylor method. It is implemented as a c library. The user must translate its initial value problem (IVP) into a pair of .c files that will then be compiled and linked against the TIDES library. The resulting binary will produce the desired output. The tools in this module can be used to automate the generation of these files from the symbolic expression of the differential equation.

```
#####  
# Copyright (C) 2014 Miguel Marco <mmarco@unizar.es>, Marcos Rodriguez  
# <marcos@unizar.es>  
#  
# Distributed under the terms of the GNU General Public License (GPL):  
#  
# http://www.gnu.org/licenses/  
#####
```

### AUTHORS:

- Miguel Marco (06-2014) - Implementation of tides solver
- Marcos Rodriguez (06-2014) - Implementation of tides solver
- Alberto Abad (06-2014) - tides solver
- Roberto Barrio (06-2014) - tides solver

### REFERENCES:

`sage.interfaces.tides.genfiles_mintides`(*integrator, driver, f, ics, initial, final, delta, tolrel=1e-16, tolabs=1e-16, output=''*)

Generate the needed files for the min\_tides library.

### INPUT:

- `integrator` – the name of the integrator file.
- `driver` – the name of the driver file.
- `f` – the function that determines the differential equation.
- `ics` – a list or tuple with the initial conditions.
- `initial` – the initial time for the integration.
- `final` – the final time for the integration.
- `delta` – the step of the output.
- `tolrel` – the relative tolerance.
- `tolabs` – the absolute tolerance.

- output – the name of the file that the compiled integrator will write to

This function creates two files, integrator and driver, that can be used later with the min\_tides library [TI].

TESTS:

```
sage: from sage.interfaces.tides import genfiles_mintides
sage: import os
sage: import shutil
sage: from sage.misc.temporary_file import tmp_dir
sage: tempdir = tmp_dir()
sage: intfile = os.path.join(tempdir, 'integrator.c')
sage: drfile = os.path.join(tempdir, 'driver.c')
sage: var('t,x,y,X,Y')
      (t, x, y, X, Y)
sage: f(t,x,y,X,Y)=[X, Y, -x/(x^2+y^2)^(3/2), -y/(x^2+y^2)^(3/2)]
sage: genfiles_mintides(intfile, drfile, f, [1,0, 0, 0.2], 0, 10, 0.1, output = 'out')
sage: fileint = open(intfile)
sage: l = fileint.readlines()
sage: fileint.close()
sage: l[5]
'    #include "minc_tides.h"\n'
sage: l[15]
'    double XX[TT+1][MO+1];\n'
sage: l[25]
'\n'
sage: l[35]
'\t\tXX[1][i+1] = XX[3][i] / (i+1.0);\n'
sage: filedr = open(drfile)
sage: l = filedr.readlines()
sage: filedr.close()
sage: l[6]
'    #include "minc_tides.h"\n'
sage: l[15]
'    double tolrel, tolabs, tini, tend, dt;\n'
sage: l[25]
'\ttolrel = 9.9999999999999998e-17 ;\n'
sage: shutil.rmtree(tempdir)
```

Check that ticket [trac ticket #17179](#) is fixed (handle expressions like  $\pi$ ):

```
sage: from sage.interfaces.tides import genfiles_mintides
sage: import os
sage: import shutil
sage: from sage.misc.temporary_file import tmp_dir
sage: tempdir = tmp_dir()
sage: intfile = os.path.join(tempdir, 'integrator.c')
sage: drfile = os.path.join(tempdir, 'driver.c')
sage: var('t,x,y,X,Y')
      (t, x, y, X, Y)
sage: f(t,x,y,X,Y)=[X, Y, -x/(x^2+y^2)^(3/2), -y/(x^2+y^2)^(3/2)]
sage: genfiles_mintides(intfile, drfile, f, [pi, 0, 0, 0.2], 0, 10, 0.1, output = 'out')
sage: fileint = open(intfile)
sage: l = fileint.readlines()
sage: fileint.close()
sage: l[30]
'\t\tXX[8][i] = pow_mc_c(XX[7],-1.5000000000000000,XX[8], i);\n'
sage: filedr = open(drfile)
sage: l = filedr.readlines()
```



```

sage: filedr.close()
sage: l[18]
'      \tv[0] = 3.1415926535897931 ; \n'
sage: shutil.rmtree(tempdir)

```

```

sage.interfaces.tides.genfiles_mpfr(integrator, driver, f, ics, initial, final, delta, parameters=None, parameter_values=None, dig=20, tolrel=1e-16, tolabs=1e-16, output='')

```

Generate the needed files for the mpfr module of the tides library.

INPUT:

- **integrator** – the name of the integrator file.
- **driver** – the name of the driver file.
- **f** – the function that determines the differential equation.
- **ics** – a list or tuple with the initial conditions.
- **initial** – the initial time for the integration.
- **final** – the final time for the integration.
- **delta** – the step of the output.
- **parameters** – the variables inside the function that should be treated as parameters.
- **parameter\_values** – the values of the parameters for the particular initial value problem.
- **dig** – the number of digits of precision that will be used in the integration
- **tolrel** – the relative tolerance.
- **tolabs** – the absolute tolerance.
- **output** – the name of the file that the compiled integrator will write to

This function creates two files, integrator and driver, that can be used later with the tides library ([TI]).

TESTS:

```

sage: from tempfile import mkdtemp
sage: from sage.interfaces.tides import genfiles_mpfr
sage: import os
sage: import shutil
sage: from sage.misc.temporary_file import tmp_dir
sage: tempdir = tmp_dir()
sage: intfile = os.path.join(tempdir, 'integrator.c')
sage: drfile = os.path.join(tempdir, 'driver.c')
sage: var('t,x,y,X,Y')
(t, x, y, X, Y)
sage: f(t,x,y,X,Y)=[X, Y, -x/(x^2+y^2)^(3/2), -y/(x^2+y^2)^(3/2)]
sage: genfiles_mpfr(intfile, drfile, f, [1,0, 0, 0.2], 0, 10, 0.1, output = 'out', dig = 50)
sage: fileint = open(intfile)
sage: l = fileint.readlines()
sage: fileint.close()
sage: l[5]
'      #include "mp_tides.h"\n'
sage: l[15]
'\tstatic int PARAMETERS = 0;\n'
sage: l[25]
'\t\ttmpfrts_var_t(itd, link[5], var[3], i);\n'
sage: l[30]

```

[illegible]

Check that ticket `trac ticket #17179` is fixed (handle expressions like `pi`):

[illegible]

```
sage.interfaces.tides.remove_constants(l1,l2)
```

Given two lists, remove the entries in the first that are real constants, and also the corresponding elements in the second one.

```
sage: from sage.interfaces.tides import subexpressions_list, remove_constants sage:
f(a)=[1+cos(7)*a] sage: l1, l2 = subexpressions_list(f) sage: l1, l2 ([sin(7), cos(7), a*cos(7), a*cos(7)
+ 1], [(‘sin’, 7), (‘cos’, 7), (‘mul’, cos(7), a), (‘add’, 1, a*cos(7))]) sage: remove_constants(l1,l2)
sage: l1, l2 ([a*cos(7), a*cos(7) + 1], [(‘mul’, cos(7), a), (‘add’, 1, a*cos(7))])
```

`sage.interfaces.tides.remove_repeated(l1, l2)`

Given two lists, remove the repeated elements in l1, and the elements in l2 that are on the same position.

EXAMPLES:

```
sage: from sage.interfaces.tides import (subexpressions_list, remove_repeated)
sage: f(a)=[1 + a^2, arcsin(a)]
sage: l1, l2 = subexpressions_list(f)
sage: l1, l2
([a^2, a^2 + 1, a^2, -a^2, -a^2 + 1, sqrt(-a^2 + 1), arcsin(a)],
 [('mul', a, a),
 ('add', 1, a^2),
 ('mul', a, a),
 ('mul', -1, a^2),
 ('add', 1, -a^2),
 ('pow', -a^2 + 1, 0.5),
 ('asin', a)])
sage: remove_repeated(l1, l2)
sage: l1, l2
([a^2, a^2 + 1, -a^2, -a^2 + 1, sqrt(-a^2 + 1), arcsin(a)],
 [('mul', a, a),
 ('add', 1, a^2),
 ('mul', -1, a^2),
 ('add', 1, -a^2),
 ('pow', -a^2 + 1, 0.5),
 ('asin', a)])
```

`sage.interfaces.tides.subexpressions_list(f, pars=None)`

Construct the lists with the intermediate steps on the evaluation of the function.

INPUT:

- `f` – a symbolic function of several components.
- `pars` – a list of the parameters that appear in the function this should be the symbolic constants that appear in `f` but are not arguments.

OUTPUT:

- a list of the intermediate subexpressions that appear in the evaluation of `f`.
- a list with the operations used to construct each of the subexpressions. each element of this list is a tuple, formed by a string describing the operation made, and the operands.

For the trigonometric functions, some extra expressions will be added. These extra expressions will be used later to compute their derivatives.

EXAMPLES:

```
sage: from sage.interfaces.tides import subexpressions_list
sage: var('x,y')
(x, y)
sage: f(x,y) = [x^2+y, cos(x)/log(y)]
sage: subexpressions_list(f)
([x^2, x^2 + y, sin(x), cos(x), log(y), cos(x)/log(y)],
 [('mul', x, x),
 ('add', y, x^2),
 ('sin', x),
 ('cos', x),
 ('log', y),
 ('div', log(y), cos(x))])
```

```
sage: f(a)=[cos(a), arctan(a)]
sage: from sage.interfaces.tides import subexpressions_list
sage: subexpressions_list(f)
([sin(a), cos(a), a^2, a^2 + 1, arctan(a)],
 [('sin', a), ('cos', a), ('mul', a, a), ('add', 1, a^2), ('atan', a)])

sage: from sage.interfaces.tides import subexpressions_list
sage: var('s,b,r')
(s, b, r)
sage: f(t,x,y,z)= [s*(y-x),x*(r-z)-y,x*y-b*z]
sage: subexpressions_list(f,[s,b,r])
([-y,
 x - y,
 s*(x - y),
 -s*(x - y),
 -z,
 r - z,
 (r - z)*x,
 -y,
 (r - z)*x - y,
 x*y,
 b*z,
 -b*z,
 x*y - b*z],
 [('mul', -1, y),
 ('add', -y, x),
 ('mul', x - y, s),
 ('mul', -1, s*(x - y)),
 ('mul', -1, z),
 ('add', -z, r),
 ('mul', x, r - z),
 ('mul', -1, y),
 ('add', -y, (r - z)*x),
 ('mul', y, x),
 ('mul', z, b),
 ('mul', -1, b*z),
 ('add', -b*z, x*y)])

sage: var('x, y')
(x, y)
sage: f(x,y)=[exp(x^2+sin(y))]
sage: from sage.interfaces.tides import *
sage: subexpressions_list(f)
([x^2, sin(y), cos(y), x^2 + sin(y), e^(x^2 + sin(y))],
 [('mul', x, x),
 ('sin', y),
 ('cos', y),
 ('add', sin(y), x^2),
 ('exp', x^2 + sin(y))])
```

## INTERFACE TO THE SAGE CLEANER

Triva Note: For the name “sage-cleaner”, think of the “The Cleaner” from Pulp Fiction:  
<http://www.frankjankowski.de/quiz/illus/keitel.jpg>

```
sage.interfaces.cleaner.cleaner(pid, cmd='')
```

Write a line to the `spawned_processes` file with the given `pid` and `cmd`.

```
sage.interfaces.cleaner.start_cleaner()
```

Start `sage-cleaner` in a new process group.



## QUITTING INTERFACES

```
sage.interfaces.quit.expect_quitall(verbose=False)
```

EXAMPLES:

```
sage: sage.interfaces.quit.expect_quitall()
```

```
sage: gp.eval('a=10')
```

```
'10'
```

```
sage: gp('a')
```

```
10
```

```
sage: sage.interfaces.quit.expect_quitall()
```

```
sage: gp('a')
```

```
a
```

```
sage: sage.interfaces.quit.expect_quitall(verbose=True)
```

```
Exiting PARI/GP interpreter with PID ... running .../local/bin/gp --fast --emacs --quiet --stack
```

```
sage.interfaces.quit.invalidate_all()
```

Invalidate all of the expect interfaces.

This is used, e.g., by the fork-based @parallel decorator.

EXAMPLES:

```
sage: a = maxima(2); b = gp(3)
```

```
sage: a, b
```

```
(2, 3)
```

```
sage: sage.interfaces.quit.invalidate_all()
```

```
sage: a
```

```
<repr(<sage.interfaces.maxima.MaximaElement at 0x...>) failed: ValueError: The maxima session in
```

```
sage: b
```

```
<repr(<sage.interfaces.gp.GpElement at 0x...>) failed: ValueError: The pari session in which thi
```

However the maxima and gp sessions should still work out, though with their state reset:

```
sage: a = maxima(2); b = gp(3) sage: a, b (2, 3)
```

```
sage.interfaces.quit.is_running(pid)
```

Return True if and only if there is a process with id pid running.

```
sage.interfaces.quit.kill_spawned_jobs(verbose=False)
```

INPUT:

- verbose – bool (default: False); if True, display a message each time a process is sent a kill signal

EXAMPLES:

```
sage: gp.eval('a=10')
```

```
'10'
```

```
sage: sage.interfaces.quit.kill_spawned_jobs(verbose=False)
```

```
sage: sage.interfaces.quit.expect_quitall()
```

```
sage: gp.eval('a=10')
'10'
sage: sage.interfaces.quit.kill_spawned_jobs(verbose=True)
Killing spawned job ...
```

After doing the above, we do the following to avoid confusion in other doctests:

```
sage: sage.interfaces.quit.expect_quitall()
```



## AN INTERFACE TO READ DATA FILES

`sage.interfaces.read_data.read_data(f, t)`

Read data from file 'f' and class 't' (one element per line), and returns a list of elements.

INPUT:

- 'f' - a file name
- 't' - a class (objects will be coerced to that class)

OUTPUT:

a list of elements of class 't'.

EXAMPLES:

```
sage: indata = tmp_filename()
sage: f = open(indata, "w")
sage: f.write("17\n42\n")
sage: f.close()
sage: l = read_data(indata, ZZ); l
[17, 42]
sage: f = open(indata, "w")
sage: f.write("1.234\n5.678\n")
sage: f.close()
sage: l = read_data(indata, RealField(17)); l
[1.234, 5.678]
```



## INDICES AND TABLES

- Index
- Module Index
- Search Page



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## i

- sage.interfaces.axiom, 15
- sage.interfaces.cleaner, 303
- sage.interfaces.ecm, 21
- sage.interfaces.expect, 9
- sage.interfaces.four\_ti\_2, 27
- sage.interfaces.fricas, 33
- sage.interfaces.frobby, 37
- sage.interfaces.gap, 41
- sage.interfaces.gap3, 53
- sage.interfaces.gfan, 63
- sage.interfaces.giac, 65
- sage.interfaces.gnuplot, 73
- sage.interfaces.gp, 75
- sage.interfaces.interface, 3
- sage.interfaces.jmoldata, 85
- sage.interfaces.kash, 87
- sage.interfaces.lie, 95
- sage.interfaces.lisp, 103
- sage.interfaces.macaulay2, 107
- sage.interfaces.magma, 117
- sage.interfaces.magma\_free, 135
- sage.interfaces.maple, 137
- sage.interfaces.mathematica, 145
- sage.interfaces.matlab, 153
- sage.interfaces.maxima, 157
- sage.interfaces.maxima\_abstract, 167
- sage.interfaces.maxima\_lib, 183
- sage.interfaces.mupad, 197
- sage.interfaces.mwrank, 201
- sage.interfaces.octave, 205
- sage.interfaces.phc, 211
- sage.interfaces.povray, 217
- sage.interfaces.psage, 219
- sage.interfaces.qepcad, 221
- sage.interfaces.qsieve, 247
- sage.interfaces.quit, 305

`sage.interfaces.r`, [251](#)  
`sage.interfaces.read_data`, [307](#)  
`sage.interfaces.rubik`, [263](#)  
`sage.interfaces.sage0`, [265](#)  
`sage.interfaces.sagespawn`, [13](#)  
`sage.interfaces.scilab`, [269](#)  
`sage.interfaces.singular`, [275](#)  
`sage.interfaces.tachyon`, [295](#)  
`sage.interfaces.tides`, [297](#)



## Symbols

`__call__()` (sage.interfaces.tachyon.TachyonRT method), 295

## A

`A()` (sage.interfaces.qepcad.qepcad\_formula\_factory method), 239  
`alexander_dual()` (sage.interfaces.frobby.Frobby method), 37  
`all_but_finitely_many()` (sage.interfaces.qepcad.qepcad\_formula\_factory method), 241  
`and_()` (sage.interfaces.qepcad.qepcad\_formula\_factory method), 241  
`answer()` (sage.interfaces.qepcad.Qepcad method), 229  
`arguments()` (sage.interfaces.maxima\_abstract.MaximaAbstractElementFunction method), 181  
`as_type()` (sage.interfaces.axiom.PanAxiomElement method), 18  
`AsciiArtString` (class in sage.interfaces.interface), 3  
`assign_names()` (sage.interfaces.magma.MagmaElement method), 129  
`AssignNames()` (sage.interfaces.magma.MagmaElement method), 128  
`associated_primes()` (sage.interfaces.frobby.Frobby method), 38  
`assume()` (sage.interfaces.qepcad.Qepcad method), 230  
`atomic()` (sage.interfaces.qepcad.qepcad\_formula\_factory method), 241  
`Attach()` (sage.interfaces.magma.Magma method), 120  
`attach()` (sage.interfaces.magma.Magma method), 121  
`attach_spec()` (sage.interfaces.magma.Magma method), 122  
`AttachSpec()` (sage.interfaces.magma.Magma method), 121  
`attrib()` (sage.interfaces.singular.SingularElement method), 286  
`attribute()` (sage.interfaces.interface.InterfaceElement method), 5  
`available_packages()` (sage.interfaces.r.R method), 254  
`Axiom` (class in sage.interfaces.axiom), 17  
`axiom_console()` (in module sage.interfaces.axiom), 20  
`AxiomElement` (class in sage.interfaces.axiom), 17  
`AxiomExpectFunction` (class in sage.interfaces.axiom), 17  
`AxiomFunctionElement` (class in sage.interfaces.axiom), 17

## B

`bar_call()` (sage.interfaces.magma.Magma method), 122  
`blackbox()` (sage.interfaces.phc.PHC method), 211  
`bool()` (sage.interfaces.gap.GapElement\_generic method), 45  
`bool()` (sage.interfaces.gp.GpElement method), 82  
`bool()` (sage.interfaces.interface.InterfaceElement method), 5  
`bool()` (sage.interfaces.lisp.LispElement method), 105

`bool()` (`sage.interfaces.maxima_abstract.MaximaAbstractElement` method), 175

## C

`C()` (`sage.interfaces.qepcad.qepcad_formula_factory` method), 239

`call()` (`sage.interfaces.four_ti_2.FourTi2` method), 27

`call()` (`sage.interfaces.interface.Interface` method), 3

`call()` (`sage.interfaces.r.R` method), 255

`cell()` (`sage.interfaces.qepcad.Qepcad` method), 230

`chdir()` (`sage.interfaces.magma.Magma` method), 123

`chdir()` (`sage.interfaces.mathematica.Mathematica` method), 150

`chdir()` (`sage.interfaces.matlab.Matlab` method), 155

`chdir()` (`sage.interfaces.maxima_abstract.MaximaAbstract` method), 167

`chdir()` (`sage.interfaces.r.R` method), 255

`circuits()` (`sage.interfaces.four_ti_2.FourTi2` method), 27

`classified_solution_dicts()` (`sage.interfaces.phc.PHC_Object` method), 214

`clean_output()` (in module `sage.interfaces.mathematica`), 151

`cleaner()` (in module `sage.interfaces.cleaner`), 303

`clear()` (`sage.interfaces.giac.Giac` method), 68

`clear()` (`sage.interfaces.interface.Interface` method), 3

`clear()` (`sage.interfaces.magma.Magma` method), 123

`clear()` (`sage.interfaces.maple.Maple` method), 140

`clear()` (`sage.interfaces.maxima.Maxima` method), 163

`clear()` (`sage.interfaces.maxima_lib.MaximaLib` method), 184

`clear()` (`sage.interfaces.octave.Octave` method), 207

`clear()` (`sage.interfaces.sage0.Sage` method), 266

`clear()` (`sage.interfaces.singular.Singular` method), 280

`clear_prompts()` (`sage.interfaces.expect.Expect` method), 9

`close()` (`sage.interfaces.sagespawn.SageSpawn` method), 13

`cls()` (`sage.interfaces.macaulay2.Macaulay2Element` method), 111

`comma()` (`sage.interfaces.axiom.PanAxiomElement` method), 19

`comma()` (`sage.interfaces.maxima_abstract.MaximaAbstractElement` method), 175

`completions()` (`sage.interfaces.giac.Giac` method), 68

`completions()` (`sage.interfaces.maple.Maple` method), 140

`completions()` (`sage.interfaces.maxima_abstract.MaximaAbstract` method), 168

`completions()` (`sage.interfaces.mupad.Mupad` method), 198

`completions()` (`sage.interfaces.r.R` method), 255

`connected_subset()` (`sage.interfaces.qepcad.qepcad_formula_factory` method), 242

`console()` (in module `sage.interfaces.expect`), 11

`console()` (`sage.interfaces.axiom.Axiom` method), 17

`console()` (`sage.interfaces.fricas.FriCAS` method), 35

`console()` (`sage.interfaces.gap.Gap` method), 43

`console()` (`sage.interfaces.gap3.Gap3` method), 59

`console()` (`sage.interfaces.giac.Giac` method), 68

`console()` (`sage.interfaces.gnuplot.Gnuplot` method), 73

`console()` (`sage.interfaces.gp.Gp` method), 77

`console()` (`sage.interfaces.interface.Interface` method), 3

`console()` (`sage.interfaces.kash.Kash` method), 93

`console()` (`sage.interfaces.lie.LiE` method), 99

`console()` (`sage.interfaces.lisp.Lisp` method), 103

`console()` (`sage.interfaces.macaulay2.Macaulay2` method), 108

console() (sage.interfaces.magma.Magma method), 123  
 console() (sage.interfaces.maple.Maple method), 140  
 console() (sage.interfaces.mathematica.Mathematica method), 150  
 console() (sage.interfaces.matlab.Matlab method), 155  
 console() (sage.interfaces.maxima\_abstract.MaximaAbstract method), 168  
 console() (sage.interfaces.mupad.Mupad method), 198  
 console() (sage.interfaces.mwrank.Mwrank\_class method), 201  
 console() (sage.interfaces.octave.Octave method), 207  
 console() (sage.interfaces.r.R method), 255  
 console() (sage.interfaces.sage0.Sage method), 266  
 console() (sage.interfaces.scilab.Scilab method), 272  
 console() (sage.interfaces.singular.Singular method), 280  
 convert\_r\_list() (sage.interfaces.r.R method), 255  
 cputime() (sage.interfaces.gap.Gap method), 43  
 cputime() (sage.interfaces.gap3.Gap3 method), 59  
 cputime() (sage.interfaces.giac.Giac method), 68  
 cputime() (sage.interfaces.gp.Gp method), 77  
 cputime() (sage.interfaces.interface.Interface method), 3  
 cputime() (sage.interfaces.macaulay2.Macaulay2 method), 108  
 cputime() (sage.interfaces.magma.Magma method), 123  
 cputime() (sage.interfaces.maple.Maple method), 140  
 cputime() (sage.interfaces.maxima\_abstract.MaximaAbstract method), 168  
 cputime() (sage.interfaces.mupad.Mupad method), 198  
 cputime() (sage.interfaces.qsieve.qsieve\_nonblock method), 248  
 cputime() (sage.interfaces.sage0.Sage method), 267  
 cputime() (sage.interfaces.singular.Singular method), 280  
 CubexSolver (class in sage.interfaces.rubik), 263  
 current\_ring() (sage.interfaces.singular.Singular method), 280  
 current\_ring\_name() (sage.interfaces.singular.Singular method), 281

## D

data\_to\_list() (in module sage.interfaces.qsieve), 247  
 de\_solve() (sage.interfaces.maxima\_abstract.MaximaAbstract method), 168  
 de\_solve\_laplace() (sage.interfaces.maxima\_abstract.MaximaAbstract method), 169  
 de\_system\_plot() (sage.interfaces.octave.Octave method), 208  
 definition() (sage.interfaces.maxima\_abstract.MaximaAbstractElementFunction method), 181  
 demo() (sage.interfaces.maxima\_abstract.MaximaAbstract method), 169  
 derivative() (sage.interfaces.maxima\_abstract.MaximaAbstractElement method), 176  
 describe() (sage.interfaces.maxima\_abstract.MaximaAbstract method), 170  
 detach() (sage.interfaces.expect.Expect method), 9  
 diff() (sage.interfaces.maxima\_abstract.MaximaAbstractElement method), 176  
 DikSolver (class in sage.interfaces.rubik), 263  
 dimension() (sage.interfaces.frobby.Frobby method), 38  
 directory() (sage.interfaces.four\_ti\_2.FourTi2 method), 27  
 display2d() (sage.interfaces.maxima.MaximaElement method), 165  
 display2d() (sage.interfaces.maxima\_lib.MaximaLibElement method), 190  
 done() (sage.interfaces.qsieve.qsieve\_nonblock method), 248  
 dot() (sage.interfaces.macaulay2.Macaulay2Element method), 111  
 dot() (sage.interfaces.maxima\_abstract.MaximaAbstractElement method), 176  
 dot\_product() (sage.interfaces.r.RElement method), 260

`dummy_integrate()` (in module `sage.interfaces.maxima_lib`), 191

## E

`E()` (`sage.interfaces.qepcad.qepcad_formula_factory` method), 239  
`ecf()` (`sage.interfaces.maxima_lib.MaximaLibElement` method), 190  
`ECM` (class in `sage.interfaces.ecm`), 21  
`eval()` (`sage.interfaces.expect.Expect` method), 9  
`eval()` (`sage.interfaces.gap.Gap_generic` method), 46  
`eval()` (`sage.interfaces.giac.Giac` method), 69  
`eval()` (`sage.interfaces.interface.Interface` method), 3  
`eval()` (`sage.interfaces.kash.Kash` method), 93  
`eval()` (`sage.interfaces.lie.LiE` method), 99  
`eval()` (`sage.interfaces.lisp.Lisp` method), 104  
`eval()` (`sage.interfaces.macaulay2.Macaulay2` method), 108  
`eval()` (`sage.interfaces.magma.Magma` method), 124  
`eval()` (`sage.interfaces.magma.MagmaElement` method), 129  
`eval()` (`sage.interfaces.magma_free.MagmaFree` method), 135  
`eval()` (`sage.interfaces.mathematica.Mathematica` method), 150  
`eval()` (`sage.interfaces.maxima_lib.MaximaLib` method), 184  
`eval()` (`sage.interfaces.mupad.Mupad` method), 199  
`eval()` (`sage.interfaces.mwrank.Mwrank_class` method), 201  
`eval()` (`sage.interfaces.p sage.PSage` method), 219  
`eval()` (`sage.interfaces.r.R` method), 256  
`eval()` (`sage.interfaces.sage0.Sage` method), 267  
`eval()` (`sage.interfaces.scilab.Scilab` method), 272  
`eval()` (`sage.interfaces.singular.Singular` method), 281  
`evaluate()` (`sage.interfaces.magma.MagmaElement` method), 129  
`exactly_k()` (`sage.interfaces.qepcad.qepcad_formula_factory` method), 242  
`example()` (`sage.interfaces.maxima_abstract.MaximaAbstract` method), 170  
`execute()` (`sage.interfaces.interface.Interface` method), 3  
`exists()` (`sage.interfaces.qepcad.qepcad_formula_factory` method), 243  
`Expect` (class in `sage.interfaces.expect`), 9  
`expect()` (`sage.interfaces.expect.Expect` method), 10  
`expect()` (`sage.interfaces.giac.Giac` method), 69  
`expect()` (`sage.interfaces.maple.Maple` method), 140  
`expect()` (`sage.interfaces.mupad.Mupad` method), 199  
`expect_peek()` (`sage.interfaces.sagespawn.SageSpawn` method), 13  
`expect_quitall()` (in module `sage.interfaces.quit`), 305  
`expect_upto()` (`sage.interfaces.sagespawn.SageSpawn` method), 13  
`ExpectElement` (class in `sage.interfaces.expect`), 11  
`ExpectFunction` (class in `sage.interfaces.expect`), 11  
`export_image()` (`sage.interfaces.jmoldata.JmolData` method), 85  
`extcode_dir()` (in module `sage.interfaces.magma`), 133  
`external_string()` (`sage.interfaces.macaulay2.Macaulay2Element` method), 111

## F

`F()` (`sage.interfaces.qepcad.qepcad_formula_factory` method), 240  
`factor()` (`sage.interfaces.ecm.ECM` method), 22  
`final_stats()` (`sage.interfaces.qepcad.Qepcad` method), 230  
`find_factor()` (`sage.interfaces.ecm.ECM` method), 23

flush() (sage.interfaces.magma.MagmaGBLogPrettyPrinter method), 133  
 flush() (sage.interfaces.singular.SingularGBLogPrettyPrinter method), 292  
 forall() (sage.interfaces.qepcad.qepcad\_formula\_factory method), 243  
 format\_cube() (sage.interfaces.rubik.CubexSolver method), 263  
 format\_cube() (sage.interfaces.rubik.DikSolver method), 263  
 format\_cube() (sage.interfaces.rubik.OptimalSolver method), 263  
 formula() (sage.interfaces.qepcad.qepcad\_formula\_factory method), 243  
 FourTi2 (class in sage.interfaces.four\_ti\_2), 27  
 FriCAS (class in sage.interfaces.fricas), 35  
 fricas\_console() (in module sage.interfaces.fricas), 35  
 FriCASElement (class in sage.interfaces.fricas), 35  
 FriCASExpectFunction (class in sage.interfaces.fricas), 35  
 FriCASFunctionElement (class in sage.interfaces.fricas), 35  
 Frobbly (class in sage.interfaces.frobbly), 37  
 function() (sage.interfaces.maxima\_abstract.MaximaAbstract method), 170  
 function\_call() (sage.interfaces.gap.Gap\_generic method), 46  
 function\_call() (sage.interfaces.interface.Interface method), 3  
 function\_call() (sage.interfaces.lie.LiE method), 99  
 function\_call() (sage.interfaces.lisp.Lisp method), 104  
 function\_call() (sage.interfaces.magma.Magma method), 125  
 function\_call() (sage.interfaces.r.R method), 256  
 FunctionElement (class in sage.interfaces.expect), 11

## G

G() (sage.interfaces.qepcad.qepcad\_formula\_factory method), 240  
 Gap (class in sage.interfaces.gap), 43  
 Gap3 (class in sage.interfaces.gap3), 58  
 gap3\_console() (in module sage.interfaces.gap3), 60  
 gap3\_version() (in module sage.interfaces.gap3), 60  
 GAP3Element (class in sage.interfaces.gap3), 56  
 GAP3Record (class in sage.interfaces.gap3), 57  
 gap\_command() (in module sage.interfaces.gap), 48  
 gap\_console() (in module sage.interfaces.gap), 48  
 Gap\_generic (class in sage.interfaces.gap), 45  
 gap\_reset\_workspace() (in module sage.interfaces.gap), 49  
 GapElement (class in sage.interfaces.gap), 45  
 GapElement\_generic (class in sage.interfaces.gap), 45  
 GapFunction (class in sage.interfaces.gap), 45  
 GapFunctionElement (class in sage.interfaces.gap), 45  
 gc\_disabled (class in sage.interfaces.expect), 11  
 gen() (sage.interfaces.interface.InterfaceElement method), 5  
 gen() (sage.interfaces.magma.MagmaElement method), 129  
 gen\_names() (sage.interfaces.magma.MagmaElement method), 130  
 generate\_docstring\_dictionary() (in module sage.interfaces.singular), 292  
 genfiles\_mintides() (in module sage.interfaces.tides), 297  
 genfiles\_mpfr() (in module sage.interfaces.tides), 299  
 gens() (sage.interfaces.magma.MagmaElement method), 130  
 get() (sage.interfaces.axiom.PanAxiom method), 18  
 get() (sage.interfaces.gap.Gap method), 44  
 get() (sage.interfaces.giac.Giac method), 69

`get()` (sage.interfaces.gp.Gp method), 78  
`get()` (sage.interfaces.interface.Interface method), 3  
`get()` (sage.interfaces.kash.Kash method), 93  
`get()` (sage.interfaces.lie.LiE method), 100  
`get()` (sage.interfaces.lisp.Lisp method), 104  
`get()` (sage.interfaces.macaulay2.Macaulay2 method), 108  
`get()` (sage.interfaces.magma.Magma method), 125  
`get()` (sage.interfaces.maple.Maple method), 141  
`get()` (sage.interfaces.mathematica.Mathematica method), 150  
`get()` (sage.interfaces.matlab.Matlab method), 155  
`get()` (sage.interfaces.maxima.Maxima method), 164  
`get()` (sage.interfaces.maxima\_lib.MaximaLib method), 184  
`get()` (sage.interfaces.mupad.Mupad method), 199  
`get()` (sage.interfaces.octave.Octave method), 208  
`get()` (sage.interfaces.p sage.PSage method), 219  
`get()` (sage.interfaces.r.R method), 256  
`get()` (sage.interfaces.sage0.Sage method), 267  
`get()` (sage.interfaces.scilab.Scilab method), 272  
`get()` (sage.interfaces.singular.Singular method), 282  
`get_classified_solution_dicts()` (in module sage.interfaces.phc), 215  
`get_default()` (sage.interfaces.gp.Gp method), 78  
`get_docstring()` (in module sage.interfaces.singular), 292  
`get_gap_memory_pool_size()` (in module sage.interfaces.gap), 49  
`get_last_params()` (sage.interfaces.ecm.ECM method), 24  
`get_magma_attribute()` (sage.interfaces.magma.MagmaElement method), 131  
`get_precision()` (sage.interfaces.gp.Gp method), 78  
`get_real_precision()` (sage.interfaces.gp.Gp method), 78  
`get_record_element()` (sage.interfaces.gap.Gap\_generic method), 47  
`get_seed()` (sage.interfaces.interface.Interface method), 3  
`get_series_precision()` (sage.interfaces.gp.Gp method), 78  
`get_solution_dicts()` (in module sage.interfaces.phc), 216  
`get_using_file()` (sage.interfaces.interface.Interface method), 4  
`get_using_file()` (sage.interfaces.interface.InterfaceElement method), 5  
`get_using_file()` (sage.interfaces.lie.LiE method), 100  
`get_variable_list()` (in module sage.interfaces.phc), 216  
`get_verbose()` (sage.interfaces.magma.Magma method), 125  
`GetVerbose()` (sage.interfaces.magma.Magma method), 121  
`Gfan` (class in sage.interfaces.gfan), 63  
`gfq_gap_to_sage()` (in module sage.interfaces.gap), 50  
`Giac` (class in sage.interfaces.giac), 67  
`giac_console()` (in module sage.interfaces.giac), 72  
`GiacElement` (class in sage.interfaces.giac), 70  
`GiacFunction` (class in sage.interfaces.giac), 71  
`GiacFunctionElement` (class in sage.interfaces.giac), 71  
`Gnuplot` (class in sage.interfaces.gnuplot), 73  
`gnuplot()` (sage.interfaces.gnuplot.Gnuplot method), 73  
`gnuplot_console()` (in module sage.interfaces.gnuplot), 74  
`Gp` (class in sage.interfaces.gp), 77  
`gp_console()` (in module sage.interfaces.gp), 82  
`gp_version()` (in module sage.interfaces.gp), 82

GpElement (class in sage.interfaces.gp), 81  
 GpFunction (class in sage.interfaces.gp), 82  
 GpFunctionElement (class in sage.interfaces.gp), 82  
 graver() (sage.interfaces.four\_ti\_2.FourTi2 method), 28  
 groebner() (sage.interfaces.four\_ti\_2.FourTi2 method), 28

## H

hasattr() (sage.interfaces.interface.InterfaceElement method), 5  
 help() (sage.interfaces.gap.Gap method), 44  
 help() (sage.interfaces.gap3.Gap3 method), 59  
 help() (sage.interfaces.giac.Giac method), 69  
 help() (sage.interfaces.gp.Gp method), 79  
 help() (sage.interfaces.interface.Interface method), 4  
 help() (sage.interfaces.interface.InterfaceFunctionElement method), 7  
 help() (sage.interfaces.kash.Kash method), 93  
 help() (sage.interfaces.lie.LiE method), 100  
 help() (sage.interfaces.lisp.Lisp method), 104  
 help() (sage.interfaces.macaulay2.Macaulay2 method), 109  
 help() (sage.interfaces.magma.Magma method), 126  
 help() (sage.interfaces.maple.Maple method), 141  
 help() (sage.interfaces.mathematica.Mathematica method), 150  
 help() (sage.interfaces.maxima\_abstract.MaximaAbstract method), 171  
 help() (sage.interfaces.r.R method), 256  
 help() (sage.interfaces.tachyon.TachyonRT method), 296  
 help\_search() (sage.interfaces.kash.Kash method), 93  
 HelpExpression (class in sage.interfaces.r), 254  
 hilbert() (sage.interfaces.four\_ti\_2.FourTi2 method), 28  
 hilbert() (sage.interfaces.frobby.Frobby method), 38

## I

ideal() (sage.interfaces.macaulay2.Macaulay2 method), 109  
 ideal() (sage.interfaces.magma.Magma method), 126  
 ideal() (sage.interfaces.magma.MagmaElement method), 131  
 ideal() (sage.interfaces.singular.Singular method), 282  
 iff() (sage.interfaces.qepcad.qepcad\_formula\_factory method), 244  
 imag() (sage.interfaces.maxima\_abstract.MaximaAbstractElement method), 177  
 implies() (sage.interfaces.qepcad.qepcad\_formula\_factory method), 244  
 index() (sage.interfaces.qepcad.QepcadCell method), 233  
 infinitely\_many() (sage.interfaces.qepcad.qepcad\_formula\_factory method), 244  
 install\_packages() (sage.interfaces.r.R method), 257  
 integral() (sage.interfaces.giac.GiacElement method), 70  
 integral() (sage.interfaces.maxima\_abstract.MaximaAbstractElement method), 177  
 integral() (sage.interfaces.maxima\_abstract.MaximaAbstractElementFunction method), 181  
 integrate() (sage.interfaces.giac.GiacElement method), 70  
 integrate() (sage.interfaces.maxima\_abstract.MaximaAbstractElement method), 177  
 integrate() (sage.interfaces.maxima\_abstract.MaximaAbstractElementFunction method), 182  
 interact() (sage.interfaces.ecm.ECM method), 24  
 interact() (sage.interfaces.gnuplot.Gnuplot method), 73  
 interact() (sage.interfaces.interface.Interface method), 4  
 Interface (class in sage.interfaces.interface), 3



InterfaceElement (class in sage.interfaces.interface), 5  
InterfaceFunction (class in sage.interfaces.interface), 6  
InterfaceFunctionElement (class in sage.interfaces.interface), 6  
interrupt() (sage.interfaces.expect.Expect method), 10  
interrupt() (sage.interfaces.gap.Gap\_generic method), 47  
intmod\_gap\_to\_sage() (in module sage.interfaces.gap), 50  
invalidate\_all() (in module sage.interfaces.quit), 305  
irreducible\_decomposition() (sage.interfaces.frobby.Frobby method), 39  
is\_AxiomElement() (in module sage.interfaces.axiom), 20  
is\_ExpectElement() (in module sage.interfaces.expect), 11  
is\_FriCASElement() (in module sage.interfaces.fricas), 36  
is\_GapElement() (in module sage.interfaces.gap), 51  
is\_GpElement() (in module sage.interfaces.gp), 83  
is\_InterfaceElement() (in module sage.interfaces.interface), 7  
is\_jvm\_available() (sage.interfaces.jmoldata.JmolData method), 86  
is\_KashElement() (in module sage.interfaces.kash), 94  
is\_LiEElement() (in module sage.interfaces.lie), 101  
is\_LispElement() (in module sage.interfaces.lisp), 105  
is\_local() (sage.interfaces.expect.Expect method), 10  
is\_locked() (sage.interfaces.p sage.PSage method), 219  
is\_locked() (sage.interfaces.p sage.PSageElement method), 220  
is\_Macaulay2Element() (in module sage.interfaces.macaulay2), 114  
is\_MagmaElement() (in module sage.interfaces.magma), 133  
is\_MaximaElement() (in module sage.interfaces.maxima), 165  
is\_MaximaLibElement() (in module sage.interfaces.maxima\_lib), 191  
is\_RElement() (in module sage.interfaces.r), 262  
is\_remote() (sage.interfaces.expect.Expect method), 10  
is\_running() (in module sage.interfaces.quit), 305  
is\_running() (sage.interfaces.expect.Expect method), 10  
is\_SingularElement() (in module sage.interfaces.singular), 292  
is\_string() (sage.interfaces.gap.GapElement\_generic method), 45  
is\_string() (sage.interfaces.gp.GpElement method), 82  
is\_string() (sage.interfaces.interface.InterfaceElement method), 6  
is\_string() (sage.interfaces.singular.SingularElement method), 287

## J

JmolData (class in sage.interfaces.jmoldata), 85

## K

Kash (class in sage.interfaces.kash), 93  
kash\_console() (in module sage.interfaces.kash), 94  
kash\_version() (in module sage.interfaces.kash), 94  
KashDocumentation (class in sage.interfaces.kash), 93  
KashElement (class in sage.interfaces.kash), 94  
kill() (sage.interfaces.gp.Gp method), 79  
kill() (sage.interfaces.lisp.Lisp method), 104  
kill\_spawned\_jobs() (in module sage.interfaces.quit), 305

## L

level() (sage.interfaces.qepcad.QepcadCell method), 233



LIB() (sage.interfaces.singular.Singular method), 280  
 lib() (sage.interfaces.singular.Singular method), 282  
 library() (sage.interfaces.r.R method), 257  
 LiE (class in sage.interfaces.lie), 99  
 lie\_console() (in module sage.interfaces.lie), 101  
 lie\_version() (in module sage.interfaces.lie), 102  
 LiEElement (class in sage.interfaces.lie), 101  
 LiEFunction (class in sage.interfaces.lie), 101  
 LiEFunctionElement (class in sage.interfaces.lie), 101  
 Lisp (class in sage.interfaces.lisp), 103  
 lisp() (sage.interfaces.maxima.Maxima method), 164  
 lisp() (sage.interfaces.maxima\_lib.MaximaLib method), 185  
 lisp\_console() (in module sage.interfaces.lisp), 105  
 LispElement (class in sage.interfaces.lisp), 105  
 LispFunction (class in sage.interfaces.lisp), 105  
 LispFunctionElement (class in sage.interfaces.lisp), 105  
 list() (sage.interfaces.qsieve.qsieve\_nonblock method), 248  
 list() (sage.interfaces.singular.Singular method), 283  
 list\_attributes() (sage.interfaces.magma.MagmaElement method), 131  
 load() (sage.interfaces.magma.Magma method), 126  
 load() (sage.interfaces.maple.Maple method), 141  
 load() (sage.interfaces.singular.Singular method), 283  
 load\_package() (sage.interfaces.gap.Gap\_generic method), 47  
 log() (sage.interfaces.qsieve.qsieve\_nonblock method), 248

## M

Macaulay2 (class in sage.interfaces.macaulay2), 108  
 macaulay2\_console() (in module sage.interfaces.macaulay2), 114  
 Macaulay2Element (class in sage.interfaces.macaulay2), 111  
 Macaulay2Function (class in sage.interfaces.macaulay2), 114  
 Magma (class in sage.interfaces.magma), 120  
 magma\_console() (in module sage.interfaces.magma), 134  
 magma\_free\_eval() (in module sage.interfaces.magma\_free), 135  
 magma\_gb\_standard\_options() (in module sage.interfaces.magma), 134  
 magma\_version() (in module sage.interfaces.magma), 134  
 MagmaElement (class in sage.interfaces.magma), 128  
 MagmaExpr (class in sage.interfaces.magma\_free), 135  
 MagmaFree (class in sage.interfaces.magma\_free), 135  
 MagmaFunction (class in sage.interfaces.magma), 133  
 MagmaFunctionElement (class in sage.interfaces.magma), 133  
 MagmaGBDefaultContext (class in sage.interfaces.magma), 133  
 MagmaGBLogPrettyPrinter (class in sage.interfaces.magma), 133  
 make\_cells() (sage.interfaces.qepcad.Qepcad method), 231  
 Maple (class in sage.interfaces.maple), 139  
 maple\_console() (in module sage.interfaces.maple), 142  
 MapleElement (class in sage.interfaces.maple), 142  
 MapleFunction (class in sage.interfaces.maple), 142  
 MapleFunctionElement (class in sage.interfaces.maple), 142  
 Mathematica (class in sage.interfaces.mathematica), 150  
 mathematica\_console() (in module sage.interfaces.mathematica), 151

MathematicaElement (class in sage.interfaces.mathematica), 151  
MathematicaFunction (class in sage.interfaces.mathematica), 151  
MathematicaFunctionElement (class in sage.interfaces.mathematica), 151  
Matlab (class in sage.interfaces.matlab), 155  
matlab\_console() (in module sage.interfaces.matlab), 156  
matlab\_version() (in module sage.interfaces.matlab), 156  
MatlabElement (class in sage.interfaces.matlab), 156  
matrix() (sage.interfaces.singular.Singular method), 283  
max\_at\_to\_sage() (in module sage.interfaces.maxima\_lib), 191  
max\_to\_sr() (in module sage.interfaces.maxima\_lib), 192  
max\_to\_string() (in module sage.interfaces.maxima\_lib), 192  
Maxima (class in sage.interfaces.maxima), 163  
maxima\_console() (in module sage.interfaces.maxima\_abstract), 182  
maxima\_version() (in module sage.interfaces.maxima\_abstract), 182  
MaximaAbstract (class in sage.interfaces.maxima\_abstract), 167  
MaximaAbstractElement (class in sage.interfaces.maxima\_abstract), 175  
MaximaAbstractElementFunction (class in sage.interfaces.maxima\_abstract), 180  
MaximaAbstractFunction (class in sage.interfaces.maxima\_abstract), 182  
MaximaAbstractFunctionElement (class in sage.interfaces.maxima\_abstract), 182  
MaximaElement (class in sage.interfaces.maxima), 164  
MaximaElementFunction (class in sage.interfaces.maxima), 165  
MaximaFunction (class in sage.interfaces.maxima), 165  
MaximaFunctionElement (class in sage.interfaces.maxima), 165  
MaximaLib (class in sage.interfaces.maxima\_lib), 183  
MaximaLibElement (class in sage.interfaces.maxima\_lib), 189  
MaximaLibElementFunction (class in sage.interfaces.maxima\_lib), 190  
MaximaLibFunction (class in sage.interfaces.maxima\_lib), 191  
MaximaLibFunctionElement (class in sage.interfaces.maxima\_lib), 191  
mdiff\_to\_sage() (in module sage.interfaces.maxima\_lib), 192  
methods() (sage.interfaces.magma.MagmaElement method), 131  
minimize() (sage.interfaces.four\_ti\_2.FourTi2 method), 28  
mixed\_volume() (sage.interfaces.phc.PHC method), 212  
mlist\_to\_sage() (in module sage.interfaces.maxima\_lib), 193  
mqapply\_to\_sage() (in module sage.interfaces.maxima\_lib), 193  
mrat\_to\_sage() (in module sage.interfaces.maxima\_lib), 193  
Mupad (class in sage.interfaces.mupad), 198  
mupad\_console() (in module sage.interfaces.mupad), 200  
MupadElement (class in sage.interfaces.mupad), 199  
MupadFunction (class in sage.interfaces.mupad), 199  
MupadFunctionElement (class in sage.interfaces.mupad), 200  
Mwrank() (in module sage.interfaces.mwrank), 201  
Mwrank\_class (class in sage.interfaces.mwrank), 201  
mwrank\_console() (in module sage.interfaces.mwrank), 202

## N

N() (sage.interfaces.mathematica.MathematicaElement method), 151  
n() (sage.interfaces.qsieve.qsieve\_nonblock method), 248  
na() (sage.interfaces.r.R method), 257  
name() (sage.interfaces.interface.Interface method), 4  
name() (sage.interfaces.interface.InterfaceElement method), 6

[new\(\)](#) (sage.interfaces.interface.Interface method), 4  
[new\(\)](#) (sage.interfaces.sage0.Sage method), 267  
[new\\_from\(\)](#) (sage.interfaces.macaulay2.Macaulay2 method), 109  
[new\\_with\\_bits\\_prec\(\)](#) (sage.interfaces.gp.Gp method), 79  
[nintegral\(\)](#) (sage.interfaces.maxima\_abstract.MaximaAbstractElement method), 178  
[not\\_\(\)](#) (sage.interfaces.qepcad.qepcad\_formula\_factory method), 244  
[number\\_of\\_children\(\)](#) (sage.interfaces.qepcad.QepcadCell method), 234  
[numer\(\)](#) (sage.interfaces.maxima\_abstract.MaximaAbstractElement method), 179

## O

[objgens\(\)](#) (sage.interfaces.magma.Magma method), 127  
[Octave](#) (class in sage.interfaces.octave), 207  
[octave\\_console\(\)](#) (in module sage.interfaces.octave), 209  
[octave\\_version\(\)](#) (in module sage.interfaces.octave), 210  
[OctaveElement](#) (class in sage.interfaces.octave), 209  
[one\\_curve\(\)](#) (sage.interfaces.ecm.ECM method), 24  
[operations\(\)](#) (sage.interfaces.gap3.GAP3Record method), 57  
[OptimalSolver](#) (class in sage.interfaces.rubik), 263  
[option\(\)](#) (sage.interfaces.singular.Singular method), 283  
[or\\_\(\)](#) (sage.interfaces.qepcad.qepcad\_formula\_factory method), 245

## P

[PanAxiom](#) (class in sage.interfaces.axiom), 18  
[PanAxiomElement](#) (class in sage.interfaces.axiom), 18  
[PanAxiomExpectFunction](#) (class in sage.interfaces.axiom), 20  
[PanAxiomFunctionElement](#) (class in sage.interfaces.axiom), 20  
[parse\\_max\\_string\(\)](#) (in module sage.interfaces.maxima\_lib), 194  
[partial\\_fraction\\_decomposition\(\)](#) (sage.interfaces.maxima\_abstract.MaximaAbstractElement method), 179  
[path\(\)](#) (sage.interfaces.expect.Expect method), 10  
[path\\_track\(\)](#) (sage.interfaces.phc.PHC method), 212  
[phase\(\)](#) (sage.interfaces.qepcad.Qepcad method), 231  
[PHC](#) (class in sage.interfaces.phc), 211  
[PHC\\_Object](#) (class in sage.interfaces.phc), 214  
[pid\(\)](#) (sage.interfaces.expect.Expect method), 10  
[pid\(\)](#) (sage.interfaces.qsieve.qsieve\_nonblock method), 248  
[plot\(\)](#) (sage.interfaces.gnuplot.Gnuplot method), 73  
[plot\(\)](#) (sage.interfaces.r.R method), 257  
[plot2d\(\)](#) (sage.interfaces.maxima\_abstract.MaximaAbstract method), 171  
[plot2d\\_parametric\(\)](#) (sage.interfaces.maxima\_abstract.MaximaAbstract method), 172  
[plot3d\(\)](#) (sage.interfaces.gnuplot.Gnuplot method), 73  
[plot3d\(\)](#) (sage.interfaces.maxima\_abstract.MaximaAbstract method), 172  
[plot3d\\_parametric\(\)](#) (sage.interfaces.gnuplot.Gnuplot method), 73  
[plot3d\\_parametric\(\)](#) (sage.interfaces.maxima\_abstract.MaximaAbstract method), 172  
[plot\\_list\(\)](#) (sage.interfaces.maxima\_abstract.MaximaAbstract method), 173  
[plot\\_multilist\(\)](#) (sage.interfaces.maxima\_abstract.MaximaAbstract method), 173  
[plot\\_paths\\_2d\(\)](#) (sage.interfaces.phc.PHC method), 212  
[png\(\)](#) (sage.interfaces.r.R method), 258  
[POVRay](#) (class in sage.interfaces.povray), 217  
[ppi\(\)](#) (sage.interfaces.four\_ti\_2.FourTi2 method), 29  
[preparse\(\)](#) (sage.interfaces.sage0.Sage method), 267

PSage (class in sage.interfaces.psage), 219  
PSageElement (class in sage.interfaces.psage), 219  
pyobject\_to\_max() (in module sage.interfaces.maxima\_lib), 194

## Q

Qepcad (class in sage.interfaces.qepcad), 229  
qepcad() (in module sage.interfaces.qepcad), 235  
qepcad\_banner() (in module sage.interfaces.qepcad), 238  
qepcad\_console() (in module sage.interfaces.qepcad), 239  
Qepcad\_expect (class in sage.interfaces.qepcad), 235  
qepcad\_formula\_factory (class in sage.interfaces.qepcad), 239  
qepcad\_version() (in module sage.interfaces.qepcad), 245  
QepcadCell (class in sage.interfaces.qepcad), 233  
QepcadFunction (class in sage.interfaces.qepcad), 235  
qformula (class in sage.interfaces.qepcad), 246  
qsieve() (in module sage.interfaces.qsieve), 247  
qsieve\_block() (in module sage.interfaces.qsieve), 248  
qsieve\_nonblock (class in sage.interfaces.qsieve), 248  
qsolve() (sage.interfaces.four\_ti\_2.FourTi2 method), 29  
quantifier() (sage.interfaces.qepcad.qepcad\_formula\_factory method), 245  
quit() (sage.interfaces.expect.Expect method), 10  
quit() (sage.interfaces.octave.Octave method), 208  
quit() (sage.interfaces.qsieve.qsieve\_nonblock method), 248  
quo() (sage.interfaces.magma.MagmaElement method), 131

## R

R (class in sage.interfaces.r), 254  
r\_console() (in module sage.interfaces.r), 262  
r\_version() (in module sage.interfaces.r), 262  
rand\_seed() (sage.interfaces.interface.Interface method), 4  
rays() (sage.interfaces.four\_ti\_2.FourTi2 method), 29  
read() (sage.interfaces.interface.Interface method), 4  
read() (sage.interfaces.lie.LiE method), 100  
read() (sage.interfaces.r.R method), 259  
read\_data() (in module sage.interfaces.read\_data), 307  
read\_matrix() (sage.interfaces.four\_ti\_2.FourTi2 method), 29  
ready() (sage.interfaces.rubik.OptimalSolver method), 263  
real() (sage.interfaces.maxima\_abstract.MaximaAbstractElement method), 179  
recfields() (sage.interfaces.gap3.GAP3Record method), 57  
recommended\_B1() (sage.interfaces.ecm.ECM method), 25  
reduce\_load() (in module sage.interfaces.gap), 51  
reduce\_load() (in module sage.interfaces.mathematica), 152  
reduce\_load() (in module sage.interfaces.singular), 292  
reduce\_load\_Axiom() (in module sage.interfaces.axiom), 20  
reduce\_load\_element() (in module sage.interfaces.sage0), 268  
reduce\_load\_fricas() (in module sage.interfaces.fricas), 36  
reduce\_load\_GAP() (in module sage.interfaces.gap), 51  
reduce\_load\_Giac() (in module sage.interfaces.giac), 72  
reduce\_load\_GP() (in module sage.interfaces.gp), 83  
reduce\_load\_Kash() (in module sage.interfaces.kash), 94

`reduce_load_lie()` (in module `sage.interfaces.lie`), 102  
`reduce_load_Lisp()` (in module `sage.interfaces.lisp`), 106  
`reduce_load_macaulay2()` (in module `sage.interfaces.macaulay2`), 114  
`reduce_load_Magma()` (in module `sage.interfaces.magma`), 134  
`reduce_load_Maple()` (in module `sage.interfaces.maple`), 143  
`reduce_load_Matlab()` (in module `sage.interfaces.matlab`), 156  
`reduce_load_Maxima()` (in module `sage.interfaces.maxima`), 165  
`reduce_load_Maxima_function()` (in module `sage.interfaces.maxima`), 165  
`reduce_load_MaximaAbstract_function()` (in module `sage.interfaces.maxima_abstract`), 182  
`reduce_load_MaximaLib()` (in module `sage.interfaces.maxima_lib`), 194  
`reduce_load_mupad()` (in module `sage.interfaces.mupad`), 200  
`reduce_load_Octave()` (in module `sage.interfaces.octave`), 210  
`reduce_load_R()` (in module `sage.interfaces.r`), 262  
`reduce_load_Sage()` (in module `sage.interfaces.sage0`), 268  
`reduce_load_Singular()` (in module `sage.interfaces.singular`), 293  
`RElement` (class in `sage.interfaces.r`), 260  
`remove_constants()` (in module `sage.interfaces.tides`), 300  
`remove_output_labels()` (in module `sage.interfaces.macaulay2`), 115  
`remove_repeated()` (in module `sage.interfaces.tides`), 300  
`repr()` (`sage.interfaces.macaulay2.Macaulay2Element` method), 111  
`require()` (`sage.interfaces.r.R` method), 259  
`restart()` (`sage.interfaces.macaulay2.Macaulay2` method), 109  
`RFunction` (class in `sage.interfaces.r`), 261  
`RFunctionElement` (class in `sage.interfaces.r`), 262  
`ring()` (`sage.interfaces.macaulay2.Macaulay2` method), 110  
`ring()` (`sage.interfaces.singular.Singular` method), 284

## S

`Sage` (class in `sage.interfaces.sage0`), 265  
`sage()` (`sage.interfaces.interface.InterfaceElement` method), 6  
`sage.interfaces.axiom` (module), 15  
`sage.interfaces.cleaner` (module), 303  
`sage.interfaces.ecm` (module), 21  
`sage.interfaces.expect` (module), 9  
`sage.interfaces.four_ti_2` (module), 27  
`sage.interfaces.fricas` (module), 33  
`sage.interfaces.frobby` (module), 37  
`sage.interfaces.gap` (module), 41  
`sage.interfaces.gap3` (module), 53  
`sage.interfaces.gfan` (module), 63  
`sage.interfaces.giac` (module), 65  
`sage.interfaces.gnuplot` (module), 73  
`sage.interfaces.gp` (module), 75  
`sage.interfaces.interface` (module), 3  
`sage.interfaces.jmoldata` (module), 85  
`sage.interfaces.kash` (module), 87  
`sage.interfaces.lie` (module), 95  
`sage.interfaces.lisp` (module), 103  
`sage.interfaces.macaulay2` (module), 107  
`sage.interfaces.magma` (module), 117

`sage.interfaces.magma_free` (module), 135  
`sage.interfaces.maple` (module), 137  
`sage.interfaces.mathematica` (module), 145  
`sage.interfaces.matlab` (module), 153  
`sage.interfaces.maxima` (module), 157  
`sage.interfaces.maxima_abstract` (module), 167  
`sage.interfaces.maxima_lib` (module), 183  
`sage.interfaces.mupad` (module), 197  
`sage.interfaces.mwrank` (module), 201  
`sage.interfaces.octave` (module), 205  
`sage.interfaces.phc` (module), 211  
`sage.interfaces.povray` (module), 217  
`sage.interfaces.psage` (module), 219  
`sage.interfaces.qepcad` (module), 221  
`sage.interfaces.qsieve` (module), 247  
`sage.interfaces.quit` (module), 305  
`sage.interfaces.r` (module), 251  
`sage.interfaces.read_data` (module), 307  
`sage.interfaces.rubik` (module), 263  
`sage.interfaces.sage0` (module), 265  
`sage.interfaces.sagespawn` (module), 13  
`sage.interfaces.scilab` (module), 269  
`sage.interfaces.singular` (module), 275  
`sage.interfaces.tachyon` (module), 295  
`sage.interfaces.tides` (module), 297  
`sage0_console()` (in module `sage.interfaces.sage0`), 268  
`sage0_version()` (in module `sage.interfaces.sage0`), 268  
`sage2matlab_matrix_string()` (`sage.interfaces.matlab.Matlab` method), 155  
`sage2octave_matrix_string()` (`sage.interfaces.octave.Octave` method), 208  
`sage2scilab_matrix_string()` (`sage.interfaces.scilab.Scilab` method), 272  
`sage_flattened_str_list()` (`sage.interfaces.singular.SingularElement` method), 287  
`sage_global_ring()` (`sage.interfaces.singular.SingularElement` method), 287  
`sage_matrix()` (`sage.interfaces.singular.SingularElement` method), 288  
`sage_poly()` (`sage.interfaces.singular.SingularElement` method), 289  
`sage_polystring()` (`sage.interfaces.macaulay2.Macaulay2Element` method), 112  
`sage_polystring()` (`sage.interfaces.singular.SingularElement` method), 290  
`sage_rat()` (in module `sage.interfaces.maxima_lib`), 194  
`sage_structured_str_list()` (`sage.interfaces.singular.SingularElement` method), 290  
`SageElement` (class in `sage.interfaces.sage0`), 268  
`SageFunction` (class in `sage.interfaces.sage0`), 268  
`SageSpawn` (class in `sage.interfaces.sagespawn`), 13  
`sample_point()` (`sage.interfaces.qepcad.QepcadCell` method), 234  
`sample_point_dict()` (`sage.interfaces.qepcad.QepcadCell` method), 234  
`save_as_start()` (`sage.interfaces.phc.PHC_Object` method), 214  
`save_image()` (`sage.interfaces.mathematica.MathematicaElement` method), 151  
`save_workspace()` (`sage.interfaces.gap.Gap` method), 44  
`Scilab` (class in `sage.interfaces.scilab`), 271  
`scilab_console()` (in module `sage.interfaces.scilab`), 274  
`scilab_version()` (in module `sage.interfaces.scilab`), 274  
`ScilabElement` (class in `sage.interfaces.scilab`), 273

set() (sage.interfaces.axiom.PanAxiom method), 18  
 set() (sage.interfaces.gap.Gap method), 44  
 set() (sage.interfaces.giac.Giac method), 70  
 set() (sage.interfaces.gp.Gp method), 79  
 set() (sage.interfaces.interface.Interface method), 5  
 set() (sage.interfaces.kash.Kash method), 93  
 set() (sage.interfaces.lie.LiE method), 100  
 set() (sage.interfaces.lisp.Lisp method), 104  
 set() (sage.interfaces.macaulay2.Macaulay2 method), 110  
 set() (sage.interfaces.magma.Magma method), 127  
 set() (sage.interfaces.maple.Maple method), 141  
 set() (sage.interfaces.mathematica.Mathematica method), 150  
 set() (sage.interfaces.matlab.Matlab method), 155  
 set() (sage.interfaces.matlab.MatlabElement method), 156  
 set() (sage.interfaces.maxima.Maxima method), 164  
 set() (sage.interfaces.maxima\_lib.MaximaLib method), 185  
 set() (sage.interfaces.mupad.Mupad method), 199  
 set() (sage.interfaces.octave.Octave method), 209  
 set() (sage.interfaces.psage.PSage method), 219  
 set() (sage.interfaces.r.R method), 259  
 set() (sage.interfaces.sage0.Sage method), 267  
 set() (sage.interfaces.scilab.Scilab method), 273  
 set() (sage.interfaces.scilab.ScilabElement method), 273  
 set() (sage.interfaces.singular.Singular method), 285  
 set\_default() (sage.interfaces.gp.Gp method), 80  
 set\_gap\_memory\_pool\_size() (in module sage.interfaces.gap), 51  
 set\_magma\_attribute() (sage.interfaces.magma.MagmaElement method), 132  
 set\_precision() (sage.interfaces.gp.Gp method), 80  
 set\_real\_precision() (sage.interfaces.gp.Gp method), 80  
 set\_ring() (sage.interfaces.singular.Singular method), 285  
 set\_ring() (sage.interfaces.singular.SingularElement method), 290  
 set\_seed() (sage.interfaces.gap.Gap method), 44  
 set\_seed() (sage.interfaces.gp.Gp method), 81  
 set\_seed() (sage.interfaces.interface.Interface method), 5  
 set\_seed() (sage.interfaces.magma.Magma method), 127  
 set\_seed() (sage.interfaces.maxima.Maxima method), 164  
 set\_seed() (sage.interfaces.octave.Octave method), 209  
 set\_seed() (sage.interfaces.r.R method), 259  
 set\_seed() (sage.interfaces.scilab.Scilab method), 273  
 set\_seed() (sage.interfaces.singular.Singular method), 285  
 set\_series\_precision() (sage.interfaces.gp.Gp method), 81  
 set\_truth() (sage.interfaces.qepcad.QepcadCell method), 234  
 set\_truth\_value() (sage.interfaces.qepcad.Qepcad method), 231  
 set\_verbose() (sage.interfaces.magma.Magma method), 127  
 setring() (sage.interfaces.singular.Singular method), 286  
 SetVerbose() (sage.interfaces.magma.Magma method), 121  
 sharp() (sage.interfaces.macaulay2.Macaulay2Element method), 112  
 show() (sage.interfaces.mathematica.MathematicaElement method), 151  
 signs() (sage.interfaces.qepcad.QepcadCell method), 235  
 SingNot (class in sage.interfaces.rubik), 264



Singular (class in sage.interfaces.singular), 279  
singular\_console() (in module sage.interfaces.singular), 293  
singular\_gb\_standard\_options() (in module sage.interfaces.singular), 293  
singular\_version() (in module sage.interfaces.singular), 294  
SingularElement (class in sage.interfaces.singular), 286  
SingularError, 291  
SingularFunction (class in sage.interfaces.singular), 291  
SingularFunctionElement (class in sage.interfaces.singular), 291  
SingularGBDefaultContext (class in sage.interfaces.singular), 291  
SingularGBLogPrettyPrinter (class in sage.interfaces.singular), 291  
solution\_dicts() (sage.interfaces.phc.PHC\_Object method), 214  
solution\_extension() (sage.interfaces.qepcad.Qepcad method), 232  
solutions() (sage.interfaces.phc.PHC\_Object method), 215  
solve() (sage.interfaces.rubik.CubexSolver method), 263  
solve() (sage.interfaces.rubik.DikSolver method), 263  
solve() (sage.interfaces.rubik.OptimalSolver method), 263  
solve\_linear() (sage.interfaces.maxima\_abstract.MaximaAbstract method), 174  
solve\_linear\_system() (sage.interfaces.octave.Octave method), 209  
source() (sage.interfaces.maple.Maple method), 141  
source() (sage.interfaces.r.R method), 259  
sr\_integral() (sage.interfaces.maxima\_lib.MaximaLib method), 185  
sr\_limit() (sage.interfaces.maxima\_lib.MaximaLib method), 187  
sr\_sum() (sage.interfaces.maxima\_lib.MaximaLib method), 188  
sr\_tlimit() (sage.interfaces.maxima\_lib.MaximaLib method), 189  
sr\_to\_max() (in module sage.interfaces.maxima\_lib), 195  
starstar() (sage.interfaces.macaulay2.Macaulay2Element method), 112  
start() (sage.interfaces.rubik.OptimalSolver method), 264  
start\_cleaner() (in module sage.interfaces.cleaner), 303  
start\_from() (sage.interfaces.phc.PHC method), 213  
stat\_model() (sage.interfaces.r.RElement method), 261  
stdout\_to\_string() (in module sage.interfaces.maxima\_lib), 195  
StdOutContext (class in sage.interfaces.expect), 11  
stop() (sage.interfaces.rubik.OptimalSolver method), 264  
str() (sage.interfaces.gap.GapElement method), 45  
str() (sage.interfaces.mathematica.MathematicaElement method), 151  
str() (sage.interfaces.maxima\_abstract.MaximaAbstractElement method), 180  
string() (sage.interfaces.singular.Singular method), 286  
strip\_answer() (sage.interfaces.matlab.Matlab method), 156  
structure\_sheaf() (sage.interfaces.macaulay2.Macaulay2Element method), 112  
sub() (sage.interfaces.magma.MagmaElement method), 132  
subexpressions\_list() (in module sage.interfaces.tides), 301  
subs() (sage.interfaces.macaulay2.Macaulay2Element method), 112  
subst() (sage.interfaces.maxima\_abstract.MaximaAbstractElement method), 180  
substitute() (sage.interfaces.macaulay2.Macaulay2Element method), 113  
sum() (sage.interfaces.giac.GiacElement method), 71

## T

TachyonRT (class in sage.interfaces.tachyon), 295  
temp\_project() (sage.interfaces.four\_ti\_2.FourTi2 method), 29  
terminate\_async() (sage.interfaces.sagespawn.SageSpawn method), 14



tilde() (sage.interfaces.r.RElement method), 261  
 time() (sage.interfaces.ecm.ECM method), 25  
 time() (sage.interfaces.qsieve.qsieve\_nonblock method), 249  
 tmpdir() (in module sage.interfaces.qsieve), 249  
 to\_poly\_solve() (sage.interfaces.maxima\_lib.MaximaLibElement method), 190  
 to\_sage() (sage.interfaces.macaulay2.Macaulay2Element method), 113  
 trait\_names() (sage.interfaces.axiom.PanAxiom method), 18  
 trait\_names() (sage.interfaces.gap.Gap method), 44  
 trait\_names() (sage.interfaces.gap.Gap\_generic method), 48  
 trait\_names() (sage.interfaces.gap.GapElement method), 45  
 trait\_names() (sage.interfaces.gap3.GAP3Record method), 58  
 trait\_names() (sage.interfaces.giac.Giac method), 70  
 trait\_names() (sage.interfaces.giac.GiacElement method), 71  
 trait\_names() (sage.interfaces.gp.Gp method), 81  
 trait\_names() (sage.interfaces.gp.GpElement method), 82  
 trait\_names() (sage.interfaces.lie.LiE method), 100  
 trait\_names() (sage.interfaces.lie.LiEElement method), 101  
 trait\_names() (sage.interfaces.lisp.Lisp method), 105  
 trait\_names() (sage.interfaces.macaulay2.Macaulay2 method), 110  
 trait\_names() (sage.interfaces.macaulay2.Macaulay2Element method), 114  
 trait\_names() (sage.interfaces.magma.Magma method), 128  
 trait\_names() (sage.interfaces.magma.MagmaElement method), 133  
 trait\_names() (sage.interfaces.maple.Maple method), 142  
 trait\_names() (sage.interfaces.maple.MapleElement method), 142  
 trait\_names() (sage.interfaces.mathematica.Mathematica method), 151  
 trait\_names() (sage.interfaces.maxima\_abstract.MaximaAbstract method), 174  
 trait\_names() (sage.interfaces.maxima\_abstract.MaximaAbstractElement method), 180  
 trait\_names() (sage.interfaces.mupad.Mupad method), 199  
 trait\_names() (sage.interfaces.mupad.MupadElement method), 199  
 trait\_names() (sage.interfaces.mupad.MupadFunction method), 199  
 trait\_names() (sage.interfaces.mupad.MupadFunctionElement method), 200  
 trait\_names() (sage.interfaces.qepcad.Qepcad method), 233  
 trait\_names() (sage.interfaces.r.R method), 260  
 trait\_names() (sage.interfaces.r.RElement method), 261  
 trait\_names() (sage.interfaces.sage0.Sage method), 267  
 trait\_names() (sage.interfaces.singular.Singular method), 286  
 trait\_names() (sage.interfaces.singular.SingularElement method), 291  
 type() (sage.interfaces.axiom.PanAxiomElement method), 19  
 type() (sage.interfaces.lie.LiEElement method), 101  
 type() (sage.interfaces.singular.SingularElement method), 291

## U

unapply() (sage.interfaces.giac.GiacElement method), 71  
 unbind() (sage.interfaces.gap.Gap\_generic method), 48  
 underscore() (sage.interfaces.macaulay2.Macaulay2Element method), 114  
 unit\_quadratic\_integer() (sage.interfaces.maxima\_abstract.MaximaAbstract method), 174  
 unparsed\_input\_form() (sage.interfaces.axiom.PanAxiomElement method), 19  
 usage() (sage.interfaces.povray.POVRay method), 217  
 usage() (sage.interfaces.tachyon.TachyonRT method), 296  
 use() (sage.interfaces.macaulay2.Macaulay2 method), 110

`user_dir()` (`sage.interfaces.expect.Expect` method), 11

## V

`validate_mwrank_input()` (in module `sage.interfaces.mwrank`), 202  
`variable_list()` (`sage.interfaces.phc.PHC_Object` method), 215  
`version()` (`sage.interfaces.gap.Gap_generic` method), 48  
`version()` (`sage.interfaces.giac.Giac` method), 70  
`version()` (`sage.interfaces.gp.Gp` method), 81  
`version()` (`sage.interfaces.kash.Kash` method), 93  
`version()` (`sage.interfaces.lie.LiE` method), 101  
`version()` (`sage.interfaces.lisp.Lisp` method), 105  
`version()` (`sage.interfaces.macaulay2.Macaulay2` method), 111  
`version()` (`sage.interfaces.magma.Magma` method), 128  
`version()` (`sage.interfaces.matlab.Matlab` method), 156  
`version()` (`sage.interfaces.maxima_abstract.MaximaAbstract` method), 175  
`version()` (`sage.interfaces.octave.Octave` method), 209  
`version()` (`sage.interfaces.r.R` method), 260  
`version()` (`sage.interfaces.sage0.Sage` method), 268  
`version()` (`sage.interfaces.scilab.Scilab` method), 273  
`version()` (`sage.interfaces.singular.Singular` method), 286

## W

`whos()` (`sage.interfaces.matlab.Matlab` method), 156  
`whos()` (`sage.interfaces.scilab.Scilab` method), 273  
`with_package()` (`sage.interfaces.maple.Maple` method), 142  
`write()` (`sage.interfaces.magma.MagmaGBLogPrettyPrinter` method), 133  
`write()` (`sage.interfaces.singular.SingularGBLogPrettyPrinter` method), 292  
`write_array()` (`sage.interfaces.four_ti_2.FourTi2` method), 30  
`write_matrix()` (`sage.interfaces.four_ti_2.FourTi2` method), 30  
`write_single_row()` (`sage.interfaces.four_ti_2.FourTi2` method), 30

## X

`X()` (`sage.interfaces.qepcad.qepcad_formula_factory` method), 240

## Z

`zsolve()` (`sage.interfaces.four_ti_2.FourTi2` method), 30