Sage Reference Manual: Discrete dynamics

Release 8.2

The Sage Development Team

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CHAPTER

ONE

CELLULAR AUTOMATA

1.1 Soliton Cellular Automata

AUTHORS:

• Travis Scrimshaw (2017-06-30): Initial version

Bases: sage.structure.sage_object.SageObject

Soliton cellular automata.

Fix an affine Lie algebra g with index I and classical index set I_0 . Fix some $r \in I_0$. A soliton cellular automaton (SCA) is a discrete (non-linear) dynamical system given as follows. The states are given by elements of a semi-infinite tensor product of Kirillov-Reshetihkin crystals $B^{r,1}$, where only a finite number of factors are not the maximal element u, which we will call the vacuum. The time evolution T_s is defined by

$$R(p \otimes u_s) = u_s \otimes T_s(p),$$

where $p=\cdots\otimes p_3\otimes p_2\otimes p_1\otimes p_0$ is a state and u_s is the maximal element of $B^{r,s}$. In more detail, we have $R(p_i\otimes u^{(i)})=u^{(i+1)}\otimes \widetilde{p}_i$ with $u^{(0)}=u_s$ and $T_s(p)=\cdots\otimes \widetilde{p}_1\otimes \widetilde{p}_0$. This is well-defined since $R(u\otimes u_s)=u_s\otimes u$ and $u^{(k)}=u_s$ for all $k\gg 1$.

INPUT:

- initial_state the list of elements, can also be a string when vacuum is 1 and n is \mathfrak{sl}_n
- cartan_type (default: 2) the value n, for \mathfrak{sl}_n , or a Cartan type
- r (default: 1) the node index r; typically this corresponds to the height of the vacuum element

EXAMPLES:

We first create an example in \mathfrak{sl}_4 (type A_3):

```
sage: B = SolitonCellularAutomata('3411111122411112223', 4)
sage: B
Soliton cellular automata of type ['A', 3, 1] and vacuum = 1
  initial state:
  34.....224....2223
  evoltuions: []
  current state:
  34.....224....2223
```

We then apply an standard evolution:

```
sage: B.evolve()
sage: B
Soliton cellular automata of type ['A', 3, 1] and vacuum = 1
  initial state:
    34.....224....223
    evoltuions: [(1, 19)]
    current state:
    ..............34.....224...223....
```

Next, we apply a smaller carrier evolution. Note that the soliton of size 4 moves only 3 steps:

```
sage: B.evolve(3)
sage: B
Soliton cellular automata of type ['A', 3, 1] and vacuum = 1
  initial state:
    34.....224....2223
    evoltuions: [(1, 19), (1, 3)]
    current state:
    ...........34....224...2223......
```

We can also use carriers corresponding to non-vacuum indices. In these cases, the carrier might not return to its initial state, which results in a message being displayed about the resulting state of the carrier:

```
sage: B.evolve(carrier_capacity=7, carrier_index=3)
Last carrier:
 1 1 1 1 1 1 1
 2 2 2 2 2 3 3
 3 3 3 3 4 4
Soliton cellular automata of type ['A', 3, 1] and vacuum = 1
 initial state:
 34.....224....2223
 evoltuions: [(1, 19), (1, 3), (3, 7)]
 current state:
 .....23....222....2223......
sage: B.evolve(carrier_capacity=3, carrier_index=2)
Last carrier:
 1 1 1
 2 2 3
sage: B
Soliton cellular automata of type ['A', 3, 1] and vacuum = 1
 initial state:
 34.....224....2223
 evoltuions: [(1, 19), (1, 3), (3, 7), (2, 3)]
 current state:
 .....22.....223....2222......
```

To summarize our current evolutions, we can use print_states():

```
t: 3
.....23...222...2223....
t: 4
.....22...223...222....
```

To run the SCA further under the standard evolutions, one can use print_states() or latex_states():

```
sage: B.print_states(15)
t: 0
t: 1
t: 2
                 ......34....224...2223......
t: 3
                   .....23....222....2223.....
t: 4
                   .....22.....223....2222......
                   ......22....223..2222......
t: 6
                  .....22..2223..222......
t: 7
                .......2222..23...222............
t: 8
             t: 9
            t: 10
            .....2222.......223.22........
t: 11
        t: 12
   t: 13
   t: 14
   .....2222......223.....22......
```

Next, we use r=2 in type A_3 . Here, we give the data as lists of values corresponding to the entries of the column of height 2 from the largest entry to smallest. Our columns are drawn in French convention:

```
sage: B = SolitonCellularAutomata([[4,1],[4,1],[2,1],[2,1],[2,1],[2,1],[3,1],[3,\rightarrow1],[3,2]], 4, 2)
```

We perform 3 evolutions and obtain the following:

```
sage: B.evolve(number=3)
sage: B
Soliton cellular automata of type ['A', 3, 1] and vacuum = 2
  initial state:
    44    333
    11....112
    evoltuions: [(2, 9), (2, 9), (2, 9)]
    current state:
        44    333
...11.112......
```

We construct Example 2.9 from [LS2017]:

```
sage: B = SolitonCellularAutomata([[2],[-3],[1],[1],[1],[4],[0],[-2],
     [1],[1],[1],[1],[3],[-4],[-3],[-3],[1]], ['D',5,2])
sage: B.print_states(10)
t: 0
                              ....23...402....3433.
t: 1
t: 2
                             .23.402..3433.....
t: 3
                          ..243.02.3433......
t: 4
                       .2403..42333......
t: 5
                    .2403...44243..
t: 6
                .2403....442.43..
t: 7
            ..2403.....442..43..
t: 8
    t: 9
    ...2403......442....43........
```

Example 3.4 from [LS2017]:

```
sage: B = SolitonCellularAutomata([['E'],[1],[1],[1],[3],[0],
....: [1],[1],[1],[1],[2],[-3],[-1],[1]], ['D',4,2])
sage: B.print_states(10)
t: 0
t: 1
t: 2
                              .....E303.21.....
t: 3
                             .....303E2.22......
t: 4
                            ....303E...222.....
t: 5
                        ....303E.....12..
t: 6
                      ..303E......1.2.
t: 7
                   ..303E....1..2.
t: 8
              ....303E.....
t: 9
```

Example 3.12 from [LS2017]:

```
. . . . . . . . . 2 . . 1 1
                                  3 2-3
t.: 1
                     . . . . . . 2 1 1 . . .
                             -3-1
                             2-2
t: 2
                  . . . . . . 1-3 . . . . . .
                       -3-1 -3
t: 3
                      2-2 2
                 . . . 1 3 . 1 . . . . . .
                        -3
2
                -3-1
                 2-2
t: 4
                           2
      . . . . . . 1 3 . . . 1 . . . . . . .
          -3-1
                        -3
           2-2
                         2
t: 5
      . . . 1 3 . . . . . 1 . . . . . . . .
```

Example 4.12 from [LS2017]:

```
sage: K = crystals.KirillovReshetikhin(['E',6,1], 1,1, 'KR')
sage: u = K.module_generators[0]
sage: x = u.f_string([1,3,4,5])
sage: y = u.f_string([1,3,4,2,5,6])
sage: a = u.f_string([1,3,4,2])
sage: B = SolitonCellularAutomata([a, u,u,u, x,y], ['E',6,1], 1)
Soliton cellular automata of type ['E', 6, 1] and vacuum = 1
 initial state:
                                           (-5, 2, 6)(-2, -6, 4)
     (-2, 5)
 evoltuions: []
 current state:
     (-2, 5)
                                            (-5, 2, 6)(-2, -6, 4)
sage: B.print_states(8)
t: 0 ...
t: 7
                          (-2, 5)(-2, -5, 4, 6) \dots (-6, 2) \dots
```

evolve(carrier_capacity=None, carrier_index=None, number=None)

Evolve self.

Time evolution T_s of a SCA state p is determined by

$$u_{r,s} \otimes T_s(p) = R(p \otimes u_{r,s}),$$

where $u_{r,s}$ is the maximal element of $B^{r,s}$.

INPUT:

- carrier_capacity (default: the number of balls in the system) the size s of carrier
- carrier_index (default: the vacuum index) the index r of the carrier
- number (optional) the number of times to perform the evolutions

To perform multiple evolutions of the SCA, carrier_capacity and carrier_index may be lists of the same length.

```
sage: B = SolitonCellularAutomata('3411111122411112223', 4)
sage: for k in range(10):
       B.evolve()
. . . . :
sage: B
Soliton cellular automata of type ['A', 3, 1] and vacuum = 1
 initial state:
 34.....224....2223
 evoltuions: [(1, 19), (1, 19), (1, 19), (1, 19), (1, 19),
             (1, 19), (1, 19), (1, 19), (1, 19), (1, 19)]
 current state:
 sage: B.reset()
sage: B.evolve(number=10); B
Soliton cellular automata of type ['A', 3, 1] and vacuum = 1
 initial state:
 34.....224....2223
 evoltuions: [(1, 19), (1, 19), (1, 19), (1, 19), (1, 19),
             (1, 19), (1, 19), (1, 19), (1, 19), (1, 19)]
 current state:
 sage: B.reset()
sage: B.evolve(carrier_capacity=[1,2,3,4,5,6,7,8,9,10]); B
Soliton cellular automata of type ['A', 3, 1] and vacuum = 1
 initial state:
 34.....224....2223
 evoltuions: [(1, 1), (1, 2), (1, 3), (1, 4), (1, 5),
            (1, 6), (1, 7), (1, 8), (1, 9), (1, 10)]
 current state:
 ......2344....222..23.......
sage: B.reset()
sage: B.evolve(carrier_index=[1,2,3])
Last carrier:
 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1
 2 2 2 2 2 2 2 2 2 2 2 2 2 2 3 4
sage: B
Soliton cellular automata of type ['A', 3, 1] and vacuum = 1
 initial state:
 34.....224....2223
 evoltuions: [(1, 19), (2, 19), (3, 19)]
 current state:
 ......22.....223...2222.....
sage: B.reset()
sage: B.evolve(carrier_capacity=[1,2,3], carrier_index=[1,2,3])
Last carrier:
 1 1
 3 4
Last carrier:
 1 1 1
 2 2 3
 3 3
sage: B
Soliton cellular automata of type ['A', 3, 1] and vacuum = 1
 initial state:
 34.....224....2223
```

```
evoltuions: [(1, 1), (2, 2), (3, 3)]
 current state:
  .....22......223.....2222..
sage: B.reset()
sage: B.evolve(1, 2, number=3)
Last carrier:
 3
Last carrier:
 1
 4
Last carrier:
 3
sage: B
Soliton cellular automata of type ['A', 3, 1] and vacuum = 1
 initial state:
 34.....224....2223
 evoltuions: [(2, 1), (2, 1), (2, 1)]
 current state:
  .24.....222.....2222.
```

latex_state_evolution (num, scale=1)

Return a latex version of the evolution process of the state num.

See also:

state_evolution(), print_state_evolution()

```
sage: B = SolitonCellularAutomata('113123', 3)
sage: B.evolve(3)
sage: B.latex_state_evolution(0)
\begin{tikzpicture}[scale=1]
\node (i0) at (0,0.9) {$1$};
\node (i1) at (-2,0.9) {$1$};
\node (i2) at (-4,0.9) {$3$};
\node (i3) at (-6,0.9) {$1$};
\node (i4) at (-8,0.9) {$2$};
\node (i5) at (-10,0.9) {$3$};
\node (t0) at (0,-1) {$1$};
\node (t1) at (-2,-1) {$3$};
\node (t2) at (-4,-1) {$2$};
\node (t3) at (-6, -1) {$3$};
\node (t4) at (-8, -1) {$1$};
\node (t5) at (-10,-1) {$1$};
\node (u0) at (1,0) {$111$};
\node (u1) at (-1,0) {$111$};
\node (u2) at (-3,0) {$113$};
\node (u3) at (-5,0) {$112$};
\node (u4) at (-7,0) {$123$};
\node (u5) at (-9,0) {$113$};
\node (u6) at (-11,0) {$111$};
\draw[->] (i0) -- (t0);
\draw[->] (u0) -- (u1);
\draw[->] (i1) -- (t1);
\draw[->] (u1) -- (u2);
```

```
\draw[->] (i2) -- (t2);
\draw[->] (u2) -- (u3);
\draw[->] (i3) -- (t3);
\draw[->] (u3) -- (u4);
\draw[->] (i4) -- (t4);
\draw[->] (u4) -- (u5);
\draw[->] (i5) -- (t5);
\draw[->] (u5) -- (u6);
\end{tikzpicture}

sage: B.latex_state_evolution(1)
\begin{tikzpicture} [scale=1]
...
\end{tikzpicture}
```

latex_states (num=None, as_array=True, box_width='5pt')

Return a latex verion of the states.

INPUT:

- num the number of states
- as_array (default: True) if True, then the states are placed inside of an array; if False, then the states are given as a word
- box_width (default: '5pt') the width of the . used to represent the vacuum state when as_array is True

If as_array is False, then the vacuum element is printed in a gray color. If as_array is True, then the vacuum is given as .

Use the box_width to help create more even spacing when a column in the output contains only vacuum elements.

```
sage: B = SolitonCellularAutomata('411122', 4)
sage: B.latex_states(8)
t = 0 & \cdots & ... & \makebox[5pt]{.} & 4 & \makebox[5pt]{.}
& \makebox[5pt]{.} & \makebox[5pt]{.} & 2 & 2 \\
t = 1 \& \cdots \& ... \& 4 \& \makebox[5pt]{.} \& \makebox[5pt]{.} & \& 2 \& 2 & ... 
t = 2 \& \cdots \& ... \& 4 \& \makebox[5pt]{.} & 2 & 2 & ... \
t = 3 & \cdots & ... & 4 & 2 & 2 & ... \\
t = 4 & \cdots & ... & 2 & 4 & 2 & ... \\
t = 5 & \cdots & ... & 2 & 4 & \makebox[5pt]{.} & 2 & ... \\
t = 6 \& \cdots \& ... \& 2 \& 4 \& \makebox[5pt]{.} \& \makebox[5pt]{.}
& 2 & ... \\
t = 7 & \cdots & \makebox[5pt]{.} & 2 & 4 & \makebox[5pt]{.}
& \makebox[5pt]{.} & \makebox[5pt]{.} & 2 & ... \\
\end{array}}
sage: B = SolitonCellularAutomata('511122', 5)
sage: B.latex_states(8, as_array=False)
{\begin{array}{c|c}
{\color{gray} 1} {\color{gray} 1} 2 2 \\
t = 1 \& \cdot cdots ... 5 {\cdot color{gray} 1} {\cdot color{gray} 1} 2 2 ... \
t = 2 & \cdots ... 5 {\color{gray} 1} 2 2 ... \\
t = 3 & \cdots ... 5 2 2 ... \\
```

```
t = 4 & \cdots ... 2 5 2 ... \\
t = 5 & \cdots ... 2 5 {\color{gray} 1} 2 ... \\
t = 6 & \cdots ... 2 5 {\color{gray} 1} {\color{gray} 1} 2 ... \\
t = 7 & \cdots {\color{gray} 1} 2 5 {\color{gray} 1}
{\color{gray} 1} {\color{gray} 1} 2 ... \\
\end{\array}}
```

print_state (num=None, vacuum_letter='.', remove_trailing_vacuums=False)

Print the state num.

INPUT:

- num (default: the current state) the state to print
- vacuum_letter (default: '.') the letter to print for the vacuum
- remove_trailing_vacuums (default: False) if True then this does not print the vacuum letters at the right end of the state

EXAMPLES:

print_state_evolution(num)

Print the evolution process of the state num.

See also:

state_evolution(), latex_state_evolution()

EXAMPLES:

```
sage: B = SolitonCellularAutomata('1113123', 3)
sage: B.evolve(3)
sage: B.evolve(3)
sage: B.print_state_evolution(0)
  1 1 1
                     3 1 2
        111 --+-- 111 --+-- 111 --+-- 113 --+-- 112 --+-- 123 --+-- 113 --+-- 111
  3
      1
                           3
                                 1
                     2
                                       1
sage: B.print_state_evolution(1)
   1 1 3
                    2
                           3
                                 1
               111 --+-- 113 --+-- 133 --+-- 123 --+-- 113 --+-- 111 --+-- 111
                                 3
         3
            2
                     1 1
                               1
                                       1
```

print states (num=None, vacuum letter='.')

Print the first num states of self.

Note: If the number of states computed for self is less than num, then this evolves the system using the default time evolution.

INPUT:

• num – the number of states to print

```
sage: B = SolitonCellularAutomata([[2],[-1],[1],[1],[1],[1],[2],[2],[3],
     [-2],[1],[1],[2],[-1],[1],[1],[1],[1],[1],[2],[3],[3],[-3],[-2]],
     ['C',3,1])
. . . . :
sage: B.print states(7)
t: 0
    .....21....2232..21.....23332
t: 1
        .....21...2232...21....23332.....
         .....21..2232....21..23332.....
t: 3
        t: 4
        ......221...232.2231....332......
    ......221...2321223......332.......
t: 6
    ..2221...321..223.....332......
sage: B = SolitonCellularAutomata([[2],[1],[1],[1],[3],[-2],[1],[1],
....: [1],[2],[2],[-3],[1],[1],[1],[1],[1],[1],[2],[3],[3],[-3]],
....: ['B',3,1])
sage: B.print_states(9, ' ')
t: 0
                      2 32 223
                                      2333
t: 1
                      2 32 223
                                   2333
t: 2
                     2 32 223
                               2333
t: 3
                   23 2223
                           2333
t.: 4
                  23 213 2333
t: 5
              2233 222 333
t: 6
           2233 23223 3
t: 7
       2233
             232 23 3
t: 8
           232 23
    2233
sage: B = SolitonCellularAutomata([[2],[-2],[1],[1],[1],[1],[2],[0],[-3],
...: [1],[1],[1],[1],[2],[2],[3],[-3],], ['D',4,2])
sage: B.print_states(10)
t: 0
       .....22....203.....2233
t: 1
```

```
.....22...203....2233....
t: 2
     t: 3
         .....22.203..2233.....
t: 4
               .....22203.2233..........
t: 5
            ......220223.233.....
t: 6
      t: 7
        .....2202..223..33......
      .....2202...223...33.....
t: 9
   ......2202....223....33.......
```

Example 4.13 from [Yamada2007]:

```
sage: B = SolitonCellularAutomata([[3],[3],[1],[1],[1],[1],[2],[2],[2]], ['D',
\hookrightarrow4,3])
sage: B.print_states(15)
t: 0
t: 1
t: 2
              t: 3
t: 5
               t: 6
                ....332......
t: 8
t: 9
t: 10
          ....20E.....
t: 11
        ....233......
t: 12
    .......2302......
t: 13
    .....23322.....
t: 14
    ..233.22.....
```

Example 4.14 from [Yamada2007]:

```
sage: B.print_states(15)
t: 0
 t: 1
  t: 2
  .....3.23.233.....
t: 3
   .....323233......
t: 4
  t: 5
 t: 6
    ......30E.3.......
t: 7
  t: 8
  t: 9
  t: 10
  t: 11
 t: 12
 t: 13
 t: 14
```

reset()

Reset self back to the initial state.

```
sage: B = SolitonCellularAutomata('34111111224', 4)
Soliton cellular automata of type ['A', 3, 1] and vacuum = 1
 initial state:
 34....224
 evoltuions: []
 current state:
 34....224
sage: B.evolve()
sage: B.evolve()
sage: B.evolve()
sage: B.evolve()
sage: B
Soliton cellular automata of type ['A', 3, 1] and vacuum = 1
 initial state:
 evoltuions: [(1, 11), (1, 11), (1, 11), (1, 11)]
 current state:
  ...34..224.....
sage: B.reset()
sage: B
```

```
Soliton cellular automata of type ['A', 3, 1] and vacuum = 1
  initial state:
  34.....224
  evoltuions: []
  current state:
  34.....224
```

state evolution(num)

Return a list of the carrier values at state num evolving to the next state.

If num is greater than the number of states, this performs the standard evolution T_k , where k is the number of balls in the system.

See also:

```
print_state_evolution(), latex_state_evolution()
```

```
sage: B = SolitonCellularAutomata('1113123', 3)
sage: B.evolve(3)
sage: B.state_evolution(0)
[[[1, 1, 1]],
[[1, 1, 1]],
[[1, 1, 1]],
[[1, 1, 3]],
[[1, 1, 2]],
[[1, 2, 3]],
[[1, 1, 3]],
[[1, 1, 1]]]
sage: B.state_evolution(2)
[[[1, 1, 1, 1, 1, 1, 1]],
 [[1, 1, 1, 1, 1, 1, 1]],
 [[1, 1, 1, 1, 1, 1, 1]],
 [[1, 1, 1, 1, 1, 1, 1]],
 [[1, 1, 1, 1, 1, 1, 1]],
 [[1, 1, 1, 1, 1, 1, 1]],
 [[1, 1, 1, 1, 1, 1, 3]],
 [[1, 1, 1, 1, 1, 3, 3]],
 [[1, 1, 1, 1, 1, 1, 3]],
 [[1, 1, 1, 1, 1, 1, 2]],
 [[1, 1, 1, 1, 1, 1, 1]],
 [[1, 1, 1, 1, 1, 1, 1]],
 [[1, 1, 1, 1, 1, 1, 1]],
 [[1, 1, 1, 1, 1, 1, 1]],
 [[1, 1, 1, 1, 1, 1, 1]]]
```

PLOTTING OF MANDELBROT AND JULIA SETS

2.1 Mandelbrot and Julia sets

Plots the Mandelbrot and Julia sets for the map $Q_c(z) = z^2 + c$ in the complex plane.

The Mandelbrot set is the set of complex numbers c for which the function $Q_c(z) = z^2 + c$ does not diverge when iterated from z = 0. This set of complex numbers can be visualized by plotting each value for c in the complex plane. The Mandelbrot set is an example of a fractal when plotted in the complex plane.

The Julia set for a given c is the set of complex numbers for which the function $Q_c(z) = z^2 + c$ is bounded under iteration.

AUTHORS:

· Ben Barros

sage.dynamics.complex_dynamics.mandel_julia.external_ray(theta, **kwds)

Draws the external ray(s) of a given angle (or list of angles) by connecting a finite number of points that were approximated using Newton's method. The algorithm used is described in a paper by Tomoki Kawahira.

REFERENCE:

[Kaw2009]

INPUT:

• theta - double or list of doubles, angles between 0 and 1 inclusive.

kwds:

- image 24-bit RGB image (optional default: None) user specified image of Mandelbrot set.
- D long (optional default: 25) depth of the approximation. As D increases, the external ray gets closer to the boundary of the Mandelbrot set. If the ray doesn't reach the boundary of the Mandelbrot set, increase D.
- S long (optional default: 10) sharpness of the approximation. Adjusts the number of points used to approximate the external ray (number of points is equal to S*D). If ray looks jagged, increase S.
- R long (optional default: 100) radial parameter. If R is large, the external ray reaches sufficiently close to infinity. If R is too small, Newton's method may not converge to the correct ray.
- prec long (optional default: 300) specifies the bits of precision used by the Complex Field when using Newton's method to compute points on the external ray.
- ray_color RGB color (optional default: [255, 255, 255]) color of the external ray(s).

OUTPUT:

24-bit RGB image of external ray(s) on the Mandelbrot set.

EXAMPLES:

```
sage: external_ray(1/3)
500x500px 24-bit RGB image
```

```
sage: external_ray(0.6, ray_color=[255, 0, 0])
500x500px 24-bit RGB image
```

```
sage: external_ray([0, 0.2, 0.4, 0.7]) # long time
500x500px 24-bit RGB image
```

```
sage: external_ray([i/5 for i in range(1,5)]) # long time
500x500px 24-bit RGB image
```

WARNING:

If you are passing in an image, make sure you specify which parameters to use when drawing the external ray. For example, the following is incorrect:

```
sage: M = mandelbrot_plot(x_center=0) # not tested
sage: external_ray(5/7, image=M) # not tested
500x500px 24-bit RGB image
```

To get the correct external ray, we adjust our parameters:

```
sage: M = mandelbrot_plot(x_center=0) # not tested
sage: external_ray(5/7, x_center=0, image=M) # not tested
500x500px 24-bit RGB image
```

Todo: The copy () function for bitmap images needs to be implemented in Sage.

```
sage.dynamics.complex_dynamics.mandel_julia.julia_plot (c=-1, **kwds)
```

Plots the Julia set of a given complex c value. Users can specify whether they would like to display the Mandelbrot side by side with the Julia set.

The Julia set of a given c value is the set of complex numbers for which the function $Q_c(z) = z^2 + c$ is bounded under iteration. The Julia set can be visualized by plotting each point in the set in the complex plane. Julia sets are examples of fractals when plotted in the complex plane.

ALGORITHM:

Define the map $Q_c(z)=z^2+c$ for some $c\in\mathbb{C}$. For every $p\in\mathbb{C}$, if $|Q_c^k(p)|>2$ for some $k\geq 0$, then $Q_c^n(p)\to\infty$. Let N be the maximum number of iterations. Compute the first N points on the orbit of p under Q_c . If for any k< N, $|Q_c^k(p)|>2$, we stop the iteration and assign a color to the point p based on how quickly p escaped to infinity under iteration of Q_c . If $|Q_c^i(p)|\leq 2$ for all $i\leq N$, we assume p is in the Julia set and assign the point p the color black.

INPUT:

• c – complex (optional - default: -1), complex point c that determines the Julia set.

kwds:

- period list (optional default: None), returns the Julia set for a random c value with the given (formal) cycle structure.
- mandelbrot boolean (optional default: True), when set to True, an image of the Mandelbrot set is appended to the right of the Julia set.

- point_color RGB color (optional default: [255, 0, 0]), color of the point c in the Mandelbrot set.
- x_center double (optional default: -1.0), Real part of center point.
- y_center double (optional default: 0.0), Imaginary part of center point.
- image_width double (optional default: 4.0), width of image in the complex plane.
- max iteration long (optional default: 500), maximum number of iterations the map $Q_c(z)$.
- pixel_count long (optional default: 500), side length of image in number of pixels.
- base_color RGB color (optional default: [40, 40, 40]), color used to determine the coloring of set.
- iteration_level long (optional default: 1), number of iterations between each color level.
- number_of_colors long (optional default: 30), number of colors used to plot image.
- interact boolean (optional default: False), controls whether plot will have interactive functionality.

OUTPUT:

24-bit RGB image of the Julia set in the complex plane.

EXAMPLES:

```
sage: julia_plot()
1001x500px 24-bit RGB image
```

To display only the Julia set, set mandelbrot to False:

```
sage: julia_plot(mandelbrot=False)
500x500px 24-bit RGB image
```

To display an interactive plot of the Julia set in the Notebook, set interact to True:

```
sage: julia_plot(interact=True)
<html>...</html>
```

To return the Julia set of a random c value with (formal) cycle structure (2,3), set period = [2,3]:

```
sage: julia_plot(period=[2,3])
1001x500px 24-bit RGB image
```

To return all of the Julia sets of c values with (formal) cycle structure (2,3):

```
sage: period = [2,3] # not tested
...: R.<c> = QQ[]
...: P.<x,y> = ProjectiveSpace(R,1)
...: f = DynamicalSystem([x^2+c*y^2, y^2])
...: L = f.dynatomic_polynomial(period).subs({x:0,y:1}).roots(ring=CC)
...: c_values = [k[0] for k in L]
...: for c in c_values:
...: julia_plot(c)
```

sage.dynamics.complex_dynamics.mandel_julia.mandelbrot_plot (**kwds)
Interactive plot of the Mandelbrot set for the map $Q_c(z) = z^2 + c$.

ALGORITHM:

Let each pixel in the image be a point $c \in \mathbb{C}$ and define the map $Q_c(z) = z^2 + c$. If $|Q_c^k(c)| > 2$ for some $k \geq 0$, it follows that $Q_c^n(c) \to \infty$. Let N be the maximum number of iterations. Compute the first N points on the orbit of 0 under Q_c . If for any k < N, $|Q_c^k(0)| > 2$, we stop the iteration and assign a color to the point c based on how quickly 0 escaped to infinity under iteration of Q_c . If $|Q_c^i(0)| \leq 2$ for all $i \leq N$, we assume c is in the Mandelbrot set and assign the point c the color black.

REFERENCE:

[Dev2005]

kwds:

- x_center double (optional default: -1.0), Real part of center point.
- y_center double (optional default: 0.0), Imaginary part of center point.
- image_width double (optional default: 4.0), width of image in the complex plane.
- max_iteration long (optional default: 500), maximum number of iterations the map Q_c(z).
- pixel_count long (optional default: 500), side length of image in number of pixels.
- base_color RGB color (optional default: [40, 40, 40]) color used to determine the coloring of set.
- iteration_level long (optional default: 1) number of iterations between each color level.
- number_of_colors long (optional default: 30) number of colors used to plot image.
- interact boolean (optional default: False), controls whether plot will have interactive functionality.

OUTPUT:

24-bit RGB image of the Mandelbrot set in the complex plane.

EXAMPLES:

```
sage: mandelbrot_plot() # long time
500x500px 24-bit RGB image
```

```
sage: mandelbrot_plot(pixel_count=1000) # long time
1000x1000px 24-bit RGB image
```

To display an interactive plot of the Mandelbrot set in the Notebook, set interact to True:

```
sage: mandelbrot_plot(interact=True)
<html>...</html>
```

```
sage: mandelbrot_plot(interact=True, x_center=-0.75, y_center=0.25,
....: image_width=1/2, number_of_colors=75)
<html>...</html>
```

CHAPTER

THREE

ABELIAN DIFFERENTIALS AND FLAT SURFACES

3.1 Strata of differentials on Riemann surfaces

Warning: This module is deprecated. You are advised to install and use the surface_dynamics package instead available at https://pypi.python.org/pypi/surface_dynamics/

The space of Abelian (or quadratic) differentials is stratified by the degrees of the zeroes (and simple poles for quadratic differentials). Each stratum has one, two or three connected components and each is associated to an (extended) Rauzy class. The <code>connected_components()</code> method (only available for Abelian stratum) give the decomposition of a stratum (which corresponds to the SAGE object <code>AbelianStratum</code>).

The work for Abelian differentials was done by Maxim Kontsevich and Anton Zorich in [KZ2003] and for quadratic differentials by Erwan Lanneau in [Lan2008]. Zorich gave an algorithm to pass from a connected component of a stratum to the associated Rauzy class (for both interval exchange transformations and linear involutions) in [Zor2008] and is implemented for Abelian stratum at different level (approximately one for each component):

- for connected stratum representative ()
- for hyperelliptic component representative ()
- for non hyperelliptic component, the algorithm is the same as for connected component
- for odd component representative ()
- for even component representative ()

The inverse operation (pass from an interval exchange transformation to the connected component) is partially written in [KZ2003] and simply named here <code>connected_component()</code>.

All the code here was first available on Mathematica [Zor].

Note: The quadratic strata are not yet implemented.

AUTHORS:

• Vincent Delecroix (2009-09-29): initial version

EXAMPLES:

Construction of a stratum from a list of singularity degrees:

```
sage: a = AbelianStratum(1,1)
doctest:warning
...
```

```
DeprecationWarning: AbelianStratum is deprecated and will be removed from Sage. You are advised to install the surface_dynamics package via:
sage -pip install surface_dynamics
If you do not have write access to the Sage installation you can alternatively do
sage -pip install surface_dynamics --user
The package surface_dynamics subsumes all flat surface related computation that are currently available in Sage. See more information at
http://www.labri.fr/perso/vdelecro/surface-dynamics/latest/
See http://trac.sagemath.org/20695 for details.
sage: a
H(1, 1)
sage: a.genus()
2
sage: a.nintervals()
5
```

```
sage: a = AbelianStratum(4,3,2,1)
sage: a
H(4, 3, 2, 1)
sage: a.genus()
6
sage: a.nintervals()
15
```

By convention, the degrees are always written in decreasing order:

```
sage: a1 = AbelianStratum(4,3,2,1)
sage: a1
H(4, 3, 2, 1)
sage: a2 = AbelianStratum(2,3,1,4)
sage: a2
H(4, 3, 2, 1)
sage: a1 == a2
True
```

It is also possible to consider stratum with an incoming or an outgoing separatrix marked (the aim of this consideration is to attach a specified degree at the left or the right of the associated interval exchange transformation):

```
sage: a_out = AbelianStratum(1, 1, marked_separatrix='out')
sage: a_out
H^out(1, 1)
sage: a_in = AbelianStratum(1, 1, marked_separatrix='in')
sage: a_in
H^in(1, 1)
sage: a_out == a_in
False
```

Get a list of strata with constraints on genus or on the number of intervals of a representative:

```
sage: for a in AbelianStrata(genus=3):
....: print(a)
doctest:warning
...
DeprecationWarning: AbelianStrata is deprecated and will be removed from Sage.
You are advised to install the surface_dynamics package via:
```

```
sage -pip install surface_dynamics
If you do not have write access to the Sage installation you can
alternatively do
    sage -pip install surface_dynamics --user
The package surface_dynamics subsumes all flat surface related
computation that are currently available in Sage. See more
information at
    http://www.labri.fr/perso/vdelecro/surface-dynamics/latest/
See http://trac.sagemath.org/20695 for details.
H(4)
H(3, 1)
H(2, 2)
H(2, 1, 1)
H(1, 1, 1, 1)
```

```
sage: for a in AbelianStrata(nintervals=5):
....:     print(a)
H^out(0, 2)
H^out(2, 0)
H^out(1, 1)
H^out(0, 0, 0, 0)
```

```
sage: for a in AbelianStrata(genus=2, nintervals=5):
...: print(a)
H^out(0, 2)
H^out(2, 0)
H^out(1, 1)
```

Obtains the connected components of a stratum:

```
sage: a = AbelianStratum(0)
sage: a.connected_components()
[H_hyp(0)]
```

```
sage: a = AbelianStratum(6)
sage: cc = a.connected_components()
sage: cc
[H_hyp(6), H_odd(6), H_even(6)]
sage: for c in cc:
         print(c)
         print(c.representative(alphabet=range(1,9)))
H_hyp(6)
1 2 3 4 5 6 7 8
8 7 6 5 4 3 2 1
H odd(6)
1 2 3 4 5 6 7 8
4 3 6 5 8 7 2 1
H_even(6)
1 2 3 4 5 6 7 8
6 5 4 3 8 7 2 1
```

```
sage: a = AbelianStratum(1, 1, 1, 1)
sage: a.connected_components()
[H_c(1, 1, 1, 1)]
sage: c = a.connected_components()[0]
sage: print(c.representative(alphabet="abcdefghi"))
```

```
abcdefghi
edcfihgba
```

The zero attached on the left of the associated Abelian permutation corresponds to the first singularity degree:

```
sage: a = AbelianStratum(4, 2, marked_separatrix='out')
sage: b = AbelianStratum(2, 4, marked_separatrix='out')
sage: a == b
False
sage: a, a.connected_components()
(H^out(4, 2), [H_odd^out(4, 2), H_even^out(4, 2)])
sage: b, b.connected_components()
(H^out(2, 4), [H_odd^out(2, 4), H_even^out(2, 4)])
sage: a_odd, a_even = a.connected_components()
sage: b_odd, b_even = b.connected_components()
```

The representatives are hence different:

```
sage: a_odd.representative(alphabet=range(1,10))
1 2 3 4 5 6 7 8 9
4 3 6 5 7 9 8 2 1
sage: b_odd.representative(alphabet=range(1,10))
1 2 3 4 5 6 7 8 9
4 3 5 7 6 9 8 2 1
```

```
sage: a_even.representative(alphabet=range(1,10))
1 2 3 4 5 6 7 8 9
6 5 4 3 7 9 8 2 1
sage: b_even.representative(alphabet=range(1,10))
1 2 3 4 5 6 7 8 9
7 6 5 4 3 9 8 2 1
```

You can retrieve the decomposition of the irreducible Abelian permutations into Rauzy diagrams from the classification of strata:

```
sage: a = AbelianStrata(nintervals=4)
sage: l = sum([stratum.connected_components() for stratum in a], [])
sage: n = [x.rauzy_diagram().cardinality() for x in l]
sage: for c,i in zip(l,n):
...: print("{}: {}".format(c, i))
H_hyp^out(2): 7
H_hyp^out(0, 0, 0): 6
sage: sum(n)
13
```

```
sage: a = AbelianStrata(nintervals=5)
sage: l = sum([stratum.connected_components() for stratum in a], [])
sage: n = [x.rauzy_diagram().cardinality() for x in l]
sage: for c,i in zip(l,n):
....: print("{} : {}".format(c, i))
H_hyp^out(0, 2) : l1
H_hyp^out(2, 0) : 35
H_hyp^out(1, 1) : 15
H_hyp^out(0, 0, 0, 0) : 10
sage: sum(n)
71
```

```
sage: a = AbelianStrata(nintervals=6)
sage: l = sum([stratum.connected_components() for stratum in a], [])
sage: n = [x.rauzy_diagram().cardinality() for x in l]
sage: for c,i in zip(l,n):
....:    print("{} : {}".format(c, i))
H_hyp^out(4) : 31
H_odd^out(4) : 134
H_hyp^out(0, 2, 0) : 66
H_hyp^out(2, 0, 0) : 105
H_hyp^out(2, 0, 0) : 105
H_hyp^out(0, 1, 1) : 20
H_hyp^out(1, 1, 0) : 90
H_hyp^out(0, 0, 0, 0, 0) : 15
sage: sum(n)
461
```

Abelian strata.

INPUT:

- genus a non negative integer or None
- nintervals a non negative integer or None
- marked_separatrix 'no' (for no marking), 'in' (for marking an incoming separatrix) or 'out' (for marking an outgoing separatrix)

EXAMPLES:

Abelian strata with a given genus:

```
sage: for s in AbelianStrata(genus=1): print(s)
H(0)
```

```
sage: for s in AbelianStrata(genus=2): print(s)
H(2)
H(1, 1)
```

```
sage: for s in AbelianStrata(genus=3): print(s)
H(4)
H(3, 1)
H(2, 2)
H(2, 1, 1)
H(1, 1, 1, 1)
```

```
sage: for s in AbelianStrata(genus=4): print(s)
H(6)
H(5, 1)
H(4, 2)
H(4, 1, 1)
H(3, 3)
H(3, 2, 1)
H(3, 1, 1, 1)
H(2, 2, 2)
H(2, 2, 2, 1, 1)
H(2, 1, 1, 1, 1, 1)
H(1, 1, 1, 1, 1, 1, 1)
```

Abelian strata with a given number of intervals:

```
sage: for s in AbelianStrata(nintervals=2): print(s)
H^out(0)
```

```
sage: for s in AbelianStrata(nintervals=3): print(s)
H^out(0, 0)
```

```
sage: for s in AbelianStrata(nintervals=4): print(s)
H^out(2)
H^out(0, 0, 0)
```

```
sage: for s in AbelianStrata(nintervals=5): print(s)
H^out(0, 2)
H^out(2, 0)
H^out(1, 1)
H^out(0, 0, 0, 0)
```

Abelian strata with both constraints:

```
sage: for s in AbelianStrata(genus=2, nintervals=4): print(s)
H^out(2)
```

```
sage: for s in AbelianStrata(genus=5, nintervals=12): print(s)
H^out(8, 0, 0)
H^out(0, 8, 0)
H^{out}(0, 7, 1)
H^{out}(1, 7, 0)
H^out(7, 1, 0)
H^{out}(0, 6, 2)
H^{out}(2, 6, 0)
H^out(6, 2, 0)
H^{out}(1, 6, 1)
H^out(6, 1, 1)
H^{out}(0, 5, 3)
H^{out}(3, 5, 0)
H^out(5, 3, 0)
H^{out}(1, 5, 2)
H^{out}(2, 5, 1)
H^{out}(5, 2, 1)
H^out(0, 4, 4)
H^out(4, 4, 0)
H^{out}(1, 4, 3)
H^{out}(3, 4, 1)
H^out(4, 3, 1)
H^out(2, 4, 2)
H^{out}(4, 2, 2)
H^{out}(2, 3, 3)
H^{out}(3, 3, 2)
```

```
class sage.dynamics.flat_surfaces.strata.AbelianStrata_all(category=None)
    Bases: sage.combinat.combinat.InfiniteAbstractCombinatorialClass
```

Abelian strata.

Strata with constraint number of intervals.

INPUT:

- nintervals an integer greater than 1
- marked_separatrix 'no', 'out' or 'in'

```
class sage.dynamics.flat_surfaces.strata.AbelianStrata_g(genus=None,
```

marked_separatrix=None)

Bases: sage.combinat.combinat.CombinatorialClass

Stratas of genus g surfaces.

INPUT:

- genus a non negative integer
- marked separatrix 'no', 'out' or 'in'

Bases: sage.combinat.combinat.CombinatorialClass

Abelian strata of prescribed genus and number of intervals.

INPUT:

- genus integer: the genus of the surfaces
- nintervals integer: the number of intervals
- marked_separatrix 'no', 'in' or 'out'

```
class sage.dynamics.flat_surfaces.strata.AbelianStratum(*l, **d)
Bases: sage.structure.sage_object.SageObject
```

Stratum of Abelian differentials.

A stratum with a marked outgoing separatrix corresponds to Rauzy diagram with left induction, a stratum with marked incoming separatrix correspond to Rauzy diagram with right induction. If there is no marked separatrix, the associated Rauzy diagram is the extended Rauzy diagram (consideration of the <code>sage.dynamics.interval_exchanges.template.Permutation.symmetric()</code> operation of Boissy-Lanneau).

When you want to specify a marked separatrix, the degree on which it is the first term of your degrees list.

INPUT:

• marked_separatrix - None (default) or 'in' (for incoming separatrix) or 'out' (for outgoing separatrix).

EXAMPLES:

Creation of an Abelian stratum and get its connected components:

```
sage: a = AbelianStratum(2, 2)
sage: a
H(2, 2)
sage: a.connected_components()
[H_hyp(2, 2), H_odd(2, 2)]
```

Specification of marked separatrix:

```
sage: a = AbelianStratum(4,2,marked_separatrix='in')
sage: a
H^in(4, 2)
sage: b = AbelianStratum(2,4,marked_separatrix='in')
sage: b
H^in(2, 4)
sage: a == b
False
```

```
sage: a = AbelianStratum(4,2,marked_separatrix='out')
sage: a
H^out(4, 2)
sage: b = AbelianStratum(2,4,marked_separatrix='out')
sage: b
H^out(2, 4)
sage: a == b
False
```

Get a representative of a connected component:

```
sage: a = AbelianStratum(2,2)
sage: a_hyp, a_odd = a.connected_components()
sage: a_hyp.representative()
1 2 3 4 5 6 7
7 6 5 4 3 2 1
sage: a_odd.representative()
0 1 2 3 4 5 6
3 2 4 6 5 1 0
```

You can choose the alphabet:

```
sage: a_odd.representative(alphabet="ABCDEFGHIJKLMNOPQRSTUVWXYZ")
A B C D E F G
D C E G F B A
```

By default, you get a reduced permutation, but you can specify that you want a labelled one:

```
sage: p_reduced = a_odd.representative()
sage: p_labelled = a_odd.representative(reduced=False)
```

connected_components()

Lists the connected components of the Stratum.

OUTPUT:

list – a list of connected components of stratum

```
sage: AbelianStratum(0).connected_components()
[H_hyp(0)]
```

```
sage: AbelianStratum(2).connected_components()
[H_hyp(2)]
sage: AbelianStratum(1,1).connected_components()
[H_hyp(1, 1)]
```

```
sage: AbelianStratum(4).connected_components()
[H_hyp(4), H_odd(4)]
sage: AbelianStratum(3,1).connected_components()
[H_c(3, 1)]
sage: AbelianStratum(2,2).connected_components()
[H_hyp(2, 2), H_odd(2, 2)]
sage: AbelianStratum(2,1,1).connected_components()
[H_c(2, 1, 1)]
sage: AbelianStratum(1,1,1,1).connected_components()
[H_c(1, 1, 1, 1)]
```

genus()

Returns the genus of the stratum.

OUTPUT:

integer - the genus

EXAMPLES:

```
sage: AbelianStratum(0).genus()
1
sage: AbelianStratum(1,1).genus()
2
sage: AbelianStratum(3,2,1).genus()
4
```

is connected()

Tests if the strata is connected.

OUTPUT:

boolean - True if it is connected else False

EXAMPLES:

```
sage: AbelianStratum(2).is_connected()
True
sage: AbelianStratum(2).connected_components()
[H_hyp(2)]
```

```
sage: AbelianStratum(2,2).is_connected()
False
sage: AbelianStratum(2,2).connected_components()
[H_hyp(2, 2), H_odd(2, 2)]
```

nintervals()

Returns the number of intervals of any iet of the strata.

OUTPUT

integer - the number of intervals for any associated iet

```
sage: AbelianStratum(0).nintervals()
2
sage: AbelianStratum(0,0).nintervals()
3
sage: AbelianStratum(2).nintervals()
```

```
4
sage: AbelianStratum(1,1).nintervals()
5
```

sage.dynamics.flat_surfaces.strata.CCA
alias of ConnectedComponentOfAbelianStratum

Connected component of Abelian stratum.

Warning: Internal class! Do not use directly!

```
sage: a = AbelianStratum(2,4,0,marked_separatrix='out')
sage: a_odd, a_even = a.connected_components()
sage: a_odd.representative().attached_out_degree()
2
sage: a_even.representative().attached_out_degree()
2
```

```
sage: a = AbelianStratum(0,4,2,marked_separatrix='out')
sage: a_odd, a_even = a.connected_components()
sage: a_odd.representative().attached_out_degree()
0
sage: a_even.representative().attached_out_degree()
0
```

```
sage: a = AbelianStratum(3,2,1,marked_separatrix='out')
sage: a_c = a.connected_components()[0]
sage: a_c.representative().attached_out_degree()
3
```

```
sage: a = AbelianStratum(2,3,1,marked_separatrix='out')
sage: a_c = a.connected_components()[0]
sage: a_c.representative().attached_out_degree()
2
```

```
sage: a = AbelianStratum(1,3,2,marked_separatrix='out')
sage: a_c = a.connected_components()[0]
sage: a_c.representative().attached_out_degree()
1
```

Tests for incoming separatrices:

```
sage: a = AbelianStratum(4,2,0,marked_separatrix='in')
sage: a_odd, a_even = a.connected_components()
sage: a_odd.representative().attached_in_degree()
4
sage: a_even.representative().attached_in_degree()
4
```

```
sage: a = AbelianStratum(2,4,0,marked_separatrix='in')
sage: a_odd, a_even = a.connected_components()
```

```
sage: a_odd.representative().attached_in_degree()
2
sage: a_even.representative().attached_in_degree()
2
```

```
sage: a = AbelianStratum(0,4,2,marked_separatrix='in')
sage: a_odd, a_even = a.connected_components()
sage: a_odd.representative().attached_in_degree()
0
sage: a_even.representative().attached_in_degree()
0
```

```
sage: a = AbelianStratum(3,2,1,marked_separatrix='in')
sage: a_c = a.connected_components()[0]
sage: a_c.representative().attached_in_degree()
3
```

```
sage: a = AbelianStratum(2,3,1,marked_separatrix='in')
sage: a_c = a.connected_components()[0]
sage: a_c.representative().attached_in_degree()
2
```

```
sage: a = AbelianStratum(1,3,2,marked_separatrix='in')
sage: a_c = a.connected_components()[0]
sage: a_c.representative().attached_in_degree()
1
```

genus()

Returns the genus of the surfaces in this connected component.

OUTPUT:

integer - the genus of the surface

EXAMPLES:

```
sage: a = AbelianStratum(6,4,2,0,0)
sage: c_odd, c_even = a.connected_components()
sage: c_odd.genus()
7
sage: c_even.genus()
7
```

```
sage: a = AbelianStratum([1] *8)
sage: c = a.connected_components()[0]
sage: c.genus()
```

nintervals()

Returns the number of intervals of the representative.

OUTPUT:

integer – the number of intervals in any representative

```
sage: a = AbelianStratum(6,4,2,0,0)
sage: c_odd, c_even = a.connected_components()
sage: c_odd.nintervals()
18
sage: c_even.nintervals()
18
```

```
sage: a = AbelianStratum([1]*8)
sage: c = a.connected_components()[0]
sage: c.nintervals()
17
```

parent()

The stratum of this component

OUTPUT:

stratum - the stratum where this component leaves

EXAMPLES:

```
sage: p = iet.Permutation('a b','b a')
doctest:warning
DeprecationWarning: Permutation is deprecated and will be removed from Sage.
You are advised to install the surface_dynamics package via:
   sage -pip install surface_dynamics
If you do not have write access to the Sage installation you can
alternatively do
   sage -pip install surface_dynamics --user
The package surface_dynamics subsumes all flat surface related
computation that are currently available in Sage. See more
information at
   http://www.labri.fr/perso/vdelecro/surface-dynamics/latest/
See http://trac.sagemath.org/20695 for details.
sage: c = p.connected_component()
sage: c.parent()
H(0)
```

rauzy_diagram(reduced=True)

Returns the Rauzy diagram associated to this connected component.

OUTPUT:

rauzy diagram - the Rauzy diagram associated to this stratum

EXAMPLES:

```
sage: c = AbelianStratum(0).connected_components()[0]
sage: r = c.rauzy_diagram()
```

representative (reduced=True, alphabet=None)

Returns the Zorich representative of this connected component.

Zorich constructs explicitely interval exchange transformations for each stratum in [Zor2008].

INPUT:

- reduced boolean (default: True): whether you obtain a reduced or labelled permutation
- alphabet an alphabet or None: whether you want to specify an alphabet for your permutation

OUTPUT:

permutation – a permutation which lives in this component

EXAMPLES:

```
sage: c = AbelianStratum(1,1,1,1).connected_components()[0]
sage: c
H_c(1, 1, 1, 1)
sage: p = c.representative(alphabet=range(9))
sage: p
0 1 2 3 4 5 6 7 8
4 3 2 5 8 7 6 1 0
sage: p.connected_component()
H_c(1, 1, 1, 1)
```

sage.dynamics.flat_surfaces.strata.EvenCCA

alias of EvenConnectedComponentOfAbelianStratum

Connected component of Abelian stratum with even spin structure.

```
Warning: Internal class! Do not use directly!
```

representative (reduced=True, alphabet=None)

Returns the Zorich representative of this connected component.

Zorich constructs explicitely interval exchange transformations for each stratum in [Zor2008].

EXAMPLES:

```
sage: c = AbelianStratum(6).connected_components()[2]
sage: c
H_even(6)
sage: p = c.representative(alphabet=range(8))
sage: p
0 1 2 3 4 5 6 7
5 4 3 2 7 6 1 0
sage: p.connected_component()
H_even(6)
```

```
sage: c = AbelianStratum(4,4).connected_components()[2]
sage: c
H_even(4, 4)
sage: p = c.representative(alphabet=range(11))
sage: p
0 1 2 3 4 5 6 7 8 9 10
5 4 3 2 6 8 7 10 9 1 0
sage: p.connected_component()
H_even(4, 4)
```

sage.dynamics.flat_surfaces.strata.HypCCA

 $alias\ of\ HypConnectedComponentOfAbelianStratum$

Hyperelliptic component of Abelian stratum.

```
Warning: Internal class! Do not use directly!
```

representative (reduced=True, alphabet=None)

Returns the Zorich representative of this connected component.

Zorich constructs explicitely interval exchange transformations for each stratum in [Zor2008].

INPUT:

- reduced boolean (default: True): whether you obtain a reduced or labelled permutation
- alphabet alphabet or None (default: None): whether you want to specify an alphabet for your representative

```
sage: c = AbelianStratum(0).connected_components()[0]
sage: c
H_hyp(0)
sage: p = c.representative(alphabet="01")
sage: p
0 1
1 0
sage: p.connected_component()
H_hyp(0)
```

```
sage: c = AbelianStratum(0,0).connected_components()[0]
sage: c
H_hyp(0, 0)
sage: p = c.representative(alphabet="abc")
sage: p
a b c
c b a
sage: p.connected_component()
H_hyp(0, 0)
```

```
sage: c = AbelianStratum(2).connected_components()[0]
sage: c
H_hyp(2)
sage: p = c.representative(alphabet="ABCD")
sage: p
A B C D
D C B A
sage: p.connected_component()
H_hyp(2)
```

```
sage: c = AbelianStratum(1,1).connected_components()[0]
sage: c
H_hyp(1, 1)
sage: p = c.representative(alphabet="01234")
sage: p
0 1 2 3 4
4 3 2 1 0
sage: p.connected_component()
H_hyp(1, 1)
```

```
sage.dynamics.flat_surfaces.strata.NonHypCCA
alias of NonHypConnectedComponentOfAbelianStratum
```

 $\textbf{class} \ \ \texttt{sage.dynamics.flat_surfaces.strata.NonHypConnectedComponentOfAbelianStratum} \ (\textit{parent})$

 $\textbf{Bases: } \textit{sage.dynamics.flat_surfaces.strata.ConnectedComponentOfAbelianStratum}$

Non hyperelliptic component of Abelian stratum.

```
Warning: Internal class! Do not use directly!
```

```
sage.dynamics.flat_surfaces.strata.OddCCA
alias of OddConnectedComponentOfAbelianStratum
```

Connected component of an Abelian stratum with odd spin parity.

```
Warning: Internal class! Do not use directly!
```

representative (reduced=True, alphabet=None)

Returns the Zorich representative of this connected component.

Zorich constructs explicitely interval exchange transformations for each stratum in [Zor2008].

EXAMPLES:

```
sage: a = AbelianStratum(6).connected_components()[1]
sage: a.representative(alphabet=range(8))
0 1 2 3 4 5 6 7
3 2 5 4 7 6 1 0
```

```
sage: a = AbelianStratum(4,4).connected_components()[1]
sage: a.representative(alphabet=range(11))
0 1 2 3 4 5 6 7 8 9 10
3 2 5 4 6 8 7 10 9 1 0
```

3.2 Strata of quadratic differentials on Riemann surfaces

Warning: This module is deprecated. You are advised to install and use the surface_dynamics package instead available at https://pypi.python.org/pypi/surface_dynamics/

```
\begin{tabular}{ll} \textbf{class} & \texttt{sage.dynamics.flat\_surfaces.quadratic\_strata.QuadraticStratum} (*l) \\ \textbf{Bases:} & \texttt{sage.structure.sage\_object.SageObject} \end{tabular}
```

Stratum of quadratic differentials.

genus()

Returns the genus.

```
sage: QuadraticStratum(-1,-1,-1,-1).genus()
0
```

INTERVAL EXCHANGE TRANSFORMATIONS AND LINEAR INVOLUTIONS

4.1 Class factories for Interval exchange transformations.

Warning: This module is deprecated. You are advised to install and use the surface_dynamics package instead available at https://pypi.python.org/pypi/surface_dynamics/

This library is designed for the usage and manipulation of interval exchange transformations and linear involutions. It defines specialized types of permutation (constructed using iet.Permutation()) some associated graph (constructed using iet.RauzyGraph()) and some maps of intervals (constructed using iet.IntervalExchangeTransformation()).

EXAMPLES:

Creation of an interval exchange transformation:

```
sage: T = iet.IntervalExchangeTransformation(('a b','b a'),(sqrt(2),1))
doctest:warning
DeprecationWarning: IntervalExchangeTransformation is deprecated and will be removed_
→from Sage.
You are advised to install the surface_dynamics package via:
    sage -pip install surface_dynamics
If you do not have write access to the Sage installation you can
alternatively do
    sage -pip install surface_dynamics --user
The package surface_dynamics subsumes all flat surface related
computation that are currently available in Sage. See more
information at
   http://www.labri.fr/perso/vdelecro/surface-dynamics/latest/
See http://trac.sagemath.org/20695 for details.
doctest:warning
DeprecationWarning: Permutation is deprecated and will be removed from Sage.
You are advised to install the surface_dynamics package via:
   sage -pip install surface_dynamics
If you do not have write access to the Sage installation you can
alternatively do
    sage -pip install surface_dynamics --user
The package surface_dynamics subsumes all flat surface related
computation that are currently available in Sage. See more
information at
```

```
http://www.labri.fr/perso/vdelecro/surface-dynamics/latest/
See http://trac.sagemath.org/20695 for details.

sage: T
Interval exchange transformation of [0, sqrt(2) + 1[ with permutation a b b a
```

It can also be initialized using permutation (group theoretic ones):

```
sage: p = Permutation([3,2,1])
sage: T = iet.IntervalExchangeTransformation(p, [1/3,2/3,1])
sage: T
Interval exchange transformation of [0, 2[ with permutation
1 2 3
3 2 1
```

For the manipulation of permutations of iet, there are special types provided by this module. All of them can be constructed using the constructor iet.Permutation. For the creation of labelled permutations of interval exchange transformation:

```
sage: p1 = iet.Permutation('a b c', 'c b a')
sage: p1
a b c
c b a
```

They can be used for initialization of an iet:

```
sage: p = iet.Permutation('a b','b a')
sage: T = iet.IntervalExchangeTransformation(p, [1,sqrt(2)])
sage: T
Interval exchange transformation of [0, sqrt(2) + 1[ with permutation
a b
b a
```

You can also, create labelled permutations of linear involutions:

```
sage: p = iet.GeneralizedPermutation('a a b', 'b c c')
doctest:warning
DeprecationWarning: GeneralizedPermutation is deprecated and will be removed from __
→Sage.
You are advised to install the surface_dynamics package via:
   sage -pip install surface_dynamics
If you do not have write access to the Sage installation you can
alternatively do
   sage -pip install surface_dynamics --user
The package surface_dynamics subsumes all flat surface related
computation that are currently available in Sage. See more
information at
   http://www.labri.fr/perso/vdelecro/surface-dynamics/latest/
See http://trac.sagemath.org/20695 for details.
sage: p
a a b
рсс
```

Sometimes it's more easy to deal with reduced permutations:

```
sage: p = iet.Permutation('a b c', 'c b a', reduced = True)
sage: p
a b c
c b a
```

Permutations with flips:

```
sage: p1 = iet.Permutation('a b c', 'c b a', flips = ['a','c'])
sage: p1
-a b -c
-c b -a
```

Creation of Rauzy diagrams:

```
sage: r = iet.RauzyDiagram('a b c', 'c b a')
doctest:warning
...
DeprecationWarning: RauzyDiagram is deprecated and will be removed from Sage.
You are advised to install the surface_dynamics package via:
    sage -pip install surface_dynamics
If you do not have write access to the Sage installation you can
alternatively do
    sage -pip install surface_dynamics --user
The package surface_dynamics subsumes all flat surface related
computation that are currently available in Sage. See more
information at
    http://www.labri.fr/perso/vdelecro/surface-dynamics/latest/
See http://trac.sagemath.org/20695 for details.
```

Reduced Rauzy diagrams are constructed using the same arguments than for permutations:

```
sage: r = iet.RauzyDiagram('a b b','c c a')
sage: r_red = iet.RauzyDiagram('a b b','c c a',reduced=True)
sage: r.cardinality()
12
sage: r_red.cardinality()
4
```

By default, Rauzy diagrams are generated by induction on the right. You can use several options to enlarge (or restrict) the diagram (try help(iet.RauzyDiagram) for more precisions):

```
sage: r1 = iet.RauzyDiagram('a b c','c b a',right_induction=True)
sage: r2 = iet.RauzyDiagram('a b c','c b a',left_right_inversion=True)
```

You can consider self similar iet using path in Rauzy diagrams and eigenvectors of the corresponding matrix:

```
sage: p = iet.Permutation("a b c d", "d c b a")
sage: d = p.rauzy_diagram()
sage: g = d.path(p, 't', 't', 'b', 't', 'b', 't', 'b')
sage: g
Path of length 8 in a Rauzy diagram
sage: g.is_loop()
True
sage: g.is_full()
True
sage: m = g.matrix()
sage: v = m.eigenvectors_right()[-1][1][0]
sage: T1 = iet.IntervalExchangeTransformation(p, v)
```

```
sage: T2 = T1.rauzy_move(iterations=8)
sage: T1.normalize(1) == T2.normalize(1)
True
```

REFERENCES:

- [BL2008]
- [DN1990]
- [Nog1985]
- [Rau1979]
- [Vee1978]
- [Zor]

AUTHORS:

• Vincent Delecroix (2009-09-29): initial version

Returns a permutation of an interval exchange transformation.

Those permutations are the combinatoric part of linear involutions and were introduced by Danthony-Nogueira [DN1990]. The full combinatoric study and precise links with strata of quadratic differentials was achieved few years later by Boissy-Lanneau [BL2008].

INPUT:

- intervals strings, list, tuples
- reduced boolean (default: False) specifies reduction. False means labelled permutation and True means reduced permutation.
- flips iterable (default: None) the letters which correspond to flipped intervals.

OUTPUT

generalized permutation – the output type depends on the data.

EXAMPLES:

Creation of labelled generalized permutations:

```
sage: iet.GeneralizedPermutation('a b b','c c a')
a b b
c c a
sage: iet.GeneralizedPermutation('a a','b b c c')
a a
b b c c
sage: iet.GeneralizedPermutation([[0,1,2,3,1],[4,2,5,3,5,4,0]])
0 1 2 3 1
4 2 5 3 5 4 0
```

Creation of reduced generalized permutations:

```
sage: iet.GeneralizedPermutation('a b b', 'c c a', reduced = True)
a b b
c c a
sage: iet.GeneralizedPermutation('a a b b', 'c c d d', reduced = True)
```

```
a a b b c c d d
```

Creation of flipped generalized permutations:

```
sage: iet.GeneralizedPermutation('a b c a', 'd c d b', flips = ['a','b'])
-a -b c -a
d c d -b
```

sage.dynamics.interval_exchanges.constructors.IET (permutation=None, lengths=None)
Constructs an Interval exchange transformation.

An interval exchange transformation (or iet) is a map from an interval to itself. It is defined on the interval except at a finite number of points (the singularities) and is a translation on each connected component of the complement of the singularities. Moreover it is a bijection on its image (or it is injective).

An interval exchange transformation is encoded by two datas. A permutation (that corresponds to the way we echange the intervals) and a vector of positive reals (that corresponds to the lengths of the complement of the singularities).

INPUT:

- permutation a permutation
- lengths a list or a dictionary of lengths

OUTPUT:

interval exchange transformation – an map of an interval

EXAMPLES:

Two initialization methods, the first using a iet.Permutation:

```
sage: p = iet.Permutation('a b c','c b a')
sage: t = iet.IntervalExchangeTransformation(p, {'a':1,'b':0.4523,'c':2.8})
```

The second is more direct:

```
sage: t = iet.IntervalExchangeTransformation(('a b','b a'),{'a':1,'b':4})
```

It's also possible to initialize the lengths only with a list:

```
sage: t = iet.IntervalExchangeTransformation(('a b c','c b a'),[0.123,0.4,2])
```

The two fundamental operations are Rauzy move and normalization:

```
sage: t = iet.IntervalExchangeTransformation(('a b c','c b a'),[0.123,0.4,2])
sage: s = t.rauzy_move()
sage: s_n = s.normalize(t.length())
sage: s_n.length() == t.length()
True
```

A not too simple example of a self similar interval exchange transformation:

```
sage: p = iet.Permutation('a b c d','d c b a')
sage: d = p.rauzy_diagram()
sage: g = d.path(p, 't', 't', 'b', 't', 'b', 't', 'b')
sage: m = g.matrix()
sage: v = m.eigenvectors_right()[-1][1][0]
```

```
sage: t = iet.IntervalExchangeTransformation(p,v)
sage: s = t.rauzy_move(iterations=8)
sage: s.normalize() == t.normalize()
True
```

Constructs an Interval exchange transformation.

An interval exchange transformation (or iet) is a map from an interval to itself. It is defined on the interval except at a finite number of points (the singularities) and is a translation on each connected component of the complement of the singularities. Moreover it is a bijection on its image (or it is injective).

An interval exchange transformation is encoded by two datas. A permutation (that corresponds to the way we echange the intervals) and a vector of positive reals (that corresponds to the lengths of the complement of the singularities).

INPUT:

- permutation a permutation
- lengths a list or a dictionary of lengths

OUTPUT:

interval exchange transformation - an map of an interval

EXAMPLES:

Two initialization methods, the first using a iet.Permutation:

```
sage: p = iet.Permutation('a b c','c b a')
sage: t = iet.IntervalExchangeTransformation(p, {'a':1,'b':0.4523,'c':2.8})
```

The second is more direct:

```
sage: t = iet.IntervalExchangeTransformation(('a b','b a'),{'a':1,'b':4})
```

It's also possible to initialize the lengths only with a list:

```
sage: t = iet.IntervalExchangeTransformation(('a b c','c b a'),[0.123,0.4,2])
```

The two fundamental operations are Rauzy move and normalization:

```
sage: t = iet.IntervalExchangeTransformation(('a b c','c b a'),[0.123,0.4,2])
sage: s = t.rauzy_move()
sage: s_n = s.normalize(t.length())
sage: s_n.length() == t.length()
True
```

A not too simple example of a self similar interval exchange transformation:

```
sage: p = iet.Permutation('a b c d','d c b a')
sage: d = p.rauzy_diagram()
sage: g = d.path(p, 't', 't', 'b', 't', 'b', 't', 'b')
sage: m = g.matrix()
sage: v = m.eigenvectors_right()[-1][1][0]
sage: t = iet.IntervalExchangeTransformation(p,v)
sage: s = t.rauzy_move(iterations=8)
sage: s.normalize() == t.normalize()
```

```
sage.dynamics.interval_exchanges.constructors.Permutation(*args, **kargs)
Returns a permutation of an interval exchange transformation.
```

Those permutations are the combinatoric part of an interval exchange transformation (IET). The combinatorial study of those objects starts with Gerard Rauzy [Rau1979] and William Veech [Vee1978].

The combinatoric part of interval exchange transformation can be taken independently from its dynamical origin. It has an important link with strata of Abelian differential (see strata)

INPUT:

- intervals string, two strings, list, tuples that can be converted to two lists
- reduced boolean (default: False) specifies reduction. False means labelled permutation and True means reduced permutation.
- flips iterable (default: None) the letters which correspond to flipped intervals.

OUTPUT:

permutation – the output type depends of the data.

EXAMPLES:

Creation of labelled permutations

```
sage: iet.Permutation('a b c d','d c b a')
a b c d
d c b a
sage: iet.Permutation([[0,1,2,3],[2,1,3,0]])
0 1 2 3
2 1 3 0
sage: iet.Permutation([[0, 'A', 'B', 1], ['B', 0, 1, 'A']))
0 A B 1
B 0 1 A
```

Creation of reduced permutations:

```
sage: iet.Permutation('a b c', 'c b a', reduced = True)
a b c
c b a
sage: iet.Permutation([0, 1, 2, 3], [1, 3, 0, 2])
0 1 2 3
1 3 0 2
```

Creation of flipped permutations:

```
sage: iet.Permutation('a b c', 'c b a', flips=['a','b'])
-a -b c
c -b -a
sage: iet.Permutation('a b c', 'c b a', flips=['a'], reduced=True)
-a b c
c b -a
```

```
sage: p = iet.Permutation('a b c','c b a')
sage: iet.Permutation(p) == p
True
sage: iet.Permutation(p, reduced=True) == p.reduced()
True
```

```
sage: p = iet.Permutation('a', 'a', flips='a', reduced=True)
sage: iet.Permutation(p) == p
True
```

```
sage: p = iet.Permutation('a b c','c b a',flips='a')
sage: iet.Permutation(p) == p
True
sage: iet.Permutation(p, reduced=True) == p.reduced()
True
```

```
sage: p = iet.Permutation('a b c','c b a',reduced=True)
sage: iet.Permutation(p) == p
True

sage: iet.Permutation('a b c','c b a',reduced='badly')
Traceback (most recent call last):
...
TypeError: reduced must be of type boolean
sage: iet.Permutation('a','a',flips='b',reduced=True)
Traceback (most recent call last):
...
ValueError: flips contains not valid letters
sage: iet.Permutation('a b c','c a a',reduced=True)
Traceback (most recent call last):
...
ValueError: letters must appear once in each interval
```

Returns an iterator over permutations.

This iterator allows you to iterate over permutations with given constraints. If you want to iterate over permutations coming from a given stratum you have to use the module <code>strata</code> and generate Rauzy diagrams from connected components.

INPUT:

- nintervals non negative integer
- irreducible boolean (default: True)
- reduced boolean (default: False)
- alphabet alphabet (default: None)

OUTPUT:

iterator - an iterator over permutations

EXAMPLES:

Generates all reduced permutations with given number of intervals:

```
sage: P = iet.Permutations_iterator(nintervals=2,alphabet="ab",reduced=True)
doctest:warning
```

```
DeprecationWarning: iet_Permutations_iterator is deprecated and will be removed.
→from Sage.
You are advised to install the surface_dynamics package via:
   sage -pip install surface_dynamics
If you do not have write access to the Sage installation you can
alternatively do
    sage -pip install surface_dynamics --user
The package surface_dynamics subsumes all flat surface related
computation that are currently available in Sage. See more
information at
    http://www.labri.fr/perso/vdelecro/surface-dynamics/latest/
See http://trac.sagemath.org/20695 for details.
sage: for p in P:
       print(p)
. . . . :
        print("* *")
. . . . :
a b
b a
* *
sage: P = iet.Permutations_iterator(nintervals=3,alphabet="abc",reduced=True)
sage: for p in P:
         print(p)
. . . . :
         print("* * *")
. . . . :
a b c
bса
* * *
a b c
c a b
a b c
c b a
* * *
```

sage.dynamics.interval_exchanges.constructors.RauzyDiagram(*args, **kargs)
Return an object coding a Rauzy diagram.

The Rauzy diagram is an oriented graph with labelled edges. The set of vertices corresponds to the permutations obtained by different operations (mainly the .rauzy_move() operations that corresponds to an induction of interval exchange transformation). The edges correspond to the action of the different operations considered.

It first appeared in the original article of Rauzy [Rau1979].

INPUT:

- intervals lists, or strings, or tuples
- reduced boolean (default: False) to precise reduction
- flips list (default: []) for flipped permutations
- right induction boolean (default: True) consideration of left induction in the diagram
- left_induction boolean (default: False) consideration of right induction in the diagram
- left_right_inversion boolean (default: False) consideration of inversion
- top_bottom_inversion boolean (default: False) consideration of reversion
- symmetric boolean (default: False) consideration of the symmetric operation

OUTPUT:

Rauzy diagram – the Rauzy diagram that corresponds to your request

EXAMPLES:

Standard Rauzy diagrams:

```
sage: iet.RauzyDiagram('a b c d', 'd b c a')
Rauzy diagram with 12 permutations
sage: iet.RauzyDiagram('a b c d', 'd b c a', reduced = True)
Rauzy diagram with 6 permutations
```

Extended Rauzy diagrams:

```
sage: iet.RauzyDiagram('a b c d', 'd b c a', symmetric=True)
Rauzy diagram with 144 permutations
```

Using Rauzy diagrams and path in Rauzy diagrams:

```
sage: r = iet.RauzyDiagram('a b c', 'c b a')
sage: r
Rauzy diagram with 3 permutations
sage: p = iet.Permutation('a b c','c b a')
sage: p in r
True
sage: g0 = r.path(p, 'top', 'bottom', 'top')
sage: g1 = r.path(p, 'bottom', 'top', 'bottom')
sage: g0.is_loop(), g1.is_loop()
(True, True)
sage: g0.is_full(), g1.is_full()
(False, False)
sage: g = g0 + g1
sage: g
Path of length 6 in a Rauzy diagram
sage: g.is_loop(), g.is_full()
(True, True)
sage: m = q.matrix()
sage: m
[1 1 1]
[2 4 1]
[2 3 2]
sage: s = q.orbit_substitution()
sage: s
WordMorphism: a->acbbc, b->acbbcbbc, c->acbc
sage: s.incidence_matrix() == m
True
```

We can then create the corresponding interval exchange transformation and comparing the orbit of 0 to the fixed point of the orbit substitution:

```
sage: v = m.eigenvectors_right()[-1][1][0]
sage: T = iet.IntervalExchangeTransformation(p, v).normalize()
sage: T
Interval exchange transformation of [0, 1[ with permutation
a b c
c b a
sage: w1 = []
sage: x = 0
sage: for i in range(20):
...: w1.append(T.in_which_interval(x))
```

```
....: x = T(x)
sage: w1 = Word(w1)
sage: w1
word: acbbcacbcacbbcbbcacb
sage: w2 = s.fixed_point('a')
sage: w2[:20]
word: acbbcacbcacbbcbbcacb
sage: w2[:20] == w1
True
```

4.2 Labelled permutations

Warning: This module is deprecated. You are advised to install and use the surface_dynamics package instead available at https://pypi.python.org/pypi/surface_dynamics/

A labelled (generalized) permutation is better suited to study the dynamic of a translation surface than a reduced one (see the module <code>sage.dynamics.interval_exchanges.reduced</code>). The latter is more adapted to the study of strata. This kind of permutation was introduced by Yoccoz [Yoc2005] (see also [MMY2003]).

In fact, there is a geometric counterpart of labelled permutations. They correspond to translation surfaces with marked outgoing separatrices (i.e. we fix a label for each of them).

Remarks that Rauzy diagram of reduced objects are significantly smaller than the one for labelled object (for the permutation a b d b e / e d c a c the labelled Rauzy diagram contains 8760 permutations, and the reduced only 73). But, as it is in geometrical way, the labelled Rauzy diagram is a covering of the reduced Rauzy diagram.

AUTHORS:

• Vincent Delecroix (2009-09-29): initial version

Bases: sage.dynamics.interval_exchanges.labelled.LabelledPermutation

General template for labelled objects

```
Warning: Internal class! Do not use directly!

list (flips=False)
```

Returns a list associated to the permutation.

```
INPUT:
```

• flips - boolean (default: False)

OUTPUT:

list – two lists of labels

EXAMPLES:

flips=None)

The list can be used to reconstruct the permutation

```
sage: p = iet.Permutation('a b c','c b a',flips='ab')
sage: p == iet.Permutation(p.list(), flips=p.flips())
True
```

```
sage: p = iet.GeneralizedPermutation('a b b c','c d d a',flips='ad')
sage: p == iet.GeneralizedPermutation(p.list(),flips=p.flips())
True
```

 $\textbf{class} \texttt{ sage.dynamics.interval_exchanges.labelled.FlippedLabelledPermutationIET} (\textit{intervals=None}, \textit{class}) \\$

alphabet=None,
flips=None)

Bases: sage.dynamics.interval_exchanges.labelled.FlippedLabelledPermutation, sage.dynamics.interval_exchanges.template.FlippedPermutationIET, sage.dynamics.interval exchanges.labelled.LabelledPermutationIET

Flipped labelled permutation from iet.

EXAMPLES:

Reducibility testing (does not depends of flips):

```
sage: p = iet.Permutation('a b c', 'c b a',flips='a')
sage: p.is_irreducible()
True
sage: q = iet.Permutation('a b c d', 'b a d c', flips='bc')
sage: q.is_irreducible()
False
```

Rauzy movability and Rauzy move:

```
sage: p = iet.Permutation('a b c', 'c b a',flips='a')
sage: p
-a b c
c b -a
sage: p.rauzy_move(1)
-c -a b
-c b -a
sage: p.rauzy_move(0)
-a b c
c -a b
```

Rauzy diagrams:

```
sage: d = iet.RauzyDiagram('a b c d','d a b c',flips='a')
doctest:warning
...
DeprecationWarning: RauzyDiagram is deprecated and will be removed from Sage.
```

```
You are advised to install the surface_dynamics package via:
sage -pip install surface_dynamics

If you do not have write access to the Sage installation you can alternatively do sage -pip install surface_dynamics --user

The package surface_dynamics subsumes all flat surface related computation that are currently available in Sage. See more information at http://www.labri.fr/perso/vdelecro/surface-dynamics/latest/
See http://trac.sagemath.org/20695 for details.
```

AUTHORS:

• Vincent Delecroix (2009-09-29): initial version

rauzy_diagram(**kargs)

Returns the Rauzy diagram associated to this permutation.

For more information, try help(iet.RauzyDiagram)

OUTPUT:

RauzyDiagram - the Rauzy diagram of self

EXAMPLES:

```
sage: p = iet.Permutation('a b c', 'c b a',flips='a')
sage: p.rauzy_diagram()
Rauzy diagram with 3 permutations
```

rauzy_move (winner=None, side=None)

Returns the Rauzy move.

INPUT:

- winner 'top' (or 't' or 0) or 'bottom' (or 'b' or 1)
- side (default: 'right') 'right' (or 'r') or 'left' (or 'l')

OUTPUT:

permutation - the Rauzy move of self

```
sage: p = iet.Permutation('a b','b a',flips='a')
sage: p.rauzy_move('top')
-a b
b -a
sage: p.rauzy_move('bottom')
-b -a
-b -a
```

```
sage: p = iet.Permutation('a b c','c b a',flips='b')
sage: p.rauzy_move('top')
a -b c
c a -b
sage: p.rauzy_move('bottom')
a c -b
c -b a
```

reduced()

The associated reduced permutation.

OUTPUT:

permutation - the associated reduced permutation

EXAMPLES:

```
sage: p = iet.Permutation('a b c','c b a',flips='a')
sage: q = iet.Permutation('a b c','c b a',flips='a',reduced=True)
sage: p.reduced() == q
True
```

class sage.dynamics.interval_exchanges.labelled.FlippedLabelledPermutationLI(intervals=None,

alphabet=None,
flips=None)

Bases: sage.dynamics.interval_exchanges.labelled.FlippedLabelledPermutation, sage.dynamics.interval_exchanges.template.FlippedPermutationLI, sage.dynamics.interval_exchanges.labelled.LabelledPermutationLI

Flipped labelled quadratic (or generalized) permutation.

EXAMPLES:

Reducibility testing:

```
sage: p = iet.GeneralizedPermutation('a b b', 'c c a', flips='a')
sage: p.is_irreducible()
True
```

Reducibility testing with associated decomposition:

```
sage: p = iet.GeneralizedPermutation('a b c a', 'b d d c', flips='ab')
sage: p.is_irreducible()
False
sage: test, decomp = p.is_irreducible(return_decomposition = True)
sage: test
False
sage: decomp
(['a'], ['c', 'a'], [], ['c'])
```

Rauzy movability and Rauzy move:

```
sage: p = iet.GeneralizedPermutation('a a b b c c', 'd d', flips='d')
sage: p.has_rauzy_move(0)
False
sage: p.has_rauzy_move(1)
True
sage: p = iet.GeneralizedPermutation('a a b','b c c',flips='c')
sage: p.has_rauzy_move(0)
True
sage: p.has_rauzy_move(1)
True
```

left_rauzy_move(winner)

Perform a Rauzy move on the left.

INPUT:

• winner - either 'top' or 'bottom' ('t' or 'b' for short)

OUTPUT:

- a permutation

EXAMPLES:

```
sage: p = iet.GeneralizedPermutation('a a b','b c c')
sage: p.left_rauzy_move(0)
a a b b
c c
sage: p.left_rauzy_move(1)
a a b
b c c
```

```
sage: p = iet.GeneralizedPermutation('a b b','c c a')
sage: p.left_rauzy_move(0)
a b b
c c a
sage: p.left_rauzy_move(1)
b b
c c a a
```

rauzy_diagram(**kargs)

Returns the associated Rauzy diagram.

For more information, try help(RauzyDiagram)

OUTPUT:

- a RauzyDiagram

EXAMPLES:

```
sage: p = iet.GeneralizedPermutation('a b b a', 'c d c d')
sage: d = p.rauzy_diagram()
```

reduced()

The associated reduced permutation.

OUTPUT

permutation – the associated reduced permutation

EXAMPLES:

```
sage: p = iet.GeneralizedPermutation('a a','b b c c',flips='a')
sage: q = iet.GeneralizedPermutation('a a','b b c c',flips='a',reduced=True)
sage: p.reduced() == q
True
```

right_rauzy_move (winner)

Perform a Rauzy move on the right (the standard one).

INPUT:

• winner - either 'top' or 'bottom' ('t' or 'b' for short)

OUTPUT:

permutation - the Rauzy move of self

EXAMPLES:

```
sage: p = iet.GeneralizedPermutation('a a b','b c c',flips='c')
sage: p.right_rauzy_move(0)
a a b
-c b -c
sage: p.right_rauzy_move(1)
a a
-b -c -b -c
```

```
sage: p = iet.GeneralizedPermutation('a b b','c c a',flips='ab')
sage: p.right_rauzy_move(0)
a -b a -b
c c
sage: p.right_rauzy_move(1)
b -a b
c c -a
```

class sage.dynamics.interval_exchanges.labelled.FlippedLabelledRauzyDiagram (p,

right_induction=True left_induction=False, left_right_inversion= top_bottom_inversion symmet-

Bases: sage.dynamics.interval_exchanges.template.FlippedRauzyDiagram, sage.dynamics.interval_exchanges.labelled.LabelledRauzyDiagram

Rauzy diagram of flipped labelled permutations

 $\textbf{class} \ \, \textbf{sage.dynamics.interval_exchanges.labelled.LabelledPermutation} \, (\textit{intervals=None}, \\ \textit{alpha-bet=None})$

Bases: sage.structure.sage_object.SageObject

General template for labelled objects.

```
Warning: Internal class! Do not use directly!
```

erase letter(letter)

Return the permutation with the specified letter removed.

OUTPUT:

permutation – the resulting permutation

```
sage: p = iet.Permutation('a b c d','c d b a')
sage: p.erase_letter('a')
b c d
c d b
sage: p.erase_letter('b')
a c d
c d a
sage: p.erase_letter('c')
a b d
```

```
d b a
sage: p.erase_letter('d')
a b c
c b a
```

```
sage: p = iet.GeneralizedPermutation('a b b','c c a')
sage: p.erase_letter('a')
b b
c c
```

Beware, there is no validity check for permutation from linear involutions:

```
sage: p = iet.GeneralizedPermutation('a b b','c c a')
sage: p.erase_letter('b')
a
c c a
```

length (interval=None)

Returns a 2-uple of lengths.

p.length() is identical to (p.length_top(), p.length_bottom()) If an interval is specified, it returns the length of the specified interval.

INPUT:

• interval - None, 'top' or 'bottom'

OUTPUT:

tuple – a 2-uple of integers

EXAMPLES:

```
sage: iet.Permutation('a b c','c b a').length()
(3, 3)
sage: iet.GeneralizedPermutation('a a','b b c c').length()
(2, 4)
sage: iet.GeneralizedPermutation('a a b b','c c').length()
(4, 2)
```

length_bottom()

Returns the number of intervals in the bottom segment.

OUTPUT:

integer - number of intervals

EXAMPLES:

```
sage: iet.Permutation('a b','b a').length_bottom()
2
sage: iet.GeneralizedPermutation('a a','b b c c').length_bottom()
4
sage: iet.GeneralizedPermutation('a a b b','c c').length_bottom()
2
```

length_top()

Returns the number of intervals in the top segment.

OUTPUT:

integer – number of intervals

EXAMPLES:

```
sage: iet.Permutation('a b c','c b a').length_top()
sage: iet.GeneralizedPermutation('a a','b b c c').length_top()
doctest:warning
. . .
DeprecationWarning: GeneralizedPermutation is deprecated and will be removed,
→from Sage.
You are advised to install the surface_dynamics package via:
   sage -pip install surface_dynamics
If you do not have write access to the Sage installation you can
alternatively do
   sage -pip install surface_dynamics --user
The package surface_dynamics subsumes all flat surface related
computation that are currently available in Sage. See more
information at
   http://www.labri.fr/perso/vdelecro/surface-dynamics/latest/
See http://trac.sagemath.org/20695 for details.
sage: iet.GeneralizedPermutation('a a b b','c c').length_top()
```

list()

Returns a list of two lists corresponding to the intervals.

OUTPUT:

list – two lists of labels

EXAMPLES:

The list of an permutation from iet:

```
sage: p1 = iet.Permutation('1 2 3', '3 1 2')
sage: p1.list()
[['1', '2', '3'], ['3', '1', '2']]
sage: p1.alphabet("abc")
sage: p1.list()
[['a', 'b', 'c'], ['c', 'a', 'b']]
```

Recovering the permutation from this list (and the alphabet):

```
sage: q1 = iet.Permutation(p1.list(),alphabet=p1.alphabet())
sage: p1 == q1
True
```

The list of a quadratic permutation:

```
sage: p2 = iet.GeneralizedPermutation('g o o', 'd d g')
sage: p2.list()
[['g', 'o', 'o'], ['d', 'd', 'g']]
```

Recovering the permutation:

```
sage: q2 = iet.GeneralizedPermutation(p2.list(),alphabet=p2.alphabet())
sage: p2 == q2
True
```

rauzy_move_loser (winner=None, side=None)

Returns the loser of a Rauzy move

INPUT:

- winner either 'top' or 'bottom' ('t' or 'b' for short)
- side either 'left' or 'right' ('l' or 'r' for short)

OUTPUT:

a label

EXAMPLES:

```
sage: p = iet.Permutation('a b c d','b d a c')
sage: p.rauzy_move_loser('top','right')
'c'
sage: p.rauzy_move_loser('bottom','right')
'd'
sage: p.rauzy_move_loser('top','left')
'b'
sage: p.rauzy_move_loser('bottom','left')
'a'
```

rauzy_move_matrix(winner=None, side='right')

Returns the Rauzy move matrix.

This matrix corresponds to the action of a Rauzy move on the vector of lengths. By convention (to get a positive matrix), the matrix is defined as the inverse transformation on the length vector.

OUTPUT:

matrix - a square matrix of positive integers

EXAMPLES:

```
sage: p = iet.Permutation('a b','b a')
sage: p.rauzy_move_matrix('t')
[1 0]
[1 1]
sage: p.rauzy_move_matrix('b')
[1 1]
[0 1]
```

```
sage: p = iet.Permutation('a b c d','b d a c')
sage: q = p.left_right_inverse()
sage: m0 = p.rauzy_move_matrix(winner='top', side='right')
sage: n0 = q.rauzy_move_matrix(winner='top', side='left')
sage: m0 == n0
True
sage: m1 = p.rauzy_move_matrix(winner='bottom', side='right')
sage: n1 = q.rauzy_move_matrix(winner='bottom', side='left')
sage: m1 == n1
True
```

rauzy move winner (winner=None, side=None)

Returns the winner of a Rauzy move.

INPUT:

• winner - either 'top' or 'bottom' ('t' or 'b' for short)

• side - either 'left' or 'right' ('l' or 'r' for short)

OUTPUT:

a label

EXAMPLES:

```
sage: p = iet.Permutation('a b c d','b d a c')
sage: p.rauzy_move_winner('top','right')
'd'
sage: p.rauzy_move_winner('bottom','right')
'c'
sage: p.rauzy_move_winner('top','left')
'a'
sage: p.rauzy_move_winner('bottom','left')
'b'
```

```
sage: p = iet.GeneralizedPermutation('a b b c','d c a e d e')
sage: p.rauzy_move_winner('top','right')
'c'
sage: p.rauzy_move_winner('bottom','right')
'e'
sage: p.rauzy_move_winner('top','left')
'a'
sage: p.rauzy_move_winner('bottom','left')
'd'
```

class sage.dynamics.interval_exchanges.labelled.LabelledPermutationIET(intervals=None,

phabet=None)

Bases: sage.dynamics.interval_exchanges.labelled.LabelledPermutation, sage.dynamics.interval_exchanges.template.PermutationIET

Labelled permutation for iet

EXAMPLES:

Reducibility testing:

```
sage: p = iet.Permutation('a b c', 'c b a')
sage: p.is_irreducible()
True

sage: q = iet.Permutation('a b c d', 'b a d c')
sage: q.is_irreducible()
False
```

Rauzy movability and Rauzy move:

```
sage: p = iet.Permutation('a b c', 'c b a')
sage: p.has_rauzy_move('top')
True
sage: p.rauzy_move('bottom')
a c b
c b a
sage: p.has_rauzy_move('top')
True
sage: p.rauzy_move('top')
```

```
a b c c a b
```

Rauzy diagram:

```
sage: p = iet.Permutation('a b c', 'c b a')
sage: d = p.rauzy_diagram()
sage: p in d
True
```

has_rauzy_move (winner=None, side=None)

Returns True if you can perform a Rauzy move.

INPUT:

- winner the winner interval ('top' or 'bottom')
- side (default: 'right') the side ('left' or 'right')

OUTPUT:

bool - True if self has a Rauzy move

EXAMPLES:

```
sage: p = iet.Permutation('a b','b a')
sage: p.has_rauzy_move()
True
```

```
sage: p = iet.Permutation('a b c','b a c')
sage: p.has_rauzy_move()
False
```

is_identity()

Returns True if self is the identity.

OUTPUT:

bool – True if self corresponds to the identity

EXAMPLES:

```
sage: iet.Permutation("a b","a b").is_identity()
True
sage: iet.Permutation("a b","b a").is_identity()
False
```

rauzy_diagram(**args)

Returns the associated Rauzy diagram.

For more information try help(iet.RauzyDiagram).

OUTPUT:

Rauzy diagram - the Rauzy diagram of the permutation

```
sage: p = iet.Permutation('a b c', 'c b a')
sage: d = p.rauzy_diagram()
```

rauzy_move (*winner=None*, *side=None*, *iteration=1*) Returns the Rauzy move.

INPUT:

- winner the winner interval ('top' or 'bottom')
- side (default: 'right') the side ('left' or 'right')

OUTPUT:

permutation – the Rauzy move of the permutation

EXAMPLES:

```
sage: p = iet.Permutation('a b','b a')
sage: p.rauzy_move('t','right')
a b
b a
sage: p.rauzy_move('b','right')
a b
b a
```

```
sage: p = iet.Permutation('a b c','c b a')
sage: p.rauzy_move('t','right')
a b c
c a b
sage: p.rauzy_move('b','right')
a c b
c b a
```

```
sage: p = iet.Permutation('a b','b a')
sage: p.rauzy_move('t','left')
a b
b a
sage: p.rauzy_move('b','left')
a b
b a
```

```
sage: p = iet.Permutation('a b c','c b a')
sage: p.rauzy_move('t','left')
a b c
b c a
sage: p.rauzy_move('b','left')
b a c
c b a
```

rauzy_move_interval_substitution(winner=None, side=None)

Returns the interval substitution associated.

INPUT:

- winner the winner interval ('top' or 'bottom')
- side (default: 'right') the side ('left' or 'right')

OUTPUT

WordMorphism – a substitution on the alphabet of the permutation

```
sage: p = iet.Permutation('a b','b a')
sage: p.rauzy_move_interval_substitution('top','right')
WordMorphism: a->a, b->ba
sage: p.rauzy_move_interval_substitution('bottom','right')
WordMorphism: a->ab, b->b
sage: p.rauzy_move_interval_substitution('top','left')
WordMorphism: a->ba, b->b
sage: p.rauzy_move_interval_substitution('bottom','left')
WordMorphism: a->a, b->ab
```

rauzy_move_orbit_substitution (winner=None, side=None)

Return the action of the rauzy_move on the orbit.

INPUT:

- i integer
- winner the winner interval ('top' or 'bottom')
- side (default: 'right') the side ('right' or 'left')

OUTPUT:

WordMorphism – a substitution on the alphabet of self

EXAMPLES:

```
sage: p = iet.Permutation('a b','b a')
sage: p.rauzy_move_orbit_substitution('top','right')
WordMorphism: a->ab, b->b
sage: p.rauzy_move_orbit_substitution('bottom','right')
WordMorphism: a->a, b->ab
sage: p.rauzy_move_orbit_substitution('top','left')
WordMorphism: a->a, b->ba
sage: p.rauzy_move_orbit_substitution('bottom','left')
WordMorphism: a->ba, b->b
```

reduced()

Returns the associated reduced abelian permutation.

OUTPUT:

a reduced permutation – the underlying reduced permutation

EXAMPLES:

```
sage: p = iet.Permutation("a b c d","d c a b")
sage: q = iet.Permutation("a b c d","d c a b",reduced=True)
sage: p.reduced() == q
True
```

class sage.dynamics.interval_exchanges.labelled.LabelledPermutationLI(intervals=None,

```
pha-
```

al-

bet=None)

Bases: sage.dynamics.interval_exchanges.labelled.LabelledPermutation, sage.dynamics.interval_exchanges.template.PermutationLI

Labelled quadratic (or generalized) permutation

Reducibility testing:

```
sage: p = iet.GeneralizedPermutation('a b b', 'c c a')
sage: p.is_irreducible()
True
```

Reducibility testing with associated decomposition:

```
sage: p = iet.GeneralizedPermutation('a b c a', 'b d d c')
sage: p.is_irreducible()
False
sage: test, decomposition = p.is_irreducible(return_decomposition = True)
sage: test
False
sage: decomposition
(['a'], ['c', 'a'], [], ['c'])
```

Rauzy movability and Rauzy move:

```
sage: p = iet.GeneralizedPermutation('a a b b c c', 'd d')
sage: p.has_rauzy_move(0)
False
sage: p.has_rauzy_move(1)
True
sage: q = p.rauzy_move(1)
sage: q
a a b b c
c d d
sage: q.has_rauzy_move(0)
True
sage: q.has_rauzy_move(1)
True
```

Rauzy diagrams:

```
sage: p = iet.GeneralizedPermutation('0 0 1 1','2 2')
sage: r = p.rauzy_diagram()
sage: p in r
True
```

has_right_rauzy_move (winner)

Test of Rauzy movability with a specified winner

A quadratic (or generalized) permutation is rauzy_movable type depending on the possible length of the last interval. It is dependent of the length equation.

INPUT:

```
• winner - 'top' (or 't' or 0) or 'bottom' (or 'b' or 1)
```

OUTPUT:

bool - True if self has a Rauzy move

```
sage: p = iet.GeneralizedPermutation('a a','b b')
sage: p.has_right_rauzy_move('top')
False
sage: p.has_right_rauzy_move('bottom')
False
```

```
sage: p = iet.GeneralizedPermutation('a a b','b c c')
sage: p.has_right_rauzy_move('top')
True
sage: p.has_right_rauzy_move('bottom')
True
```

```
sage: p = iet.GeneralizedPermutation('a a','b b c c')
sage: p.has_right_rauzy_move('top')
True
sage: p.has_right_rauzy_move('bottom')
False
```

```
sage: p = iet.GeneralizedPermutation('a a b b','c c')
sage: p.has_right_rauzy_move('top')
False
sage: p.has_right_rauzy_move('bottom')
True
```

left_rauzy_move(winner)

Perform a Rauzy move on the left.

INPUT:

• winner - 'top' or 'bottom'

OUTPUT

permutation – the Rauzy move of self

EXAMPLES:

```
sage: p = iet.GeneralizedPermutation('a a b','b c c')
sage: p.left_rauzy_move(0)
a a b b
c c
sage: p.left_rauzy_move(1)
a a b
b c c
```

```
sage: p = iet.GeneralizedPermutation('a b b','c c a')
sage: p.left_rauzy_move(0)
a b b
c c a
sage: p.left_rauzy_move(1)
b b
c c a a
```

rauzy_diagram(**kargs)

Returns the associated RauzyDiagram.

OUTPUT:

Rauzy diagram – the Rauzy diagram of the permutation

```
sage: p = iet.GeneralizedPermutation('a b c b', 'c d d a')
sage: d = p.rauzy_diagram()
```

```
sage: p in d
True
```

For more information, try help(iet.RauzyDiagram)

reduced()

Returns the associated reduced quadratic permutations.

OUTPUT:

permutation – the underlying reduced permutation

EXAMPLES:

```
sage: p = iet.GeneralizedPermutation('a a','b b c c')
sage: q = p.reduced()
sage: q
a a
b b c c
sage: p.rauzy_move(0).reduced() == q.rauzy_move(0)
True
```

right_rauzy_move (winner)

Perform a Rauzy move on the right (the standard one).

INPUT:

• winner - 'top' (or 't' or 0) or 'bottom' (or 'b' or 1)

OUTPUT:

boolean - True if self has a Rauzy move

EXAMPLES:

```
sage: p = iet.GeneralizedPermutation('a a b','b c c')
sage: p.right_rauzy_move(0)
a a b
b c c
sage: p.right_rauzy_move(1)
a a
b b c c
```

```
sage: p = iet.GeneralizedPermutation('a b b','c c a')
sage: p.right_rauzy_move(0)
a a b b
c c
sage: p.right_rauzy_move(1)
a b b
c c a
```

sage.dynamics.interval_exchanges.labelled.LabelledPermutationsIET_iterator (nintervals=None,

irreducible=True,
alphabet=None)

Returns an iterator over labelled permutations.

INPUT:

- nintervals integer or None
- irreducible boolean (default: True)
- alphabet something that should be converted to an alphabet of at least nintervals letters

OUTPUT:

iterator – an iterator over permutations

 $\textbf{Bases: } \textit{sage.dynamics.interval_exchanges.template.RauzyDiagram}$

Template for Rauzy diagrams of labelled permutations.

```
Warning: DO NOT USE
```

class Path (parent, *data)

```
Bases: sage.dynamics.interval_exchanges.template.RauzyDiagram.Path
```

Path in Labelled Rauzy diagram.

dual_substitution()

Returns the substitution of intervals obtained.

OUTPUT:

WordMorphism – the word morphism corresponding to the interval

EXAMPLES:

```
sage: p = iet.Permutation('a b','b a')
sage: r = p.rauzy_diagram()
sage: p0 = r.path(p,0)
sage: s0 = p0.interval_substitution()
sage: s0
WordMorphism: a->a, b->ba
sage: p1 = r.path(p,1)
sage: s1 = p1.interval_substitution()
sage: s1
WordMorphism: a->ab, b->b
sage: (p0 + p1).interval_substitution() == s1 * s0
True
sage: (p1 + p0).interval_substitution() == s0 * s1
```

interval_substitution()

Returns the substitution of intervals obtained.

OUTPUT:

WordMorphism – the word morphism corresponding to the interval

```
sage: p = iet.Permutation('a b','b a')
sage: r = p.rauzy_diagram()
sage: p0 = r.path(p,0)
sage: s0 = p0.interval_substitution()
sage: s0
WordMorphism: a->a, b->ba
sage: p1 = r.path(p,1)
sage: s1 = p1.interval_substitution()
sage: s1
WordMorphism: a->ab, b->b
sage: (p0 + p1).interval_substitution() == s1 * s0
True
sage: (p1 + p0).interval_substitution() == s0 * s1
```

is full()

Tests the fullness.

A path is full if all intervals win at least one time.

OUTPUT:

boolean - True if the path is full and False else

EXAMPLES:

```
sage: p = iet.Permutation('a b c','c b a')
sage: r = p.rauzy_diagram()
sage: g0 = r.path(p,'t','b','t')
sage: g1 = r.path(p,'b','t','b')
sage: g0.is_full()
False
sage: g1.is_full()
False
sage: (g0 + g1).is_full()
True
sage: (g1 + g0).is_full()
```

matrix()

Returns the matrix associated to a path.

The matrix associated to a Rauzy induction, is the linear application that allows to recover the lengths of self from the lengths of the induced.

OUTPUT:

matrix - a square matrix of integers

```
sage: p = iet.Permutation('a1 a2','a2 a1')
sage: d = p.rauzy_diagram()
sage: g = d.path(p,'top')
sage: g.matrix()
[1 0]
[1 1]
sage: g = d.path(p,'bottom')
sage: g.matrix()
[1 1]
[0 1]
```

```
sage: p = iet.Permutation('a b c','c b a')
sage: d = p.rauzy_diagram()
sage: g = d.path(p)
sage: g.matrix() == identity_matrix(3)
True
sage: g = d.path(p,'top')
sage: g.matrix()
[1 0 0]
[0 1 0]
[1 0 1]
sage: g = d.path(p,'bottom')
sage: g.matrix()
[1 0 1]
[0 1 0]
[0 1 0]
[0 1 0]
```

orbit_substitution()

Return the substitution on the orbit of the left extremity.

OUTPUT:

WordMorphism – the word morphism corresponding to the orbit

EXAMPLES:

```
sage: p = iet.Permutation('a b','b a')
sage: d = p.rauzy_diagram()
sage: g0 = d.path(p,'top')
sage: s0 = g0.orbit_substitution()
sage: s0
WordMorphism: a->ab, b->b
sage: g1 = d.path(p,'bottom')
sage: s1 = g1.orbit_substitution()
sage: s1
WordMorphism: a->a, b->ab
sage: (g0 + g1).orbit_substitution() == s0 * s1
True
sage: (g1 + g0).orbit_substitution() == s1 * s0
True
```

substitution()

Return the substitution on the orbit of the left extremity.

OUTPUT:

WordMorphism – the word morphism corresponding to the orbit

```
sage: p = iet.Permutation('a b','b a')
sage: d = p.rauzy_diagram()
sage: g0 = d.path(p,'top')
sage: s0 = g0.orbit_substitution()
sage: s0
WordMorphism: a->ab, b->b
sage: g1 = d.path(p,'bottom')
sage: s1 = g1.orbit_substitution()
sage: s1
```

```
WordMorphism: a->a, b->ab
sage: (g0 + g1).orbit_substitution() == s0 * s1
True
sage: (g1 + g0).orbit_substitution() == s1 * s0
True
```

edge_to_interval_substitution(p=None, edge_type=None)

Returns the interval substitution associated to an edge

OUTPUT:

WordMorphism – the WordMorphism corresponding to the edge

EXAMPLES:

```
sage: p = iet.Permutation('a b c','c b a')
sage: r = p.rauzy_diagram()
sage: r.edge_to_interval_substitution(None, None)
WordMorphism: a->a, b->b, c->c
sage: r.edge_to_interval_substitution(p,0)
WordMorphism: a->a, b->b, c->ca
sage: r.edge_to_interval_substitution(p,1)
WordMorphism: a->ac, b->b, c->c
```

edge_to_orbit_substitution (p=None, edge_type=None)

Returns the interval substitution associated to an edge

OUTPUT:

WordMorphism – the word morphism corresponding to the edge

EXAMPLES:

```
sage: p = iet.Permutation('a b c','c b a')
sage: r = p.rauzy_diagram()
sage: r.edge_to_orbit_substitution(None, None)
WordMorphism: a->a, b->b, c->c
sage: r.edge_to_orbit_substitution(p,0)
WordMorphism: a->ac, b->b, c->c
sage: r.edge_to_orbit_substitution(p,1)
WordMorphism: a->a, b->b, c->ac
```

full_loop_iterator (start=None, max_length=1)

Returns an iterator over all full path starting at start.

INPUT:

- start the start point
- max_length a limit on the length of the paths

OUTPUT:

iterator – iterator over full loops

```
sage: p = iet.Permutation('a b','b a')
sage: r = p.rauzy_diagram()
sage: for g in r.full_loop_iterator(p,2):
...:     print(g.matrix())
...:     print("*****")
```

```
[1 1]
[1 2]
*****
[2 1]
[1 1]
*****
```

full_nloop_iterator (start=None, length=1)

Returns an iterator over all full loops of given length.

INPUT:

- start the initial permutation
- length the length to consider

OUTPUT:

iterator – an iterator over the full loops of given length

EXAMPLES:

```
sage: p = iet.Permutation('a b','b a')
sage: d = p.rauzy_diagram()
sage: for g in d.full_nloop_iterator(p,2):
...:     print(g.matrix())
...:     print("*****")
[1 1]
[1 2]
******
[2 1]
[1 1]
*****
```

4.3 Reduced permutations

Warning: This module is deprecated. You are advised to install and use the surface_dynamics package instead available at https://pypi.python.org/pypi/surface_dynamics/

A reduced (generalized) permutation is better suited to study strata of Abelian (or quadratic) holomorphic forms on Riemann surfaces. The Rauzy diagram is an invariant of such a component. Corentin Boissy proved the identification of Rauzy diagrams with connected components of stratas. But the geometry of the diagram and the relation with the strata is not yet totally understood.

AUTHORS:

• Vincent Delecroix (2000-09-29): initial version

Bases: sage.dynamics.interval_exchanges.reduced.ReducedPermutation

Flipped Reduced Permutation.

Warning: Internal class! Do not use directly!

INPUT:

- intervals a list of two lists
- flips the flipped letters
- alphabet an alphabet

right_rauzy_move (winner)

Performs a Rauzy move on the right.

EXAMPLES:

```
sage: p = iet.Permutation('a b c','c b a',reduced=True,flips='c')
sage: p.right_rauzy_move('top')
-a b -c
-a -c b
```

 $\textbf{class} \texttt{ sage.dynamics.interval_exchanges.reduced.FlippedReducedPermutationIET} (\textit{intervals=None}, \textit{class}) \\$

flips=None, alpha-

 $\bet=None) \\ \textbf{Bases:} \quad sage.dynamics.interval_exchanges.reduced.FlippedReducedPermutation,} \\ sage.dynamics.interval_exchanges.template.FlippedPermutationIET, \quad sage. \\ dynamics.interval_exchanges.reduced.ReducedPermutationIET$

Flipped Reduced Permutation from iet

EXAMPLES

```
sage: p = iet.Permutation('a b c', 'c b a', flips=['a'], reduced=True)
sage: p.rauzy_move(1)
-a -b c
-a c -b
```

list (flips=False)

Returns a list representation of self.

INPUT:

• flips - boolean (default: False) if True the output contains 2-uple of (label, flip)

EXAMPLES:

```
sage: p = iet.Permutation('a b','b a',reduced=True,flips='b')
sage: p.list(flips=True)
[[('a', 1), ('b', -1)], [('b', -1), ('a', 1)]]
sage: p.list(flips=False)
[['a', 'b'], ['b', 'a']]
sage: p.alphabet([0,1])
sage: p.list(flips=True)
[[(0, 1), (1, -1)], [(1, -1), (0, 1)]]
sage: p.list(flips=False)
[[0, 1], [1, 0]]
```

One can recover the initial permutation from this list:

```
sage: p = iet.Permutation('a b','b a',reduced=True,flips='a')
sage: iet.Permutation(p.list(), flips=p.flips(), reduced=True) == p
True
```

rauzy_diagram(**kargs)

Returns the associated Rauzy diagram.

EXAMPLES:

```
sage: p = iet.Permutation('a b','b a',reduced=True,flips='a')
sage: r = p.rauzy_diagram()
sage: p in r
True
```

class sage.dynamics.interval_exchanges.reduced.FlippedReducedPermutationLI (intervals=None,

flips=None, al-

phabet=None)

Bases: sage.dynamics.interval_exchanges.reduced.FlippedReducedPermutation, sage.dynamics.interval_exchanges.template.FlippedPermutationLI, sage.dynamics.interval_exchanges.reduced.ReducedPermutationLI

Flipped Reduced Permutation from li

EXAMPLES:

Creation using the GeneralizedPermutation function:

```
sage: p = iet.GeneralizedPermutation('a a b', 'b c c', reduced=True, flips='a')
```

list (flips=False)

Returns a list representation of self.

INPUT:

• flips - boolean (default: False) return the list with flips

EXAMPLES:

```
sage: p = iet.GeneralizedPermutation('a a','b b',reduced=True,flips='a')
sage: p.list(flips=True)
[[('a', -1), ('a', -1)], [('b', 1), ('b', 1)]]
sage: p.list(flips=False)
[['a', 'a'], ['b', 'b']]

sage: p = iet.GeneralizedPermutation('a a b','b c c',reduced=True,flips='abc')
sage: p.list(flips=True)
[[('a', -1), ('a', -1), ('b', -1)], [('b', -1), ('c', -1), ('c', -1)]]
sage: p.list(flips=False)
[['a', 'a', 'b'], ['b', 'c', 'c']]
```

one can rebuild the permutation from the list:

```
sage: p = iet.GeneralizedPermutation('a a b','b c c',flips='a',reduced=True)
sage: iet.GeneralizedPermutation(p.list(),flips=p.flips(),reduced=True) == p
True
```

rauzy_diagram(**kargs)

Returns the associated Rauzy diagram.

For more explanation and a list of arguments try help(iet.RauzyDiagram)

EXAMPLES:

```
sage: p = iet.GeneralizedPermutation('a a b','c c b',reduced=True)
sage: r = p.rauzy_diagram()
sage: p in r
True
```

class sage.dynamics.interval_exchanges.reduced.FlippedReducedRauzyDiagram(p,

right_induction=True, left_induction=False, left_right_inversion=Fal top_bottom_inversion=Falsymmetric=False)

Bases: sage.dynamics.interval_exchanges.template.FlippedRauzyDiagram, sage.dynamics.interval_exchanges.reduced.ReducedRauzyDiagram

Rauzy diagram of flipped reduced permutations.

Bases: sage.structure.sage_object.SageObject

Template for reduced objects.

```
Warning: Internal class! Do not use directly!
```

INPUT:

- intervals a list of two list of labels
- alphabet (default: None) any object that can be used to initialize an Alphabet or None. In this latter case, the letter of the intervals are used to generate one.

erase_letter(letter)

Erases a letter.

INPUT:

• letter - a letter which is a label of an interval of self

EXAMPLES:

```
sage: p = iet.Permutation('a b c','c b a')
sage: p.erase_letter('a')
b c
c b
```

```
sage: p = iet.GeneralizedPermutation('a b b','c c a')
sage: p.erase_letter('a')
b b
c c
```

left_rauzy_move (winner)

Performs a Rauzy move on the left.

EXAMPLES:

```
sage: p = iet.Permutation('a b c','c b a',reduced=True)
sage: p.left_rauzy_move(0)
a b c
b c a
sage: p.right_rauzy_move(1)
a b c
b c a
```

```
sage: p = iet.GeneralizedPermutation('a a','b b c c',reduced=True)
sage: p.left_rauzy_move(0)
a a b
b c c
```

length (interval=None)

Returns the 2-uple of lengths.

p.length() is identical to (p.length_top(), p.length_bottom()) If an interval is specified, it returns the length of the specified interval.

INPUT:

• interval - None, 'top' (or 't' or 0) or 'bottom' (or 'b' or 1)

OUTPUT:

integer or 2-uple of integers - the corresponding lengths

EXAMPLES:

```
sage: p = iet.Permutation('a b c','c b a')
sage: p.length()
(3, 3)
sage: p = iet.GeneralizedPermutation('a a b','c d c b d')
sage: p.length()
(3, 5)
```

length bottom()

Returns the number of intervals in the bottom segment.

OUTPUT:

integer - the length of the bottom segment

EXAMPLES:

```
sage: p = iet.Permutation('a b c','c b a')
sage: p.length_bottom()
3
sage: p = iet.GeneralizedPermutation('a a b','c d c b d')
sage: p.length_bottom()
5
```

length_top()

Returns the number of intervals in the top segment.

OUTPUT:

integer – the length of the top segment

```
sage: p = iet.Permutation('a b c','c b a')
sage: p.length_top()
3
sage: p = iet.GeneralizedPermutation('a a b','c d c b d')
sage: p.length_top()
3
sage: p = iet.GeneralizedPermutation('a b c b d c d', 'e a e')
sage: p.length_top()
7
```

right_rauzy_move (winner)

Performs a Rauzy move on the right.

EXAMPLES:

```
sage: p = iet.Permutation('a b c','c b a',reduced=True)
sage: p.right_rauzy_move(0)
a b c
c a b
sage: p.right_rauzy_move(1)
a b c
b c a
```

```
sage: p = iet.GeneralizedPermutation('a a','b b c c',reduced=True)
sage: p.right_rauzy_move(0)
a b b
c c a
```

 ${\bf class} \ \, {\bf sage.dynamics.interval_exchanges.reduced.ReducedPermutationIET} \, ({\it intervals=None}, \\ alpha-$

bet=None)

Bases: sage.dynamics.interval_exchanges.reduced.ReducedPermutation, sage.dynamics.interval_exchanges.template.PermutationIET

Reduced permutation from iet

Permutation from iet without numerotation of intervals. For initialization, you should use GeneralizedPermutation which is the class factory for all permutation types.

EXAMPLES:

Equality testing (no equality of letters but just of ordering):

```
sage: p = iet.Permutation('a b c', 'c b a', reduced = True)
sage: q = iet.Permutation('p q r', 'r q p', reduced = True)
sage: p == q
True
```

Reducibility testing:

```
sage: p = iet.Permutation('a b c', 'c b a', reduced = True)
sage: p.is_irreducible()
True
```

```
sage: q = iet.Permutation('a b c d', 'b a d c', reduced = True)
sage: q.is_irreducible()
False
```

Rauzy movability and Rauzy move:

```
sage: p = iet.Permutation('a b c', 'c b a', reduced = True)
sage: p.has_rauzy_move(1)
True
sage: p.rauzy_move(1)
a b c
b c a
```

Rauzy diagrams:

```
sage: p = iet.Permutation('a b c d', 'd a b c')
sage: p_red = iet.Permutation('a b c d', 'd a b c', reduced = True)
sage: d = p.rauzy_diagram()
sage: d_red = p_red.rauzy_diagram()
sage: p.rauzy_move(0) in d
True
sage: d.cardinality(), d_red.cardinality()
(12, 6)
```

has_rauzy_move (winner, side='right')

Tests if the permutation is rauzy_movable on the left.

EXAMPLES:

```
sage: p = iet.Permutation('a b c','a c b',reduced=True)
sage: p.has_rauzy_move(0,'right')
True
sage: p.has_rauzy_move(0,'left')
False
sage: p.has_rauzy_move(1,'right')
True
sage: p.has_rauzy_move(1,'left')
False
```

```
sage: p = iet.Permutation('a b c d','c a b d',reduced=True)
sage: p.has_rauzy_move(0,'right')
False
sage: p.has_rauzy_move(0,'left')
True
sage: p.has_rauzy_move(1,'right')
False
sage: p.has_rauzy_move(1,'left')
True
```

is_identity()

Returns True if self is the identity.

EXAMPLES:

```
sage: iet.Permutation("a b","a b",reduced=True).is_identity()
True
sage: iet.Permutation("a b","b a",reduced=True).is_identity()
False
```

list()

Returns a list of two list that represents the permutation.

```
sage: p = iet.GeneralizedPermutation('a b','b a',reduced=True)
sage: p.list() == [p[0], p[1]]
True
sage: p.list() == [['a', 'b'], ['b', 'a']]
True
```

```
sage: p = iet.GeneralizedPermutation('a b c', 'b c a', reduced=True)
sage: iet.GeneralizedPermutation(p.list(), reduced=True) == p
True
```

rauzy_diagram(**kargs)

Returns the associated Rauzy diagram.

OUTPUT:

A Rauzy diagram

EXAMPLES:

```
sage: p = iet.Permutation('a b c d', 'd a b c',reduced=True)
sage: d = p.rauzy_diagram()
sage: p.rauzy_move(0) in d
True
sage: p.rauzy_move(1) in d
True
```

For more information, try help RauzyDiagram

rauzy_move_relabel (winner, side='right')

Returns the relabelization obtained from this move.

EXAMPLES:

```
sage: p = iet.Permutation('a b c d','d c b a')
sage: q = p.reduced()
sage: p_t = p.rauzy_move('t')
sage: q_t = q.rauzy_move('t')
sage: s_t = q.rauzy_move_relabel('t')
sage: s_t
WordMorphism: a->a, b->b, c->c, d->d
sage: list(map(s_t, p_t[0])) == list(map(Word, q_t[0]))
True
sage: list(map(s_t, p_t[1])) == list(map(Word, q_t[1]))
True
sage: p_b = p.rauzy_move('b')
sage: q_b = q.rauzy_move('b')
sage: s_b = q.rauzy_move_relabel('b')
sage: s_b
WordMorphism: a->a, b->d, c->b, d->c
sage: list(map(s_b, q_b[0])) == list(map(Word, p_b[0]))
sage: list(map(s_b, q_b[1])) == list(map(Word, p_b[1]))
True
```

bet=None)

Bases: sage.dynamics.interval_exchanges.reduced.ReducedPermutation, sage.dynamics.interval_exchanges.template.PermutationLI

Reduced quadratic (or generalized) permutation.

EXAMPLES:

Reducibility testing:

```
sage: p = iet.GeneralizedPermutation('a b b', 'c c a', reduced = True)
sage: p.is_irreducible()
True
```

```
sage: p = iet.GeneralizedPermutation('a b c a', 'b d d c', reduced = True)
sage: p.is_irreducible()
False
sage: test, decomposition = p.is_irreducible(return_decomposition = True)
sage: test
False
sage: decomposition
(['a'], ['c', 'a'], [], ['c'])
```

Rauzy movability and Rauzy move:

```
sage: p = iet.GeneralizedPermutation('a b b', 'c c a', reduced = True)
sage: p.has_rauzy_move(0)
True
sage: p.rauzy_move(0)
a a b b
c c
sage: p.rauzy_move(0).has_rauzy_move(0)
False
sage: p.rauzy_move(1)
a b b
c c a
```

Rauzy diagrams:

```
sage: p_red = iet.GeneralizedPermutation('a b b', 'c c a', reduced = True)
sage: d_red = p_red.rauzy_diagram()
sage: d_red.cardinality()
4
```

list()

The permutations as a list of two lists.

EXAMPLES:

```
sage: p = iet.GeneralizedPermutation('a b b', 'c c a', reduced = True)
sage: list(p)
[['a', 'b', 'b'], ['c', 'c', 'a']]
```

rauzy_diagram(**kargs)

Returns the associated Rauzy diagram.

The Rauzy diagram of a permutation corresponds to all permutations that we could obtain from this one by Rauzy move. The set obtained is a labelled Graph. The label of vertices being 0 or 1 depending on the type.

OUTPUT:

Rauzy diagram – the graph of permutations obtained by rauzy induction

EXAMPLES:

```
sage: p = iet.Permutation('a b c d', 'd a b c')
sage: d = p.rauzy_diagram()
```

sage.dynamics.interval_exchanges.reduced.ReducedPermutationsIET_iterator (nintervals=None, ir-reducible=True, alphabet=None)

Returns an iterator over reduced permutations

INPUT:

- nintervals integer or None
- irreducible boolean
- alphabet something that should be converted to an alphabet of at least nintervals letters

class sage.dynamics.interval_exchanges.reduced.ReducedRauzyDiagram (p,

right_induction=True, left_induction=False, left_right_inversion=False, top_bottom_inversion=False, symmetric=False)

Bases: sage.dynamics.interval_exchanges.template.RauzyDiagram

Rauzy diagram of reduced permutations

sage.dynamics.interval_exchanges.reduced.alphabetized_atwin (twin, alphabet)
Alphabetization of a twin of iet.

```
sage: twin = [[0,1],[0,1]]
sage: alphabet = Alphabet("ab")
sage: alphabetized_atwin(twin, alphabet)
[['a', 'b'], ['a', 'b']]
```

```
sage: twin = [[1,0],[1,0]]
sage: alphabet = Alphabet([0,1])
sage: alphabetized_atwin(twin, alphabet)
[[0, 1], [1, 0]]
```

```
sage: twin = [[1,2,3,0],[3,0,1,2]]
sage: alphabet = Alphabet("abcd")
sage: alphabetized_atwin(twin,alphabet)
[['a', 'b', 'c', 'd'], ['d', 'a', 'b', 'c']]
```

sage.dynamics.interval_exchanges.reduced.alphabetized_qtwin (twin, alphabet)
Alphabetization of a qtwin.

```
sage: twin = [[(1,0),(1,1)],[(0,0),(0,1)]]
sage: alphabet = Alphabet("ab")
sage: alphabetized_qtwin(twin,alphabet)
[['a', 'b'], ['a', 'b']]
```

```
sage: twin = [[(1,1), (1,0)],[(0,1), (0,0)]]
sage: alphabet=Alphabet("AB")
sage: alphabetized_qtwin(twin,alphabet)
[['A', 'B'], ['B', 'A']]
sage: alphabet=Alphabet("BA")
sage: alphabetized_qtwin(twin,alphabet)
[['B', 'A'], ['A', 'B']]
```

```
sage: twin = [[(0,1),(0,0)],[(1,1),(1,0)]]
sage: alphabet=Alphabet("ab")
sage: alphabetized_qtwin(twin,alphabet)
[['a', 'a'], ['b', 'b']]
```

```
sage: twin = [[(0,2),(1,1),(0,0)],[(1,2),(0,1),(1,0)]]
sage: alphabet=Alphabet("abc")
sage: alphabetized_qtwin(twin,alphabet)
[['a', 'b', 'a'], ['c', 'b', 'c']]
```

sage.dynamics.interval_exchanges.reduced.labelize_flip(couple)
Returns a string from a 2-uple couple of the form (name, flip).

4.4 Permutations template

Warning: This module is deprecated. You are advised to install and use the surface_dynamics package instead available at https://pypi.python.org/pypi/surface_dynamics/

This file define high level operations on permutations (alphabet, the different rauzy induction, ...) shared by reduced and labeled permutations.

AUTHORS:

• Vincent Delecroix (2008-12-20): initial version

Todo:

- construct as options different string representations for a permutation
 - the two intervals: str
 - the two intervals on one line: str_one_line
 - the separatrix diagram: str_separatrix_diagram
 - twin[0] and twin[1] for reduced permutation
 - nothing (useful for Rauzy diagram)

```
class sage.dynamics.interval_exchanges.template.FlippedPermutation
    Bases: sage.dynamics.interval_exchanges.template.Permutation
```

Template for flipped generalized permutations.

Warning: Internal class! Do not use directly!

AUTHORS:

• Vincent Delecroix (2008-12-20): initial version

```
str (sep='\n')
```

String representation.

class sage.dynamics.interval_exchanges.template.FlippedPermutationIET

Bases: sage.dynamics.interval_exchanges.template.FlippedPermutation, sage.dynamics.interval_exchanges.template.PermutationIET

Template for flipped Abelian permutations.

Warning: Internal class! Do not use directly!

AUTHORS:

• Vincent Delecroix (2008-12-20): initial version

flips()

Returns the list of flips.

EXAMPLES:

```
sage: p = iet.Permutation('a b c','c b a',flips='ac')
sage: p.flips()
['a', 'c']
```

class sage.dynamics.interval_exchanges.template.FlippedPermutationLI

Bases: sage.dynamics.interval_exchanges.template.FlippedPermutation, sage.dynamics.interval exchanges.template.PermutationLI

Template for flipped quadratic permutations.

Warning: Internal class! Do not use directly!

AUTHORS:

• Vincent Delecroix (2008-12-20): initial version

flips

Returns the list of flipped intervals.

```
sage: p = iet.GeneralizedPermutation('a a','b b',flips='a')
sage: p.flips()
['a']
sage: p = iet.GeneralizedPermutation('a a','b b',flips='b',reduced=True)
sage: p.flips()
['b']
```

```
class sage.dynamics.interval_exchanges.template.FlippedRauzyDiagram (p,
```

right_induction=True, left_induction=False, left_right_inversion=False, top_bottom_inversion=False, symmetric=False)

Bases: sage.dynamics.interval_exchanges.template.RauzyDiagram

Template for flipped Rauzy diagrams.

```
Warning: Internal class! Do not use directly!
```

AUTHORS:

• Vincent Delecroix (2009-09-29): initial version

```
complete(p, reducible=False)
```

Completion of the Rauzy diagram

Add all successors of p for defined operations in edge_types. Could be used for generating non (strongly) connected Rauzy diagrams. Sometimes, for flipped permutations, the maximal connected graph in all permutations is not strongly connected. Finding such components needs to call most than once the .complete() method.

INPUT:

- p a permutation
- reducible put or not reducible permutations

EXAMPLES:

```
sage: p = iet.Permutation('a b c','c b a',flips='a')
sage: d = p.rauzy_diagram()
sage: d
Rauzy diagram with 3 permutations
sage: p = iet.Permutation('a b c','c b a',flips='b')
sage: d.complete(p)
sage: d
Rauzy diagram with 8 permutations
sage: p = iet.Permutation('a b c','c b a',flips='a')
sage: d.complete(p)
sage: d
Rauzy diagram with 8 permutations
```

```
class sage.dynamics.interval_exchanges.template.Permutation
    Bases: sage.structure.sage_object.SageObject
```

Template for all permutations.

```
Warning: Internal class! Do not use directly!
```

This class implement generic algorithm (stratum, connected component, ...) and unfies all its children.

```
alphabet (data=None)
```

Manages the alphabet of self.

If there is no argument, the method returns the alphabet used. If the argument could be converted to an alphabet, this alphabet will be used.

INPUT:

• data - None or something that could be converted to an alphabet

OUTPUT:

- either None or the current alphabet

EXAMPLES:

```
sage: p = iet.Permutation('a b','a b')
sage: p.alphabet([0,1])
True
sage: p
0 1
0 1
sage: p.alphabet("cd")
sage: p.alphabet("cd")
True
sage: p.alphabet() == Alphabet(['c','d'])
True
sage: p
c d
c d
```

has_rauzy_move (winner='top', side=None)

Tests the legality of a Rauzy move.

INPUT:

- winner 'top' or 'bottom' corresponding to the interval
- side 'left' or 'right' (default)

OUTPUT:

– a boolean

```
sage: p = iet.Permutation('a b','a b')
sage: p.has_rauzy_move('top','right')
False
sage: p.has_rauzy_move('bottom','right')
False
sage: p.has_rauzy_move('top','left')
False
sage: p.has_rauzy_move('bottom','left')
False
```

```
sage: p = iet.Permutation('a b c','b a c')
sage: p.has_rauzy_move('top','right')
False
sage: p.has_rauzy_move('bottom', 'right')
False
sage: p.has_rauzy_move('top','left')
True
sage: p.has_rauzy_move('bottom','left')
True
```

```
sage: p = iet.Permutation('a b','b a')
sage: p.has_rauzy_move('top','right')
True
sage: p.has_rauzy_move('bottom','right')
True
sage: p.has_rauzy_move('top','left')
True
sage: p.has_rauzy_move('bottom','left')
True
```

horizontal_inverse()

Returns the top-bottom inverse.

You can use also use the shorter .tb_inverse().

OUTPUT:

- a permutation

EXAMPLES:

```
sage: p = iet.Permutation('a b','b a')
sage: p.top_bottom_inverse()
b a
a b
sage: p = iet.Permutation('a b','b a',reduced=True)
sage: p.top_bottom_inverse() == p
True
```

```
sage: p = iet.Permutation('a b c d','c d a b')
sage: p.top_bottom_inverse()
c d a b
a b c d
```

left_right_inverse()

Returns the left-right inverse.

You can also use the shorter .lr inverse()

OUTPUT:

- a permutation

```
sage: p = iet.Permutation('a b c','c a b')
sage: p.left_right_inverse()
c b a
b a c
sage: p = iet.Permutation('a b c d','c d a b')
sage: p.left_right_inverse()
d c b a
b a d c
```

```
sage: p = iet.GeneralizedPermutation('a a','b b c c')
sage: p.left_right_inverse()
a a
c c b b
```

```
sage: p = iet.Permutation('a b c','c b a',reduced=True)
sage: p.left_right_inverse() == p
True
sage: p = iet.Permutation('a b c','c a b',reduced=True)
sage: q = p.left_right_inverse()
sage: q == p
False
sage: q
a b c
b c a
```

```
sage: p = iet.GeneralizedPermutation('a a','b b c c',reduced=True)
sage: p.left_right_inverse() == p
True
sage: p = iet.GeneralizedPermutation('a b b','c c a',reduced=True)
sage: q = p.left_right_inverse()
sage: q == p
False
sage: q
a a b
b c c
```

letters()

Returns the list of letters of the alphabet used for representation.

The letters used are not necessarily the whole alphabet (for example if the alphabet is infinite).

OUTPUT:

- a list of labels

EXAMPLES:

```
sage: p = iet.Permutation([1,2],[2,1])
sage: p.alphabet(Alphabet(name="NN"))
sage: p
0 1
1 0
sage: p.letters()
[0, 1]
```

lr_inverse()

Returns the left-right inverse.

You can also use the shorter .lr_inverse()

OUTPUT:

- a permutation

```
sage: p = iet.Permutation('a b c','c a b')
sage: p.left_right_inverse()
c b a
b a c
sage: p = iet.Permutation('a b c d','c d a b')
sage: p.left_right_inverse()
d c b a
b a d c
```

```
sage: p = iet.GeneralizedPermutation('a a','b b c c')
sage: p.left_right_inverse()
a a
c c b b
```

```
sage: p = iet.Permutation('a b c','c b a',reduced=True)
sage: p.left_right_inverse() == p
True
sage: p = iet.Permutation('a b c','c a b',reduced=True)
sage: q = p.left_right_inverse()
sage: q == p
False
sage: q
a b c
b c a
```

```
sage: p = iet.GeneralizedPermutation('a a','b b c c',reduced=True)
sage: p.left_right_inverse() == p
True
sage: p = iet.GeneralizedPermutation('a b b','c c a',reduced=True)
sage: q = p.left_right_inverse()
sage: q == p
False
sage: q
a a b
b c c
```

rauzy_move (winner, side='right', iteration=1)

Returns the permutation after a Rauzy move.

INPUT:

- winner 'top' or 'bottom' interval
- side 'right' or 'left' (default: 'right') corresponding to the side on which the Rauzy move must be performed.
- iteration a non negative integer

OUTPUT:

• a permutation

```
sage: p = iet.Permutation('a b c','c b a')
sage: p.rauzy_move(winner=0, side='right')
a b c
c a b
sage: p.rauzy_move(winner=1, side='right')
a c b
c b a
sage: p.rauzy_move(winner=0, side='left')
a b c
b c a
sage: p.rauzy_move(winner=1, side='left')
b a c
c b a
```

$str(sep='\n')$

A string representation of the generalized permutation.

INPUT:

• sep - (default: 'n') a separator for the two intervals

OUTPUT:

string – the string that represents the permutation

EXAMPLES:

For permutations of iet:

```
sage: p = iet.Permutation('a b c','c b a')
sage: p.str()
'a b c\nc b a'
sage: p.str(sep=' | ')
'a b c | c b a'
```

..the permutation can be rebuilt from the standard string:

```
sage: p == iet.Permutation(p.str())
True
```

For permutations of li:

```
sage: p = iet.GeneralizedPermutation('a b b','c c a')
doctest:warning
DeprecationWarning: GeneralizedPermutation is deprecated and will be removed_
→from Sage.
You are advised to install the surface_dynamics package via:
sage -pip install surface_dynamics
If you do not have write access to the Sage installation you can
alternatively do
sage -pip install surface_dynamics --user
The package surface_dynamics subsumes all flat surface related
computation that are currently available in Sage. See more
information at
http://www.labri.fr/perso/vdelecro/surface-dynamics/latest/
See http://trac.sagemath.org/20695 for details.
sage: p.str()
'a b b\nc c a'
sage: p.str(sep=' | ')
'abb|cca'
```

..the generalized permutation can be rebuilt from the standard string:

```
sage: p == iet.GeneralizedPermutation(p.str())
True
```

symmetric()

Returns the symmetric permutation.

The symmetric permutation is the composition of the top-bottom inversion and the left-right inversion (which are geometrically orientation reversing).

OUTPUT:

- a permutation

```
sage: p = iet.Permutation("a b c","c b a")
sage: p.symmetric()
a b c
c b a
sage: q = iet.Permutation("a b c d","b d a c")
sage: q.symmetric()
c a d b
d c b a
```

```
sage: p = iet.Permutation('a b c d','c a d b')
sage: q = p.symmetric()
sage: q1 = p.tb_inverse().lr_inverse()
sage: q2 = p.lr_inverse().tb_inverse()
sage: q == q1
True
sage: q == q2
True
```

```
sage: p = iet.GeneralizedPermutation('a a b','b c c',reduced=True)
sage: q = p.symmetric()
sage: q1 = p.tb_inverse().lr_inverse()
sage: q2 = p.lr_inverse().tb_inverse()
sage: q == q1
True
sage: q == q2
True
```

```
sage: p = iet.GeneralizedPermutation('a a b','b c c',reduced=True,flips='a')
sage: q = p.symmetric()
sage: q1 = p.tb_inverse().lr_inverse()
sage: q2 = p.lr_inverse().tb_inverse()
sage: q == q1
True
sage: q == q2
True
```

tb_inverse()

Returns the top-bottom inverse.

You can use also use the shorter .tb_inverse().

OUTPUT:

- a permutation

```
sage: p = iet.Permutation('a b','b a')
sage: p.top_bottom_inverse()
b a
a b
sage: p = iet.Permutation('a b','b a',reduced=True)
sage: p.top_bottom_inverse() == p
True
```

```
sage: p = iet.Permutation('a b c d','c d a b')
sage: p.top_bottom_inverse()
```

```
c d a b a b c d
```

top_bottom_inverse()

Returns the top-bottom inverse.

You can use also use the shorter .tb_inverse().

OUTPUT:

- a permutation

EXAMPLES:

```
sage: p = iet.Permutation('a b','b a')
sage: p.top_bottom_inverse()
b a
a b
sage: p = iet.Permutation('a b','b a',reduced=True)
sage: p.top_bottom_inverse() == p
True
```

```
sage: p = iet.Permutation('a b c d','c d a b')
sage: p.top_bottom_inverse()
c d a b
a b c d
```

vertical_inverse()

Returns the left-right inverse.

You can also use the shorter .lr_inverse()

OUTPUT:

- a permutation

```
sage: p = iet.Permutation('a b c','c a b')
sage: p.left_right_inverse()
c b a
b a c
sage: p = iet.Permutation('a b c d','c d a b')
sage: p.left_right_inverse()
d c b a
b a d c
```

```
sage: p = iet.GeneralizedPermutation('a a','b b c c')
sage: p.left_right_inverse()
a a
c c b b
```

```
sage: p = iet.Permutation('a b c','c b a',reduced=True)
sage: p.left_right_inverse() == p
True
sage: p = iet.Permutation('a b c','c a b',reduced=True)
sage: q = p.left_right_inverse()
sage: q == p
False
```

```
sage: q
a b c
b c a
```

```
sage: p = iet.GeneralizedPermutation('a a','b b c c',reduced=True)
sage: p.left_right_inverse() == p
True
sage: p = iet.GeneralizedPermutation('a b b','c c a',reduced=True)
sage: q = p.left_right_inverse()
sage: q == p
False
sage: q
a a b
b c c
```

```
class sage.dynamics.interval_exchanges.template.PermutationIET
    Bases: sage.dynamics.interval_exchanges.template.Permutation
```

Template for permutation from Interval Exchange Transformation.

```
Warning: Internal class! Do not use directly!
```

AUTHOR:

• Vincent Delecroix (2008-12-20): initial version

arf invariant()

Returns the Arf invariant of the suspension of self.

OUTPUT:

integer -0 or 1

EXAMPLES:

Permutations from the odd and even component of H(2,2,2):

```
sage: a = range(10)
sage: b1 = [3,2,4,6,5,7,9,8,1,0]
sage: b0 = [6,5,4,3,2,7,9,8,1,0]
sage: p1 = iet.Permutation(a,b1)
sage: p1.arf_invariant()
1
sage: p0 = iet.Permutation(a,b0)
sage: p0.arf_invariant()
```

Permutations from the odd and even component of H(4,4):

```
sage: a = range(11)
sage: b1 = [3,2,5,4,6,8,7,10,9,1,0]
sage: b0 = [5,4,3,2,6,8,7,10,9,1,0]
sage: p1 = iet.Permutation(a,b1)
sage: p1.arf_invariant()
1
sage: p0 = iet.Permutation(a,b0)
sage: p0.arf_invariant()
```

REFERENCES:

attached_in_degree()

Returns the degree of the singularity at the right of the interval.

OUTPUT:

- a positive integer

EXAMPLES:

```
sage: p1 = iet.Permutation('a b c d e f g','d c g f e b a')
sage: p2 = iet.Permutation('a b c d e f g','e d c g f b a')
sage: p1.attached_in_degree()
1
sage: p2.attached_in_degree()
3
```

attached_out_degree()

Returns the degree of the singularity at the left of the interval.

OUTPUT:

- a positive integer

EXAMPLES:

```
sage: p1 = iet.Permutation('a b c d e f g','d c g f e b a')
sage: p2 = iet.Permutation('a b c d e f g','e d c g f b a')
sage: p1.attached_out_degree()
3
sage: p2.attached_out_degree()
1
```

attached_type()

Return the singularity degree attached on the left and the right.

OUTPUT:

([degre], angle_parity) - if the same singularity is attached on the left and right

([left_degree, right_degree], 0) - the degrees at the left and the right which are different singularitites

EXAMPLES:

With two intervals:

```
sage: p = iet.Permutation('a b','b a')
sage: p.attached_type()
([0], 1)
```

With three intervals:

```
sage: p = iet.Permutation('a b c', 'b c a')
sage: p.attached_type()
([0], 1)

sage: p = iet.Permutation('a b c', 'c a b')
sage: p.attached_type()
([0], 1)
```

```
sage: p = iet.Permutation('a b c','c b a')
sage: p.attached_type()
([0, 0], 0)
```

With four intervals:

```
sage: p = iet.Permutation('1 2 3 4','4 3 2 1')
sage: p.attached_type()
([2], 0)
```

connected_component (marked_separatrix='no')

Returns a connected components of a stratum.

EXAMPLES:

Permutations from the stratum H(6):

```
sage: a = range(8)
sage: b_hyp = [7,6,5,4,3,2,1,0]
sage: b_odd = [3,2,5,4,7,6,1,0]
sage: b_even = [5,4,3,2,7,6,1,0]
sage: p_hyp = iet.Permutation(a, b_hyp)
sage: p_odd = iet.Permutation(a, b_odd)
sage: p_even = iet.Permutation(a, b_even)
sage: p_hyp.connected_component()
H_hyp(6)
sage: p_odd.connected_component()
H_odd(6)
sage: p_even.connected_component()
H_even(6)
```

Permutations from the stratum H(4,4):

```
sage: a = range(11)
sage: b_hyp = [10,9,8,7,6,5,4,3,2,1,0]
sage: b_odd = [3,2,5,4,6,8,7,10,9,1,0]
sage: b_{even} = [5, 4, 3, 2, 6, 8, 7, 10, 9, 1, 0]
sage: p_hyp = iet.Permutation(a,b_hyp)
sage: p_odd = iet.Permutation(a,b_odd)
sage: p_even = iet.Permutation(a,b_even)
sage: p_hyp.stratum() == AbelianStratum(4,4)
True
sage: p_hyp.connected_component()
H_hyp(4, 4)
sage: p_odd.stratum() == AbelianStratum(4,4)
True
sage: p_odd.connected_component()
H_odd(4, 4)
sage: p_even.stratum() == AbelianStratum(4,4)
True
sage: p_even.connected_component()
H_{even(4, 4)}
```

As for stratum you can specify that you want to attach the singularity on the left of the interval using the option marked_separatrix:

```
sage: a = [1,2,3,4,5,6,7,8,9]
sage: b4_odd = [4,3,6,5,7,9,8,2,1]
```

```
sage: b4_{even} = [6, 5, 4, 3, 7, 9, 8, 2, 1]
sage: b2\_odd = [4,3,5,7,6,9,8,2,1]
sage: b2_even = [7,6,5,4,3,9,8,2,1]
sage: p4_odd = iet.Permutation(a,b4_odd)
sage: p4_even = iet.Permutation(a,b4_even)
sage: p2_odd = iet.Permutation(a,b2_odd)
sage: p2_even = iet.Permutation(a,b2_even)
sage: p4_odd.connected_component(marked_separatrix='out')
H_odd^out(4, 2)
sage: p4_even.connected_component (marked_separatrix='out')
H_{even}^{out}(4, 2)
sage: p2_odd.connected_component(marked_separatrix='out')
H_odd^out(2, 4)
sage: p2_even.connected_component (marked_separatrix='out')
H_{even}^{out}(2, 4)
sage: p2_odd.connected_component() == p4_odd.connected_component()
sage: p2_odd.connected_component('out') == p4_odd.connected_component('out')
False
```

cylindric()

Returns a permutation in the Rauzy class such that

```
twin[0][-1] == 0 twin[1][-1] == 0
```

decompose()

Returns the decomposition of self.

OUTPUT:

- a list of permutations

EXAMPLES:

```
sage: p = iet.Permutation('a b c','c b a').decompose()[0]
sage: p
a b c
c b a
```

```
sage: p1,p2,p3 = iet.Permutation('a b c d e','b a c e d').decompose()
sage: p1
a b
b a
sage: p2
c
c
sage: p3
d e
e d
```

erase_marked_points()

Returns a permutation equivalent to self but without marked points.

```
sage: a = iet.Permutation('a b1 b2 c d', 'd c b1 b2 a')
sage: a.erase_marked_points()
a b1 c d
d c b1 a
```

genus()

Returns the genus corresponding to any suspension of the permutation.

OUTPUT:

- a positive integer

EXAMPLES:

```
sage: p = iet.Permutation('a b c', 'c b a')
sage: p.genus()
1
```

```
sage: p = iet.Permutation('a b c d','d c b a')
sage: p.genus()
2
```

REFERENCES: Veech

intersection matrix()

Returns the intersection matrix.

This d * d antisymmetric matrix is given by the rule :

$$m_{ij} = \begin{cases} 1 & i < j \text{ and } \pi(i) > \pi(j) \\ -1 & i > j \text{ and } \pi(i) < \pi(j) \\ 0 & \text{else} \end{cases}$$

OUTPUT:

· a matrix

EXAMPLES:

```
sage: p = iet.Permutation('a b c d','d c b a')
sage: p.intersection_matrix()
[ 0 1 1 1]
[-1 0 1 1]
[-1 -1 0 1]
[-1 -1 0 0]
```

```
sage: p = iet.Permutation('1 2 3 4 5','5 3 2 4 1')
sage: p.intersection_matrix()
[ 0 1 1 1 1]
[-1 0 1 0 1]
[-1 -1 0 0 0 1]
[-1 -1 -1 -1 -1 0]
```

is_cylindric()

Returns True if the permutation is Rauzy_1n.

A permutation is cylindric if 1 and n are exchanged.

```
sage: iet.Permutation('1 2 3','3 2 1').is_cylindric()
True
sage: iet.Permutation('1 2 3','2 1 3').is_cylindric()
False
```

is_hyperelliptic()

Returns True if the permutation is in the class of the symmetric permutations (with eventual marked points).

This is equivalent to say that the suspension lives in an hyperelliptic stratum of Abelian differentials $H_hyp(2g-2)$ or $H_hyp(g-1, g-1)$ with some marked points.

EXAMPLES:

```
sage: iet.Permutation('a b c d','d c b a').is_hyperelliptic()
True
sage: iet.Permutation('0 1 2 3 4 5','5 2 1 4 3 0').is_hyperelliptic()
False
```

REFERENCES:

Gerard Rauzy, "Echanges d'intervalles et transformations induites", Acta Arith. 34, no. 3, 203-212, 1980

M. Kontsevich, A. Zorich "Connected components of the moduli space of Abelian differentials with prescribed singularities" Invent. math. 153, 631-678 (2003)

is_irreducible (return_decomposition=False)

Tests the irreducibility.

An abelian permutation p = (p0,p1) is reducible if: set(p0[:i]) = set(p1[:i]) for an i < len(p0)

OUTPUT:

· a boolean

EXAMPLES:

```
sage: p = iet.Permutation('a b c', 'c b a')
sage: p.is_irreducible()
True

sage: p = iet.Permutation('a b c', 'b a c')
sage: p.is_irreducible()
False
```

order_of_rauzy_action (winner, side=None)

Returns the order of the action of a Rauzy move.

INPUT:

- winner string 'top' or 'bottom'
- side string 'left' or 'right'

OUTPUT:

An integer corresponding to the order of the Rauzy action.

```
sage: p = iet.Permutation('a b c d','d a c b')
sage: p.order_of_rauzy_action('top', 'right')
3
sage: p.order_of_rauzy_action('bottom', 'right')
2
sage: p.order_of_rauzy_action('top', 'left')
1
sage: p.order_of_rauzy_action('bottom', 'left')
3
```

separatrix diagram(side=False)

Returns the separatrix diagram of the permutation.

INPUT:

• side - boolean

OUTPUT:

a list of lists

EXAMPLES:

```
sage: iet.Permutation([0, 1], [1, 0]).separatrix_diagram()
[[(1, 0), (1, 0)]]
```

```
sage: iet.Permutation('a b c d','d c b a').separatrix_diagram()
[[('d', 'a'), 'b', 'c', ('d', 'a'), 'b', 'c']]
```

stratum (marked_separatrix='no')

Returns the strata in which any suspension of this permutation lives.

OUTPUT:

• a stratum of Abelian differentials

EXAMPLES:

```
sage: p = iet.Permutation('a b c', 'c b a')
sage: p.stratum()
doctest:warning
DeprecationWarning: AbelianStratum is deprecated and will be removed from,
You are advised to install the surface_dynamics package via:
   sage -pip install surface_dynamics
If you do not have write access to the Sage installation you can
alternatively do
   sage -pip install surface_dynamics --user
The package surface_dynamics subsumes all flat surface related
computation that are currently available in Sage. See more
information at
   http://www.labri.fr/perso/vdelecro/surface-dynamics/latest/
See http://trac.sagemath.org/20695 for details.
H(0, 0)
sage: p = iet.Permutation('a b c d', 'd a b c')
sage: p.stratum()
H(0, 0, 0)
sage: p = iet.Permutation(range(9), [8,5,2,7,4,1,6,3,0])
sage: p.stratum()
H(1, 1, 1, 1)
```

You can specify that you want to attach the singularity on the left (or on the right) with the option marked_separatrix:

```
sage: a = 'a b c d e f g h i j'
sage: b3 = 'd c g f e j i h b a'
sage: b2 = 'd c e g f j i h b a'
sage: b1 = 'e d c g f h j i b a'
```

```
sage: p3 = iet.Permutation(a, b3)
sage: p3.stratum()
H(3, 2, 1)
sage: p3.stratum(marked_separatrix='out')
H^out(3, 2, 1)
sage: p2 = iet.Permutation(a, b2)
sage: p2.stratum()
H(3, 2, 1)
sage: p2.stratum(marked_separatrix='out')
H^out(2, 3, 1)
sage: p1 = iet.Permutation(a, b1)
sage: p1.stratum()
H(3, 2, 1)
sage: p1.stratum(marked_separatrix='out')
H^out(1, 3, 2)
```

AUTHORS:

• Vincent Delecroix (2008-12-20)

to permutation()

Returns the permutation as an element of the symmetric group.

EXAMPLES:

```
sage: p = iet.Permutation('a b c','c b a')
sage: p.to_permutation()
[3, 2, 1]
```

```
sage: p = Permutation([2,4,1,3])
sage: q = iet.Permutation(p)
sage: q.to_permutation() == p
True
```

class sage.dynamics.interval_exchanges.template.PermutationLI

Bases: sage.dynamics.interval_exchanges.template.Permutation

Template for quadratic permutation.

```
Warning: Internal class! Do not use directly!
```

AUTHOR:

• Vincent Delecroix (2008-12-20): initial version

has_right_rauzy_move(winner)

Test of Rauzy movability (with an eventual specified choice of winner)

A quadratic (or generalized) permutation is rauzy_movable type depending on the possible length of the last interval. It's dependent of the length equation.

INPUT:

• winner - the integer 'top' or 'bottom'

```
sage: p = iet.GeneralizedPermutation('a a','b b')
sage: p.has_right_rauzy_move('top')
False
sage: p.has_right_rauzy_move('bottom')
False
```

```
sage: p = iet.GeneralizedPermutation('a a b','b c c')
sage: p.has_right_rauzy_move('top')
True
sage: p.has_right_rauzy_move('bottom')
True
```

```
sage: p = iet.GeneralizedPermutation('a a','b b c c')
sage: p.has_right_rauzy_move('top')
True
sage: p.has_right_rauzy_move('bottom')
False
```

```
sage: p = iet.GeneralizedPermutation('a a b b','c c')
sage: p.has_right_rauzy_move('top')
False
sage: p.has_right_rauzy_move('bottom')
True
```

is_irreducible (return_decomposition=False)

Test of reducibility

A quadratic (or generalized) permutation is reducible if there exists a decomposition

$$A1uB1|...|B1uA2$$

$$A1uB2|...|B2uA2$$

where no corners is empty, or exactly one corner is empty and it is on the left, or two and they are both on the right or on the left. The definition is due to [BL2008] where they prove that the property of being irreducible is stable under Rauzy induction.

INPUT:

• return_decomposition - boolean (default: False) - if True, and the permutation is reducible, returns also the blocs A1 u B1, B1 u A2, A1 u B2 and B2 u A2 of a decomposition as above.

OUTPUT:

If return_decomposition is True, returns a 2-uple (test,decomposition) where test is the preceding test and decomposition is a 4-uple (A11,A12,A21,A22) where:

```
A11 = A1 u B1 A12 = B1 u A2 A21 = A1 u B2 A22 = B2 u A2
```

```
sage: GP = iet.GeneralizedPermutation

sage: GP('a a','b b').is_irreducible()
False
sage: GP('a a b','b c c').is_irreducible()
True
sage: GP('1 2 3 4 5 1','5 6 6 4 3 2').is_irreducible()
True
```

AUTHORS:

• Vincent Delecroix (2008-12-20)

class sage.dynamics.interval_exchanges.template.RauzyDiagram(p,

right_induction=True, left_induction=False, left_right_inversion=False, top_bottom_inversion=False, symmetric=False)

Bases: sage.structure.sage_object.SageObject

Template for Rauzy diagrams.

```
Warning: Internal class! Do not use directly!
```

AUTHORS:

• Vincent Delecroix (2008-12-20): initial version

```
class Path (parent, *data)
```

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

Path in Rauzy diagram.

A path in a Rauzy diagram corresponds to a subsimplex of the simplex of lengths. This correspondance is obtained via the Rauzy induction. To a idoc IET we can associate a unique path in a Rauzy diagram. This establishes a correspondance between infinite full path in Rauzy diagram and equivalence topologic class of IET.

append(edge type)

Append an edge to the path.

EXAMPLES:

```
sage: p = iet.Permutation('a b c','c b a')
sage: r = p.rauzy_diagram()
sage: g = r.path(p)
sage: g.append('top')
sage: g
Path of length 1 in a Rauzy diagram
sage: g.append('bottom')
sage: g
Path of length 2 in a Rauzy diagram
```

composition (function, composition=None)

Compose an edges function on a path

INPUT:

- path either a Path or a tuple describing a path
- function function must be of the form
- composition the composition function

AUTHOR:

• Vincent Delecroix (2009-09-29)

```
sage: p = iet.Permutation('a b','b a')
sage: r = p.rauzy_diagram()
sage: def f(i,t):
```

```
if t is None: return []
....: return [t]
sage: g = r.path(p)
sage: g.composition(f,list.__add__)
[]
sage: g = r.path(p,0,1)
sage: g.composition(f, list.__add__)
[0, 1]
```

edge_types()

Returns the edge types of the path.

EXAMPLES:

```
sage: p = iet.Permutation('a b c','c b a')
sage: r = p.rauzy_diagram()
sage: g = r.path(p, 0, 1)
sage: g.edge_types()
[0, 1]
```

end()

Returns the last vertex of the path.

EXAMPLES:

```
sage: p = iet.Permutation('a b c','c b a')
sage: r = p.rauzy_diagram()
sage: g1 = r.path(p, 't', 'b', 't')
sage: g1.end() == p
True
sage: g2 = r.path(p, 'b', 't', 'b')
sage: g2.end() == p
True
```

extend (path)

Extends self with another path.

EXAMPLES:

```
sage: p = iet.Permutation('a b c d','d c b a')
sage: r = p.rauzy_diagram()
sage: g1 = r.path(p,'t','t')
sage: g2 = r.path(p.rauzy_move('t',iteration=2),'b','b')
sage: g = r.path(p,'t','t','b','b')
sage: g == g1 + g2
True
sage: g = copy(g1)
sage: g.extend(g2)
sage: g == g1 + g2
True
```

is_loop()

Tests whether the path is a loop (start point = end point).

```
sage: p = iet.Permutation('a b','b a')
sage: r = p.rauzy_diagram()
sage: r.path(p).is_loop()
```

```
True sage: r.path(p,0,1,0,0).is_loop()
True
```

losers()

Returns a list of the loosers on the path.

EXAMPLES:

```
sage: p = iet.Permutation('a b c','c b a')
sage: r = p.rauzy_diagram()
sage: g0 = r.path(p,'t','b','t')
sage: g0.losers()
['a', 'c', 'b']
sage: g1 = r.path(p,'b','t','b')
sage: g1.losers()
['c', 'a', 'b']
```

pop()

Pops the queue of the path

OUTPUT:

a path corresponding to the last edge

EXAMPLES:

```
sage: p = iet.Permutation('a b','b a')
sage: r = p.rauzy_diagram()
sage: g = r.path(p,0,1,0)
sage: g0,g1,g2,g3 = g[0], g[1], g[2], g[3]
sage: g.pop() == r.path(g2,0)
True
sage: g == r.path(g0,0,1)
True
sage: g.pop() == r.path(g1,1)
True
sage: g == r.path(g0,0)
True
sage: g.pop() == r.path(g0,0)
True
sage: g == r.path(g0)
True
sage: g == r.path(g0)
True
sage: g.pop() == r.path(g0)
True
```

right_composition (function, composition=None)

Compose an edges function on a path

INPUT:

- function function must be of the form (indice,type) -> element. Moreover function(None,None) must be an identity element for initialization.
- composition the composition function for the function. * if None (default None)

start()

Returns the first vertex of the path.

```
sage: p = iet.Permutation('a b c','c b a')
sage: r = p.rauzy_diagram()
sage: g = r.path(p, 't', 'b')
sage: g.start() == p
True
```

winners()

Returns the winner list associated to the edge of the path.

EXAMPLES:

```
sage: p = iet.Permutation('a b','b a')
sage: r = p.rauzy_diagram()
sage: r.path(p).winners()
[]
sage: r.path(p,0).winners()
['b']
sage: r.path(p,1).winners()
['a']
```

alphabet (data=None)

cardinality()

Returns the number of permutations in this Rauzy diagram.

OUTPUT

ullet integer - the number of vertices in the diagram

EXAMPLES:

```
sage: r = iet.RauzyDiagram('a b','b a')
sage: r.cardinality()
1
sage: r = iet.RauzyDiagram('a b c','c b a')
sage: r.cardinality()
3
sage: r = iet.RauzyDiagram('a b c d','d c b a')
sage: r.cardinality()
7
```

complete(p)

Completion of the Rauzy diagram.

Add to the Rauzy diagram all permutations that are obtained by successive operations defined by edge_types(). The permutation must be of the same type and the same length as the one used for the creation.

INPUT:

• p - a permutation of Interval exchange transformation

Rauzy diagram is the reunion of all permutations that could be obtained with successive rauzy moves. This function just use the functions __getitem__ and has_rauzy_move and rauzy_move which must be defined for child and their corresponding permutation types.

edge_iterator()

Returns an iterator over the edges of the graph.

```
sage: p = iet.Permutation('a b','b a')
sage: r = p.rauzy_diagram()
sage: for e in r.edge_iterator():
...:     print(e[0].str(sep='/') + ' --> ' + e[1].str(sep='/'))
a b/b a --> a b/b a
a b/b a --> a b/b a
```

edge_to_loser (p=None, edge_type=None)

Return the corresponding loser

edge_to_matrix(p=None, edge_type=None)

Return the corresponding matrix

INPUT:

- p a permutation
- edge_type 0 or 1 corresponding to the type of the edge

OUTPUT:

A matrix

EXAMPLES:

```
sage: p = iet.Permutation('a b c','c b a')
sage: d = p.rauzy_diagram()
sage: d.edge_to_matrix(p,1)
[1 0 1]
[0 1 0]
[0 0 1]
```

edge_to_winner (p=None, edge_type=None)

Return the corresponding winner

edge_types()

Print information about edges.

EXAMPLES:

```
sage: r = iet.RauzyDiagram('a b', 'b a')
sage: r.edge_types()
0: rauzy_move(0, -1)
1: rauzy_move(1, -1)
```

```
sage: r = iet.RauzyDiagram('a b', 'b a', left_induction=True)
sage: r.edge_types()
0: rauzy_move(0, -1)
1: rauzy_move(1, -1)
2: rauzy_move(0, 0)
3: rauzy_move(1, 0)
```

```
sage: r = iet.RauzyDiagram('a b',' b a',symmetric=True)
sage: r.edge_types()
0: rauzy_move(0, -1)
1: rauzy_move(1, -1)
2: symmetric()
```

edge_types_index (data)

Try to convert the data as an edge type.

INPUT:

• data - a string

OUTPUT:

integer

EXAMPLES:

For a standard Rauzy diagram (only right induction) the 0 index corresponds to the 'top' induction and the index 1 corresponds to the 'bottom' one:

```
sage: p = iet.Permutation('a b c','c b a')
sage: r = p.rauzy_diagram()
sage: r.edge_types_index('top')
0
sage: r[p][0] == p.rauzy_move('top')
True
sage: r.edge_types_index('bottom')
1
sage: r[p][1] == p.rauzy_move('bottom')
True
```

The special operations (inversion and symmetry) always appears after the different Rauzy inductions:

```
sage: p = iet.Permutation('a b c','c b a')
sage: r = p.rauzy_diagram(symmetric=True)
sage: r.edge_types_index('symmetric')
2
sage: r[p][2] == p.symmetric()
True
```

This function always try to resolve conflictuous name. If it's impossible a ValueError is raised:

```
sage: p = iet.Permutation('a b c','c b a')
sage: r = p.rauzy_diagram(left_induction=True)
sage: r.edge_types_index('top')
Traceback (most recent call last):
...
ValueError: left and right inductions must be differentiated
sage: r.edge_types_index('top_right')
0
sage: r[p][0] == p.rauzy_move(0)
True
sage: r.edge_types_index('bottom_left')
3
sage: r[p][3] == p.rauzy_move('bottom', 'left')
True
```

```
sage: p = iet.Permutation('a b c','c b a')
sage: r = p.rauzy_diagram(left_right_inversion=True,top_bottom_inversion=True)
sage: r.edge_types_index('inversion')
Traceback (most recent call last):
...
ValueError: left-right and top-bottom inversions must be differentiated
sage: r.edge_types_index('lr_inverse')
2
sage: p.lr_inverse() == r[p][2]
True
```

```
sage: r.edge_types_index('tb_inverse')
3
sage: p.tb_inverse() == r[p][3]
True
```

Short names are accepted:

```
sage: p = iet.Permutation('a b c','c b a')
sage: r = p.rauzy_diagram(right_induction='top',top_bottom_inversion=True)
sage: r.edge_types_index('top_rauzy_move')
0
sage: r.edge_types_index('t')
0
sage: r.edge_types_index('tb')
1
sage: r.edge_types_index('inversion')
1
sage: r.edge_types_index('inverse')
1
sage: r.edge_types_index('inverse')
1
sage: r.edge_types_index('i')
```

edges (labels=True)

Returns a list of the edges.

EXAMPLES:

```
sage: r = iet.RauzyDiagram('a b','b a')
sage: len(r.edges())
2
```

graph()

Returns the Rauzy diagram as a Graph object

The graph returned is more precisely a DiGraph (directed graph) with loops and multiedges allowed.

EXAMPLES:

```
sage: r = iet.RauzyDiagram('a b c','c b a')
sage: r
Rauzy diagram with 3 permutations
sage: r.graph()
Looped multi-digraph on 3 vertices
```

letters()

Returns the letters used by the RauzyDiagram.

```
sage: r = iet.RauzyDiagram('a b','b a')
sage: r.alphabet()
{'a', 'b'}
sage: r.letters()
['a', 'b']
sage: r.alphabet('ABCDEF')
sage: r.alphabet()
{'A', 'B', 'C', 'D', 'E', 'F'}
sage: r.letters()
['A', 'B']
```

path (*data)

Returns a path over this Rauzy diagram.

INPUT:

- initial_vertex the initial vertex (starting point of the path)
- data a sequence of edges

EXAMPLES:

```
sage: p = iet.Permutation('a b c','c b a')
sage: r = p.rauzy_diagram()
sage: g = r.path(p, 'top', 'bottom')
```

vertex_iterator()

Returns an iterator over the vertices

EXAMPLES:

```
sage: r = iet.RauzyDiagram('a b','b a')
sage: for p in r.vertex_iterator(): print(p)
a b
b a
```

```
sage: r = iet.RauzyDiagram('a b c d','d c b a')
sage: from six.moves import filter
sage: r_1n = filter(lambda x: x.is_cylindric(), r)
sage: for p in r_1n: print(p)
a b c d
d c b a
```

vertices()

Returns a list of the vertices.

EXAMPLES:

```
sage: r = iet.RauzyDiagram('a b','b a')
sage: for p in r.vertices(): print(p)
a b
b a
```

sage.dynamics.interval_exchanges.template.interval_conversion (interval=None) Converts the argument in 0 or 1.

INPUT:

• winner - 'top' (or 't' or 0) or bottom (or 'b' or 1)

OUTPUT:

integer -0 or 1

```
sage: from sage.dynamics.interval_exchanges.template import interval_conversion
sage: interval_conversion('top')
0
sage: interval_conversion('t')
0
sage: interval_conversion(0)
0
sage: interval_conversion('bottom')
```

```
sage: interval conversion('b')
     sage: interval_conversion(1)
sage.dynamics.interval_exchanges.template.labelize_flip(couple)
     Returns a string from a 2-uple couple of the form (name, flip).
sage.dynamics.interval_exchanges.template.side_conversion(side=None)
     Converts the argument in 0 or -1.
     INPUT:
       • side - either 'left' (or 'l' or 0) or 'right' (or 'r' or -1)
     OUTPUT:
     integer -0 or -1
     sage: from sage.dynamics.interval_exchanges.template import side_conversion
     sage: side_conversion('left')
     sage: side_conversion('l')
     sage: side_conversion(0)
     sage: side_conversion('right')
     sage: side_conversion('r')
     -1
     sage: side_conversion(1)
     -1
     sage: side_conversion(-1)
     -1
sage.dynamics.interval exchanges.template.twin list iet(a=None)
     Returns the twin list of intervals.
     The twin intervals is the correspondance between positions of labels in such way that a[interval][position] is
     a[1-interval][twin[interval][position]]
     INPUT:
        • a - two lists of labels
     OUTPUT:
     list – a list of two lists of integers
```

 $\verb|sage.dynamics.interval_exchanges.template.twin_list_li|(a=None)$

Returns the twin list of intervals

INPUT:

• a - two lists of labels

OUTPUT:

list – a list of two lists of couples of integers

4.5 Interval Exchange Transformations and Linear Involution

Warning: This module is deprecated. You are advised to install and use the surface_dynamics package instead available at https://pypi.python.org/pypi/surface_dynamics/

An interval exchange transformation is a map defined on an interval (see help(iet.IntervalExchangeTransformation) for a more complete help.

EXAMPLES:

Initialization of a simple iet with integer lengths:

```
sage: T = iet.IntervalExchangeTransformation(Permutation([3,2,1]), [3,1,2])
doctest:warning
DeprecationWarning: IntervalExchangeTransformation is deprecated and will be removed_
→from Sage.
You are advised to install the surface_dynamics package via:
sage -pip install surface_dynamics
If you do not have write access to the Sage installation you can
alternatively do
sage -pip install surface_dynamics --user
The package surface_dynamics subsumes all flat surface related
computation that are currently available in Sage. See more
information at
http://www.labri.fr/perso/vdelecro/surface-dynamics/latest/
See http://trac.sagemath.org/20695 for details.
doctest:warning
DeprecationWarning: Permutation is deprecated and will be removed from Sage.
You are advised to install the surface_dynamics package via:
sage -pip install surface_dynamics
If you do not have write access to the Sage installation you can
alternatively do
sage -pip install surface_dynamics --user
The package surface_dynamics subsumes all flat surface related
computation that are currently available in Sage. See more
information at
http://www.labri.fr/perso/vdelecro/surface-dynamics/latest/
See http://trac.sagemath.org/20695 for details.
Interval exchange transformation of [0, 6[ with permutation
1 2 3
3 2 1
```

Rotation corresponds to iet with two intervals:

```
sage: p = iet.Permutation('a b', 'b a')
sage: T = iet.IntervalExchangeTransformation(p, [1, (sqrt(5)-1)/2])
sage: print(T.in_which_interval(0))
a
sage: print(T.in_which_interval(T(0)))
a
sage: print(T.in_which_interval(T(T(0))))
b
sage: print(T.in_which_interval(T(T(0))))
```

```
a
```

There are two plotting methods for iet:

```
sage: p = iet.Permutation('a b c','c b a')
sage: T = iet.IntervalExchangeTransformation(p, [1, 2, 3])
```

Bases: sage.structure.sage_object.SageObject

Interval exchange transformation

INPUT:

- permutation a permutation (LabelledPermutationIET)
- lengths the list of lengths

EXAMPLES:

Direct initialization:

```
sage: p = iet.IET(('a b c','c b a'), {'a':1,'b':1,'c':1})
sage: p.permutation()
a b c
c b a
sage: p.lengths()
[1, 1, 1]
```

Initialization from a iet.Permutation:

```
sage: perm = iet.Permutation('a b c','c b a')
sage: l = [0.5,1,1.2]
sage: t = iet.IET(perm,l)
sage: t.permutation() == perm
True
sage: t.lengths() == l
True
```

Initialization from a Permutation:

```
sage: p = Permutation([3,2,1])
sage: iet.IET(p, [1,1,1])
Interval exchange transformation of [0, 3[ with permutation
1 2 3
3 2 1
```

If it is not possible to convert lengths to real values an error is raised:

```
sage: iet.IntervalExchangeTransformation(('a b','b a'),['e','f'])
Traceback (most recent call last):
...
TypeError: unable to convert 'e' to a float
```

The value for the lengths must be positive:

```
sage: iet.IET(('a b','b a'),[-1,-1])
Traceback (most recent call last):
```

```
ValueError: lengths must be positive
```

domain_singularities()

Returns the list of singularities of T

OUTPUT:

list – positive reals that corresponds to singularities in the top interval

EXAMPLES:

```
sage: t = iet.IET(("a b","b a"), [1, sqrt(2)])
sage: t.domain_singularities()
[0, 1, sqrt(2) + 1]
```

in_which_interval (x, interval=0)

Returns the letter for which x is in this interval.

INPUT:

- x a positive number
- interval (default: 'top') 'top' or 'bottom'

OUTPUT:

label – a label corresponding to an interval

```
sage: t = iet.IntervalExchangeTransformation(('a b c','c b a'),[1,1,1])
sage: t.in_which_interval(0)
'a'
sage: t.in_which_interval(0.3)
'a'
sage: t.in_which_interval(1)
'b'
sage: t.in_which_interval(1.9)
'b'
sage: t.in_which_interval(2)
'c'
sage: t.in_which_interval(2.1)
'c'
sage: t.in_which_interval(3)
Traceback (most recent call last):
...
ValueError: your value does not lie in [0;1[
```

inverse()

Returns the inverse iet.

OUTPUT:

iet – the inverse interval exchange transformation

```
sage: p = iet.Permutation("a b","b a")
sage: s = iet.IET(p, [1,sqrt(2)-1])
sage: t = s.inverse()
sage: t.permutation()
b a
a b
```

```
sage: t.lengths()
[1, sqrt(2) - 1]
sage: t*s
Interval exchange transformation of [0, sqrt(2)[ with permutation
aa bb
aa bb
```

We can verify with the method .is_identity():

```
sage: p = iet.Permutation("a b c d","d a c b")
sage: s = iet.IET(p, [1, sqrt(2), sqrt(3), sqrt(5)])
sage: (s * s.inverse()).is_identity()
True
sage: (s.inverse() * s).is_identity()
True
```

is_identity()

Returns True if self is the identity.

OUTPUT:

boolean - the answer

EXAMPLES:

```
sage: p = iet.Permutation("a b","b a")
sage: q = iet.Permutation("c d","d c")
sage: s = iet.IET(p, [1,5])
sage: t = iet.IET(q, [5,1])
sage: (s*t).is_identity()
True
sage: (t*s).is_identity()
True
```

length()

Returns the total length of the interval.

OUTPUT:

real - the length of the interval

EXAMPLES:

```
sage: t = iet.IntervalExchangeTransformation(('a b', 'b a'),[1,1])
sage: t.length()
2
```

lengths()

Returns the list of lengths associated to this iet.

OUTPUT:

list – the list of lengths of subinterval

```
sage: p = iet.IntervalExchangeTransformation(('a b','b a'),[1,3])
sage: p.lengths()
[1, 3]
```

normalize(total=1)

Returns a interval exchange transformation of normalized lengths.

The normalization consists in multiplying all lengths by a constant in such way that their sum is given by total (default is 1).

INPUT:

• total - (default: 1) The total length of the interval

OUTPUT:

iet - the normalized iet

EXAMPLES:

```
sage: t = iet.IntervalExchangeTransformation(('a b', 'b a'), [1,3])
sage: t.length()
4
sage: s = t.normalize(2)
sage: s.length()
2
sage: s.lengths()
[1/2, 3/2]
```

permutation()

Returns the permutation associated to this iet.

OUTPUT:

permutation - the permutation associated to this iet

EXAMPLES:

```
sage: perm = iet.Permutation('a b c','c b a')
sage: p = iet.IntervalExchangeTransformation(perm, (1,2,1))
sage: p.permutation() == perm
True
```

plot (position=(0, 0), vertical_alignment='center', horizontal_alignment='left', interval_height=0.1, labels_height=0.05, fontsize=14, labels=True, colors=None)
Returns a picture of the interval exchange transformation.

INPUT:

- position a 2-uple of the position
- horizontal_alignment left (default), center or right
- labels boolean (default: True)
- fontsize the size of the label

OUTPUT:

2d plot – a plot of the two intervals (domain and range)

```
sage: t = iet.IntervalExchangeTransformation(('a b','b a'),[1,1])
sage: t.plot_two_intervals()
Graphics object consisting of 8 graphics primitives
```

plot function(**d)

Return a plot of the interval exchange transformation as a function.

INPUT:

• Any option that is accepted by line2d

OUTPUT:

2d plot – a plot of the iet as a function

EXAMPLES:

```
sage: t = iet.IntervalExchangeTransformation(('a b c d','d a c b'),[1,1,1,1])
sage: t.plot_function(rgbcolor=(0,1,0))
Graphics object consisting of 4 graphics primitives
```

Returns a picture of the interval exchange transformation.

INPUT:

- position a 2-uple of the position
- horizontal_alignment left (default), center or right
- labels boolean (default: True)
- fontsize the size of the label

OUTPUT:

2d plot – a plot of the two intervals (domain and range)

EXAMPLES:

```
sage: t = iet.IntervalExchangeTransformation(('a b','b a'),[1,1])
sage: t.plot_two_intervals()
Graphics object consisting of 8 graphics primitives
```

range_singularities()

Returns the list of singularities of T^{-1}

OUTPUT:

list – real numbers that are singular for T^{-1}

EXAMPLES:

```
sage: t = iet.IET(("a b", "b a"), [1, sqrt(2)])
sage: t.range_singularities()
[0, sqrt(2), sqrt(2) + 1]
```

rauzy_move (side='right', iterations=1)

Performs a Rauzy move.

INPUT:

- side 'left' (or 'l' or 0) or 'right' (or 'r' or 1)
- iterations integer (default :1) the number of iteration of Rauzy moves to perform

OUTPUT:

iet – the Rauzy move of self

EXAMPLES:

```
sage: phi = QQbar((sqrt(5)-1)/2)
sage: t1 = iet.IntervalExchangeTransformation(('a b','b a'),[1,phi])
sage: t2 = t1.rauzy_move().normalize(t1.length())
sage: 12 = t2.lengths()
sage: 11 = t1.lengths()
sage: 12[0] == 11[1] and 12[1] == 11[0]
True
```

show()

Shows a picture of the interval exchange transformation

EXAMPLES:

```
sage: phi = QQbar((sqrt(5)-1)/2)
sage: t = iet.IntervalExchangeTransformation(('a b', 'b a'),[1,phi])
sage: t.show()
```

singularities()

The list of singularities of T and T^{-1} .

OUTPUT

list - two lists of positive numbers which corresponds to extremities of subintervals

```
sage: t = iet.IntervalExchangeTransformation(('a b','b a'),[1/2,3/2])
sage: t.singularities()
[[0, 1/2, 2], [0, 3/2, 2]]
```



CHAPTER

FIVE

SANDPILES

Functions and classes for mathematical sandpiles.

Version: 2.4

AUTHOR:

• David Perkinson (June 4, 2015) Upgraded from version 2.3 to 2.4.

MAJOR CHANGES

- 1. Eliminated dependence on 4ti2, substituting the use of Polyhedron methods. Thus, no optional packages are necessary.
- 2. Fixed bug in Sandpile. __init__ so that now multigraphs are handled correctly.
- 3. Created sandpiles to handle examples of Sandpiles in analogy with graphs, simplicial_complexes, and polytopes. In the process, we implemented a much faster way of producing the sandpile grid graph.
- 4. Added support for open and closed sandpile Markov chains.
- 5. Added support for Weierstrass points.
- 6. Implemented the Cori-Le Borgne algorithm for computing ranks of divisors on complete graphs.

NEW METHODS

Sandpile: avalanche_polynomial, genus, group_gens, help, jacobian_representatives, markov_chain, picard_representatives, smith_form, stable_configs, stationary_density, tutte_polynomial.

SandpileConfig: burst size, help.

SandpileDivisor: help, is_linearly_equivalent, is_q_reduced, is_weierstrass_pt, polytope, polytope_integer_pts, q_reduced, rank, simulate_threshold, stabilize, weierstrass_div, weierstrass_gap_seq, weierstrass_pts, weierstrass_rank_seq.

MINOR CHANGES

- The sink argument to Sandpile.__init__ now defaults to the first vertex.
- A SandpileConfig or SandpileDivisor may now be multiplied by an integer.
- Sped up __add__ method for SandpileConfig and SandpileDivisor.
- Enhanced string representation of a Sandpile (via _repr_ and the name methods).
- Recurrents for complete graphs and cycle graphs are computed more quickly.
- The stabilization code for SandpileConfig has been made more efficient.
- Added optional probability distribution arguments to add_random methods.

- Marshall Hampton (2010-1-10) modified for inclusion as a module within Sage library.
- David Perkinson (2010-12-14) added show3d(), fixed bug in resolution(), replaced elementary_divisors() with invariant_factors(), added show() for SandpileConfig and SandpileDivisor.
- David Perkinson (2010-9-18): removed is_undirected, added show(), added verbose arguments to several functions to display SandpileConfigs and divisors as lists of integers
- David Perkinson (2010-12-19): created separate SandpileConfig, SandpileDivisor, and Sandpile classes
- David Perkinson (2009-07-15): switched to using config_to_list instead of .values(), thus fixing a few bugs when not using integer labels for vertices.
- David Perkinson (2009): many undocumented improvements
- David Perkinson (2008-12-27): initial version

EXAMPLES:

For general help, enter Sandpile.help(), SandpileConfig.help(), and SandpileDivisor.help(). Miscellaneous examples appear below.

A weighted directed graph given as a Python dictionary:

The associated sandpile with 0 chosen as the sink:

```
sage: S = Sandpile(g,0)
```

Or just:

```
sage: S = Sandpile(g)
```

A picture of the graph:

```
sage: S.show() # long time
```

The relevant Laplacian matrices:

```
sage: S.laplacian()
[ 0  0  0  0  0  0]
[-1  3 -1 -1  0]
[ 0  -1  3 -1 -1]
[ 0  -1 -1  3 -1]
[ 0  0 -1 -1  2]
sage: S.reduced_laplacian()
[ 3  -1  -1  0]
[-1  3  -1 -1]
[ -1  -1  3  -1]
[ 0  -1  -1  2]
```

The number of elements of the sandpile group for S:

```
sage: S.group_order()
8
```

The structure of the sandpile group:

```
sage: S.invariant_factors()
[1, 1, 1, 8]
```

The elements of the sandpile group for S:

```
sage: S.recurrents()
[{1: 2, 2: 2, 3: 2, 4: 1},
    {1: 2, 2: 2, 3: 2, 4: 0},
    {1: 2, 2: 1, 3: 2, 4: 0},
    {1: 2, 2: 2, 3: 0, 4: 1},
    {1: 2, 2: 0, 3: 2, 4: 1},
    {1: 2, 2: 2, 3: 1, 4: 0},
    {1: 2, 2: 2, 3: 1, 4: 0},
    {1: 2, 2: 2, 3: 1, 4: 1}]
```

The maximal stable element (2 grains of sand on vertices 1, 2, and 3, and 1 grain of sand on vertex 4:

```
sage: S.max_stable()
{1: 2, 2: 2, 3: 2, 4: 1}
sage: S.max_stable().values()
[2, 2, 2, 1]
```

The identity of the sandpile group for S:

```
sage: S.identity()
{1: 2, 2: 2, 3: 2, 4: 0}
```

An arbitrary sandpile configuration:

```
sage: c = SandpileConfig(S,[1,0,4,-3])
sage: c.equivalent_recurrent()
{1: 2, 2: 2, 3: 2, 4: 0}
```

Some group operations:

```
sage: m = S.max_stable()
sage: i = S.identity()
sage: m.values()
[2, 2, 2, 1]
sage: i.values()
[2, 2, 2, 0]
sage: m + i
              # coordinate-wise sum
{1: 4, 2: 4, 3: 4, 4: 1}
sage: m - i
{1: 0, 2: 0, 3: 0, 4: 1}
sage: m & i # add, then stabilize
{1: 2, 2: 2, 3: 2, 4: 1}
sage: e = m + m
sage: e
{1: 4, 2: 4, 3: 4, 4: 2}
sage: ~e # stabilize
{1: 2, 2: 2, 3: 2, 4: 0}
sage: a = -m
sage: a & m
{1: 0, 2: 0, 3: 0, 4: 0}
sage: a * m # add, then find the equivalent recurrent
{1: 2, 2: 2, 3: 2, 4: 0}
sage: a^3 # a*a*a
```

```
{1: 2, 2: 2, 3: 2, 4: 1}
sage: a^(-1) == m
True
sage: a < m # every coordinate of a is < that of m
True
```

Firing an unstable vertex returns resulting configuration:

```
sage: c = S.max_stable() + S.identity()
sage: c.fire_vertex(1)
{1: 1, 2: 5, 3: 5, 4: 1}
sage: c
{1: 4, 2: 4, 3: 4, 4: 1}
```

Fire all unstable vertices:

```
sage: c.unstable()
[1, 2, 3]
sage: c.fire_unstable()
{1: 3, 2: 3, 3: 3, 4: 3}
```

Stabilize c, returning the resulting configuration and the firing vector:

```
sage: c.stabilize(True)
[{1: 2, 2: 2, 3: 2, 4: 1}, {1: 6, 2: 8, 3: 8, 4: 8}]
sage: c
{1: 4, 2: 4, 3: 4, 4: 1}
sage: S.max_stable() & S.identity() == c.stabilize()
True
```

The number of superstable configurations of each degree:

```
sage: S.h_vector()
[1, 3, 4]
sage: S.postulation()
2
```

the saturated homogeneous toppling ideal:

```
sage: S.ideal()

Ideal (x1 - x0, x3*x2 - x0^2, x4^2 - x0^2, x2^3 - x4*x3*x0, x4*x2^2 - x3^2*x0, x3^3 - x4*x2*x0, x4*x3^2 - x2^2*x0) of Multivariate Polynomial Ring in x4, x3, x2, x1, x0, where Rational Field
```

its minimal free resolution:

```
sage: S.resolution()
'R^1 <-- R^15 <-- R^13 <-- R^4'</pre>
```

and its Betti numbers:

```
sage: S.betti()
              2
                  3
            1
                       4
  0:
      1
            1
            2
  1:
                2
  2:
           4
               13
                    13
                         4
```

```
total: 1 7 15 13 4
```

Some various ways of creating Sandpiles:

```
sage: S = sandpiles.Complete(4) # for more options enter ``sandpile.TAB``
sage: S = sandpiles.Wheel(6)
```

A multidigraph with loops (vertices 0, 1, 2; for example, there is a directed edge from vertex 2 to vertex 1 of weight 3, which can be thought of as three directed edges of the form (2,3). There is also a single loop at vertex 2 and an edge (2,0) of weight 2):

```
sage: S = Sandpile({0:[1,2], 1:[0,0,2], 2:[0,0,1,1,1,2], 3:[2]})
```

Using the graph library (vertex 1 is specified as the sink; omitting this would make the sink vertex 0 by default):

```
sage: S = Sandpile(graphs.PetersenGraph(),1)
```

Distribution of avalanche sizes:

Working with sandpile divisors:

```
sage: S = sandpiles.Complete(4)
sage: D = SandpileDivisor(S, [0,0,0,5])
sage: E = D.stabilize(); E
\{0: 1, 1: 1, 2: 1, 3: 2\}
sage: D.is_linearly_equivalent(E)
True
sage: D.g_reduced()
\{0: 4, 1: 0, 2: 0, 3: 1\}
sage: S = sandpiles.Complete(4)
sage: D = SandpileDivisor(S, [0,0,0,5])
sage: E = D.stabilize(); E
\{0: 1, 1: 1, 2: 1, 3: 2\}
sage: D.is_linearly_equivalent(E)
sage: D.q_reduced()
\{0: 4, 1: 0, 2: 0, 3: 1\}
sage: D.rank()
sage: sorted(D.effective_div(), key=str)
[\{0: 0, 1: 0, 2: 0, 3: 5\},
\{0: 0, 1: 0, 2: 4, 3: 1\},\
\{0: 0, 1: 4, 2: 0, 3: 1\},\
\{0: 1, 1: 1, 2: 1, 3: 2\},\
```

```
{0: 4, 1: 0, 2: 0, 3: 1}]
sage: sorted(D.effective_div(False))
[[0, 0, 0, 5], [0, 0, 4, 1], [0, 4, 0, 1], [1, 1, 1, 2], [4, 0, 0, 1]]
sage: D.rank()
sage: D.rank(True)
(2, \{0: 2, 1: 1, 2: 0, 3: 0\})
sage: E = D.rank(True)[1] # E proves the rank is not 3
sage: E.values()
[2, 1, 0, 0]
sage: E.deg()
sage: rank(D - E)
sage: (D - E).effective_div()
[]
sage: D.weierstrass_pts()
(0, 1, 2, 3)
sage: D.weierstrass_rank_seq(0)
(2, 1, 0, 0, 0, -1)
sage: D.weierstrass_pts()
(0, 1, 2, 3)
sage: D.weierstrass_rank_seq(0)
(2, 1, 0, 0, 0, -1)
```

```
class sage.sandpiles.sandpile (g, sink=None)
```

Bases: sage.graphs.digraph.DiGraph

Class for Dhar's abelian sandpile model.

all_k_config(k)

The constant configuration with all values set to k.

INPUT:

k-integer

OUTPUT:

SandpileConfig

EXAMPLES:

```
sage: s = sandpiles.Diamond()
sage: s.all_k_config(7)
{1: 7, 2: 7, 3: 7}
```

$\mathbf{all} _\mathbf{k} _\mathbf{div} \left(k \right)$

The divisor with all values set to k.

INPUT:

k – integer

OUTPUT:

SandpileDivisor

```
sage: S = sandpiles.House()
sage: S.all_k_div(7)
{0: 7, 1: 7, 2: 7, 3: 7, 4: 7}
```

avalanche_polynomial (multivariable=True)

The avalanche polynomial. See NOTE for details.

INPUT:

```
multivariable - (default: True) boolean
```

OUTPUT:

polynomial

EXAMPLES:

```
sage: s = sandpiles.Complete(4)
sage: s.avalanche_polynomial()
9*x0*x1*x2 + 2*x0*x1 + 2*x0*x2 + 2*x1*x2 + 3*x0 + 3*x1 + 3*x2 + 24
sage: s.avalanche_polynomial(False)
9*x0^3 + 6*x0^2 + 9*x0 + 24
```

Note: For each nonsink vertex v, let x_v be an indeterminate. If (r,v) is a pair consisting of a recurrent r and nonsink vertex v, then for each nonsink vertex w, let n_w be the number of times vertex w fires in the stabilization of r+v. Let M(r,v) be the monomial $\prod_w x_w^{n_w}$, i.e., the exponent records the vector of n_w as w ranges over the nonsink vertices. The avalanche polynomial is then the sum of M(r,v) as r ranges over the recurrents and v ranges over the nonsink vertices. If multivariable is False, then set all the indeterminates equal to each other (and, thus, only count the number of vertex firings in the stabilizations, forgetting which particular vertices fired).

betti (verbose=True)

The Betti table for the homogeneous toppling ideal. If verbose is True, it prints the standard Betti table, otherwise, it returns a less formatted table.

INPUT:

```
verbose - (default: True) boolean
```

OUTPUT:

Betti numbers for the sandpile

EXAMPLES:

```
sage: S = sandpiles.Diamond()
sage: S.betti()
           0
                         2
                               3
                  1
           1
    0:
    1:
                  2
    2:
                  4
                               4
total:
                  6
           1
sage: S.betti(False)
[1, 6, 9, 4]
```

betti_complexes()

The support-complexes with non-trivial homology. (See NOTE.)

OUTPUT:

list (of pairs [divisors, corresponding simplicial complex])

EXAMPLES:

```
sage: S = Sandpile({0:{},1:{}0:1, 2:1, 3:4},2:{}3:5},3:{}1:1, 2:1},0)
sage: p = S.betti_complexes()
sage: p[0]
[{0: -8, 1: 5, 2: 4, 3: 1}, Simplicial complex with vertex set (1, 2, 3) and.
\rightarrow facets {(1, 2), (3,)}]
sage: S.resolution()
'R^1 <-- R^5 <-- R^1'
sage: S.betti()
               1
                     2.
   ______
   0: 1
   1:
              5
                   5
   2:
total: 1 5 5
sage: len(p)
11
sage: p[0][1].homology()
{0: Z, 1: 0}
sage: p[-1][1].homology()
{0: 0, 1: 0, 2: Z}
```

Note: A support-complex is the simplicial complex formed from the supports of the divisors in a linear system.

burning_config()

The minimal burning configuration.

OUTPUT:

dict (configuration)

EXAMPLES:

Note: The burning configuration and script are computed using a modified version of Speer's script algorithm. This is a generalization to directed multigraphs of Dhar's burning algorithm.

A burning configuration is a nonnegative integer-linear combination of the rows of the reduced Laplacian matrix having nonnegative entries and such that every vertex has a path from some vertex in its support. The corresponding burning script gives the integer-linear combination needed to obtain the burning configuration. So if b is the burning configuration, σ is its script, and \tilde{L} is the reduced Laplacian, then $\sigma \cdot \tilde{L} = b$. The minimal burning configuration is the one with the minimal script (its components are no larger than the components of any other script for a burning configuration).

The following are equivalent for a configuration c with burning configuration b having script σ :

- c is recurrent;
- c + b stabilizes to c;
- the firing vector for the stabilization of c + b is σ .

burning_script()

A script for the minimal burning configuration.

OUTPUT:

dict

EXAMPLES:

```
sage: g = {0:{},1:{0:1,3:1,4:1},2:{0:1,3:1,5:1},\
3:{2:1,5:1},4:{1:1,3:1},5:{2:1,3:1}}
sage: S = Sandpile(g,0)
sage: S.burning_config()
{1: 2, 2: 0, 3: 1, 4: 1, 5: 0}
sage: S.burning_config().values()
[2, 0, 1, 1, 0]
sage: S.burning_script()
{1: 1, 2: 3, 3: 5, 4: 1, 5: 4}
sage: script = S.burning_script().values()
sage: script
[1, 3, 5, 1, 4]
sage: matrix(script)*S.reduced_laplacian()
[2 0 1 1 0]
```

Note: The burning configuration and script are computed using a modified version of Speer's script algorithm. This is a generalization to directed multigraphs of Dhar's burning algorithm.

A burning configuration is a nonnegative integer-linear combination of the rows of the reduced Laplacian matrix having nonnegative entries and such that every vertex has a path from some vertex in its support. The corresponding burning script gives the integer-linear combination needed to obtain the burning configuration. So if b is the burning configuration, s is its script, and $L_{\rm red}$ is the reduced Laplacian, then $s \cdot L_{\rm red} = b$. The minimal burning configuration is the one with the minimal script (its components are no larger than the components of any other script for a burning configuration).

The following are equivalent for a configuration c with burning configuration b having script s:

- c is recurrent;
- c + b stabilizes to c;
- the firing vector for the stabilization of c+b is s.

canonical_divisor()

The canonical divisor. This is the divisor with deg(v) - 2 grains of sand on each vertex (not counting

loops). Only for undirected graphs.

OUTPUT:

SandpileDivisor

EXAMPLES:

```
sage: S = sandpiles.Complete(4)
sage: S.canonical_divisor()
{0: 1, 1: 1, 2: 1, 3: 1}
sage: s = Sandpile({0:[1,1],1:[0,0,1,1,1]},0)
sage: s.canonical_divisor() # loops are disregarded
{0: 0, 1: 0}
```

Warning: The underlying graph must be undirected.

dict()

A dictionary of dictionaries representing a directed graph.

OUTPUT:

dict

EXAMPLES:

```
sage: S = sandpiles.Diamond()
sage: S.dict()
{0: {1: 1, 2: 1},
    1: {0: 1, 2: 1, 3: 1},
    2: {0: 1, 1: 1, 3: 1},
    3: {1: 1, 2: 1}}
sage: S.sink()
0
```

genus()

The genus: (# non-loop edges) - (# vertices) + 1. Only defined for undirected graphs.

OUTPUT:

integer

EXAMPLES:

```
sage: sandpiles.Complete(4).genus()
3
sage: sandpiles.Cycle(5).genus()
1
```

groebner()

A Groebner basis for the homogeneous toppling ideal. It is computed with respect to the standard sandpile ordering (see ring).

OUTPUT:

Groebner basis

group_gens (verbose=True)

A minimal list of generators for the sandpile group. If verbose is False then the generators are represented as lists of integers.

INPUT:

verbose - (default: True) boolean

OUTPUT:

list of SandpileConfig (or of lists of integers if verbose is False)

EXAMPLES:

```
sage: s = sandpiles.Cycle(5)
sage: s.group_gens()
[{1: 1, 2: 1, 3: 1, 4: 0}]
sage: s.group_gens()[0].order()
5
sage: s = sandpiles.Complete(5)
sage: s.group_gens(False)
[[2, 2, 3, 2], [2, 3, 2, 2], [3, 2, 2, 2]]
sage: [i.order() for i in s.group_gens()]
[5, 5, 5]
sage: s.invariant_factors()
[1, 5, 5, 5]
```

group_order()

The size of the sandpile group.

OUTPUT:

integer

EXAMPLES:

```
sage: S = sandpiles.House()
sage: S.group_order()
11
```

h_vector()

The number of superstable configurations in each degree. Equivalently, this is the list of first differences of the Hilbert function of the (homogeneous) toppling ideal.

OUTPUT:

list of nonnegative integers

```
sage: s = sandpiles.Grid(2,2)
sage: s.hilbert_function()
[1, 5, 15, 35, 66, 106, 146, 178, 192]
sage: s.h_vector()
[1, 4, 10, 20, 31, 40, 40, 32, 14]
```

static help(verbose=True)

List of Sandpile-specific methods (not inherited from Graph). If verbose, include short descriptions.

INPUT:

verbose - (default: True) boolean

OUTPUT:

printed string

```
sage: Sandpile.help() # long time
For detailed help with any method FOO listed below,
enter "Sandpile.FOO?" or enter "S.FOO?" for any Sandpile S.
                        -- The constant configuration with all values set to...
all_k_config
۰k.
all_k_div
                       -- The divisor with all values set to k.
avalanche_polynomial -- The avalanche polynomial.
                        -- The Betti table for the homogeneous toppling.
bet.t.i
⇒ideal.
betti_complexes
                    -- The support-complexes with non-trivial homology.
                       -- The minimal burning configuration.
burning_config
burning_script
                       -- A script for the minimal burning configuration.
canonical_divisor
                        -- The canonical divisor.
dict
                        -- A dictionary of dictionaries representing a...
→directed graph.
                        -- The genus: (# non-loop edges) - (# vertices) + 1.
genus
                       -- A Groebner basis for the homogeneous toppling.
groebner
⊶ideal.
                       -- A minimal list of generators for the sandpile.
group_gens
⊶group.
                        -- The size of the sandpile group.
group_order
h_vector
                        -- The number of superstable configurations in each_
-degree.
help
                        -- List of Sandpile-specific methods (not inherited...
→from "Graph").
                        -- The Hilbert function of the homogeneous toppling.
hilbert function
→ideal.
                        -- The saturated homogeneous toppling ideal.
ideal
identity
                        -- The identity configuration.
                        -- The in-degree of a vertex or a list of all in-
in_degree
⇔degrees.
invariant_factors
                       -- The invariant factors of the sandpile group.
is_undirected
                       -- Is the underlying graph undirected?
jacobian_representatives -- Representatives for the elements of the Jacobian,
⇔group.
laplacian
                        -- The Laplacian matrix of the graph.
markov_chain
                       -- The sandpile Markov chain for configurations or_
⊶divisors.
                      The maximal stable configuration.The maximal stable divisor.
max_stable
max_stable_div
                       -- The maximal superstable configurations.
max_superstables
min_recurrents
                       -- The minimal recurrent elements.
nonsink_vertices
                       -- The nonsink vertices.
nonspecial_divisors
                       -- The nonspecial divisors.
                       -- The out-degree of a vertex or a list of all out-
out_degree
⊶degrees.
```

```
picard_representatives
                        -- Representatives of the divisor classes of degree...
\rightarrowd in the Picard group.
points
                        -- Generators for the multiplicative group of zeros.
\rightarrowof the sandpile ideal.
postulation
                       -- The postulation number of the toppling ideal.
                       -- The recurrent configurations.
recurrents
reduced_laplacian
                       -- The reduced Laplacian matrix of the graph.
                       -- A copy of the sandpile with vertex names permuted.
reorder_vertices
resolution
                        -- A minimal free resolution of the homogeneous.
→toppling ideal.
ring
                        -- The ring containing the homogeneous toppling.
⇒ideal.
show
                        -- Draw the underlying graph.
show3d
                        -- Draw the underlying graph.
                        -- The sink vertex.
sink
smith_form
                        -- The Smith normal form for the Laplacian.
solve
                        -- Approximations of the complex affine zeros of the
⇔sandpile ideal.
                        -- Generator for all stable configurations.
stable_configs
stationary_density
                       -- The stationary density of the sandpile.
                       -- The superstable configurations.
superstables
symmetric_recurrents
                       -- The symmetric recurrent configurations.
tutte_polynomial
                       -- The Tutte polynomial of the underlying graph.
                       -- The unsaturated, homogeneous toppling ideal.
unsaturated_ideal
                        -- The version number of Sage Sandpiles.
version
zero_config
                        -- The all-zero configuration.
                        -- The all-zero divisor.
zero_div
```

hilbert function()

The Hilbert function of the homogeneous toppling ideal.

OUTPUT:

list of nonnegative integers

EXAMPLES:

```
sage: s = sandpiles.Wheel(5)
sage: s.hilbert_function()
[1, 5, 15, 31, 45]
sage: s.h_vector()
[1, 4, 10, 16, 14]
```

ideal (gens=False)

The saturated homogeneous toppling ideal. If gens is True, the generators for the ideal are returned instead.

INPUT:

```
gens - (default: False) boolean
```

OUTPUT:

ideal or, optionally, the generators of an ideal

identity (verbose=True)

The identity configuration. If verbose is False, the configuration are converted to a list of integers.

INPUT:

verbose - (default: True) boolean

OUTPUT:

SandpileConfig or a list of integers If verbose is False, the configuration are converted to a list of integers.

EXAMPLES:

```
sage: s = sandpiles.Diamond()
sage: s.identity()
{1: 2, 2: 2, 3: 0}
sage: s.identity(False)
[2, 2, 0]
sage: s.identity() & s.max_stable() == s.max_stable()
True
```

in_degree (v=None)

The in-degree of a vertex or a list of all in-degrees.

INPUT:

∨ – (optional) vertex name

OUTPUT:

integer or dict

EXAMPLES:

```
sage: s = sandpiles.House()
sage: s.in_degree()
{0: 2, 1: 2, 2: 3, 3: 3, 4: 2}
sage: s.in_degree(2)
3
```

invariant_factors()

The invariant factors of the sandpile group.

OUTPUT:

list of integers

```
sage: s = sandpiles.Grid(2,2)
sage: s.invariant_factors()
[1, 1, 8, 24]
```

is undirected()

Is the underlying graph undirected? True if (u, v) is and edge if and only if (v, u) is an edge, each edge with the same weight.

OUTPUT:

boolean

EXAMPLES:

```
sage: sandpiles.Complete(4).is_undirected()
True
sage: s = Sandpile({0:[1,2], 1:[0,2], 2:[0]}, 0)
sage: s.is_undirected()
False
```

jacobian_representatives (verbose=True)

Representatives for the elements of the Jacobian group. If verbose is False, then lists representing the divisors are returned.

INPUT:

```
verbose - (default: True) boolean
```

OUTPUT:

list of SandpileDivisor (or of lists representing divisors)

EXAMPLES:

For an undirected graph, divisors of the form s - deg(s) *sink as s varies over the superstables forms a distinct set of representatives for the Jacobian group.:

```
sage: s = sandpiles.Complete(3)
sage: s.superstables(False)
[[0, 0], [0, 1], [1, 0]]
sage: s.jacobian_representatives(False)
[[0, 0, 0], [-1, 0, 1], [-1, 1, 0]]
```

If the graph is directed, the representatives described above may by equivalent modulo the rowspan of the Laplacian matrix:

```
sage: s = Sandpile({0: {1: 1, 2: 2}, 1: {0: 2, 2: 4}, 2: {0: 4, 1: 2}},0)
sage: s.group_order()
28
sage: s.jacobian_representatives()
[{0: -5, 1: 3, 2: 2}, {0: -4, 1: 3, 2: 1}]
```

Let τ be the nonnegative generator of the kernel of the transpose of the Laplacian, and let tau_s be it sink component, then the sandpile group is isomorphic to the direct sum of the cyclic group of order τ_s and the Jacobian group. In the example above, we have:

```
sage: s.laplacian().left_kernel()
Free module of degree 3 and rank 1 over Integer Ring
Echelon basis matrix:
[14 5 8]
```

Note: The Jacobian group is the set of all divisors of degree zero modulo the integer rowspan of the Laplacian matrix.

laplacian()

The Laplacian matrix of the graph. Its rows encode the vertex firing rules.

OUTPUT:

matrix

EXAMPLES:

```
sage: G = sandpiles.Diamond()
sage: G.laplacian()
[ 2 -1 -1   0]
[-1   3 -1 -1]
[-1 -1   3 -1]
[ 0 -1 -1   2]
```

Warning: The function laplacian_matrix should be avoided. It returns the indegree version of the Laplacian.

markov_chain (state, distrib=None)

The sandpile Markov chain for configurations or divisors. The chain starts at state. See NOTE for details.

INPUT:

- state SandpileConfig, SandpileDivisor, or list representing one of these
- distrib (optional) list of nonnegative numbers summing to 1 (representing a prob. dist.)

OUTPUT:

generator for Markov chain (see NOTE)

```
sage: s = sandpiles.Complete(4)
sage: m = s.markov\_chain([0,0,0])
sage: next(m)
                      # random
{1: 0, 2: 0, 3: 0}
sage: next(m).values() # random
[0, 0, 0]
sage: next(m).values() # random
[0, 0, 0]
sage: next(m).values() # random
[0, 0, 0]
sage: next(m).values() # random
[0, 1, 0]
sage: next(m).values() # random
[0, 2, 0]
sage: next(m).values() # random
[0, 2, 1]
sage: next(m).values() # random
[1, 2, 1]
sage: next(m).values() # random
[2, 2, 1]
sage: m = s.markov_chain(s.zero_div(), [0.1, 0.1, 0.1, 0.7])
sage: next(m).values() # random
[0, 0, 0, 1]
sage: next(m).values() # random
```

```
[0, 0, 1, 1]
sage: next(m).values() # random
[0, 0, 1, 2]
sage: next(m).values() # random
[1, 1, 2, 0]
sage: next(m).values() # random
[1, 1, 2, 1]
sage: next(m).values() # random
[1, 1, 2, 2]
sage: next(m).values() # random
[1, 1, 2, 3]
sage: next(m).values() # random
[1, 1, 2, 4]
sage: next(m).values() # random
[1, 1, 3, 4]
```

Note: The closed sandpile Markov chain has state space consisting of the configurations on a sandpile. It transitions from a state by choosing a vertex at random (according to the probability distribution distrib), dropping a grain of sand at that vertex, and stabilizing. If the chosen vertex is the sink, the chain stays at the current state.

The open sandpile Markov chain has state space consisting of the recurrent elements, i.e., the state space is the sandpile group. It transitions from the configuration c by choosing a vertex v at random according to distrib. The next state is the stabilization of c+v. If v is the sink vertex, then the stabilization of c+v is defined to be c.

Note that in either case, if distrib is specified, its length is equal to the total number of vertices (including the sink).

REFERENCES:

• [Lev2014]

max stable()

The maximal stable configuration.

OUTPUT:

SandpileConfig (the maximal stable configuration)

EXAMPLES:

```
sage: S = sandpiles.House()
sage: S.max_stable()
{1: 1, 2: 2, 3: 2, 4: 1}
```

max stable div()

The maximal stable divisor.

OUTPUT:

SandpileDivisor (the maximal stable divisor)

```
sage: s = sandpiles.Diamond()
sage: s.max_stable_div()
{0: 1, 1: 2, 2: 2, 3: 1}
```

```
sage: s.out_degree()
{0: 2, 1: 3, 2: 3, 3: 2}
```

max_superstables (verbose=True)

The maximal superstable configurations. If the underlying graph is undirected, these are the superstables of highest degree. If verbose is False, the configurations are converted to lists of integers.

INPUT:

```
verbose - (default: True) boolean
```

OUTPUT:

tuple of SandpileConfig

EXAMPLES:

```
sage: s = sandpiles.Diamond()
sage: s.superstables(False)
[[0, 0, 0],
  [0, 0, 1],
  [1, 0, 1],
  [0, 2, 0],
  [2, 0, 0],
  [0, 1, 1],
  [1, 0, 0],
  [0, 1, 1],
  [1, 0, 0],
  [0, 1, 0]]
sage: s.max_superstables(False)
[[1, 0, 1], [0, 2, 0], [2, 0, 0], [0, 1, 1]]
sage: s.h_vector()
[1, 3, 4]
```

min_recurrents (verbose=True)

The minimal recurrent elements. If the underlying graph is undirected, these are the recurrent elements of least degree. If verbose is False, the configurations are converted to lists of integers.

INPUT:

```
verbose - (default: True) boolean
```

OUTPUT:

list of SandpileConfig

```
sage: s = sandpiles.Diamond()
sage: s.recurrents(False)
[[2, 2, 1],
       [2, 2, 0],
       [1, 2, 0],
       [2, 0, 1],
       [0, 2, 1],
       [2, 1, 0],
       [1, 2, 1],
       [2, 1, 0],
       [1, 2, 1],
       [2, 0, 1], [0, 2, 1], [2, 1, 0]]
sage: s.min_recurrents(False)
[[1, 2, 0], [2, 0, 1], [0, 2, 1], [2, 1, 0]]
sage: [i.deg() for i in s.recurrents()]
[5, 4, 3, 3, 3, 3, 4, 4]
```

nonsink vertices()

The nonsink vertices.

OUTPUT:

list of vertices

EXAMPLES:

```
sage: s = sandpiles.Grid(2,3)
sage: s.nonsink_vertices()
[(1, 1), (1, 2), (1, 3), (2, 1), (2, 2), (2, 3)]
```

nonspecial_divisors (verbose=True)

The nonspecial divisors. Only for undirected graphs. (See NOTE.)

INPUT:

verbose - (default: True) boolean

OUTPUT:

list (of divisors)

EXAMPLES:

```
sage: S = sandpiles.Complete(4)
sage: ns = S.nonspecial_divisors()
sage: D = ns[0]
sage: D.values()
[-1, 0, 1, 2]
sage: D.deg()
2
sage: [i.effective_div() for i in ns]
[[], [], [], [], [], []]
```

Note: The "nonspecial divisors" are those divisors of degree g-1 with empty linear system. The term is only defined for undirected graphs. Here, g=|E|-|V|+1 is the genus of the graph (not counting loops as part of |E|). If verbose is False, the divisors are converted to lists of integers.

Warning: The underlying graph must be undirected.

out_degree (v=None)

The out-degree of a vertex or a list of all out-degrees.

INPUT:

 \lor - (optional) vertex name

OUTPUT:

integer or dict

```
sage: s = sandpiles.House()
sage: s.out_degree()
{0: 2, 1: 2, 2: 3, 3: 3, 4: 2}
```

```
sage: s.out_degree(2)
3
```

picard_representatives (d, verbose=True)

Representatives of the divisor classes of degree d in the Picard group. (Also see the documentation for jacobian_representatives.)

INPUT:

- d integer
- verbose (default: True) boolean

OUTPUT:

list of SandpileDivisors (or lists representing divisors)

EXAMPLES:

```
sage: s = sandpiles.Complete(3)
sage: s.superstables(False)
[[0, 0], [0, 1], [1, 0]]
sage: s.jacobian_representatives(False)
[[0, 0, 0], [-1, 0, 1], [-1, 1, 0]]
sage: s.picard_representatives(3,False)
[[3, 0, 0], [2, 0, 1], [2, 1, 0]]
```

points()

Generators for the multiplicative group of zeros of the sandpile ideal.

OUTPUT:

list of complex numbers

EXAMPLES:

The sandpile group in this example is cyclic, and hence there is a single generator for the group of solutions.

```
sage: S = sandpiles.Complete(4)
sage: S.points()
[[1, I, -I], [I, 1, -I]]
```

postulation()

The postulation number of the toppling ideal. This is the largest weight of a superstable configuration of the graph.

OUTPUT:

nonnegative integer

EXAMPLES:

```
sage: s = sandpiles.Complete(4)
sage: s.postulation()
3
```

recurrents (verbose=True)

The recurrent configurations. If verbose is False, the configurations are converted to lists of integers.

INPUT:

```
verbose - (default: True) boolean
```

OUTPUT:

list of recurrent configurations

EXAMPLES:

```
sage: r = Sandpile(graphs.HouseXGraph(),0).recurrents()
sage: r[:3]
[\{1: 2, 2: 3, 3: 3, 4: 1\}, \{1: 1, 2: 3, 3: 3, 4: 0\}, \{1: 1, 2: 3, 3: 3, 4: 1\}]
sage: sandpiles.Complete(4).recurrents(False)
[[2, 2, 2],
 [2, 2, 1],
 [2, 1, 2],
 [1, 2, 2],
 [2, 2, 0],
[2, 0, 2],
[0, 2, 2],
[2, 1, 1],
[1, 2, 1],
[1, 1, 2],
[2, 1, 0],
[2, 0, 1],
[1, 2, 0],
[1, 0, 2],
 [0, 2, 1],
[0, 1, 2]]
sage: sandpiles.Cycle(4).recurrents(False)
[[1, 1, 1], [0, 1, 1], [1, 0, 1], [1, 1, 0]]
```

reduced_laplacian()

The reduced Laplacian matrix of the graph.

OUTPUT:

matrix

EXAMPLES:

```
sage: S = sandpiles.Diamond()
sage: S.laplacian()
[ 2 -1 -1   0]
[-1   3 -1 -1]
[-1 -1   3 -1]
[ 0 -1 -1   2]
sage: S.reduced_laplacian()
[ 3 -1 -1]
[-1   3 -1]
[-1   -1   2]
```

Note: This is the Laplacian matrix with the row and column indexed by the sink vertex removed.

reorder_vertices()

A copy of the sandpile with vertex names permuted.

After reordering, vertex u comes before vertex v in the list of vertices if u is closer to the sink.

OUTPUT:

Sandpile

EXAMPLES:

```
sage: S = Sandpile({0:[1], 2:[0,1], 1:[2]})
sage: S.dict()
{0: {1: 1}, 1: {2: 1}, 2: {0: 1, 1: 1}}
sage: T = S.reorder_vertices()
```

The vertices 1 and 2 have been swapped:

```
sage: T.dict()
{0: {1: 1}, 1: {0: 1, 2: 1}, 2: {0: 1}}
```

resolution(verbose=False)

A minimal free resolution of the homogeneous toppling ideal. If verbose is True, then all of the mappings are returned. Otherwise, the resolution is summarized.

INPUT:

verbose - (default: False) boolean

OUTPUT:

free resolution of the toppling ideal

EXAMPLES:

```
sage: S = Sandpile({0: {}, 1: {0: 1, 2: 1, 3: 4}, 2: {3: 5}, 3: {1: 1, 2: 1}},
sage: S.resolution() # a Gorenstein sandpile graph
'R^1 <-- R^5 <-- R^1'
sage: S.resolution(True)
[x1^2 - x3*x0 x3*x1 - x2*x0 x3^2 - x2*x1 x2*x3 - x0^2 x2^2 - x1*x0],
[ x3  x2  0  x0  0]  [ x2^2 - x1*x0]
[-x1 -x3  x2  0 -x0]  [-x2*x3 + x0^2]
[ x0 x1 0 x2 0] [-x3^2 + x2*x1]
[ 0 0 -x1 -x3 x2] [x3*x1 - x2*x0]
[ 0 0 \times0 \times1 -\times3], [ \times1^2 -\times3*\times0]
sage: r = S.resolution(True)
sage: r[0]*r[1]
[0 0 0 0 0]
sage: r[1]*r[2]
[0]
[0]
[0]
[0]
[0]
```

ring()

The ring containing the homogeneous toppling ideal.

OUTPUT:

ring

```
sage: S = sandpiles.Diamond()
sage: S.ring()
```

```
Multivariate Polynomial Ring in x3, x2, x1, x0 over Rational Field sage: S.ring().gens() (x3, x2, x1, x0)
```

Note: The indeterminate xi corresponds to the *i*-th vertex as listed my the method vertices. The term-ordering is degrevlex with indeterminates ordered according to their distance from the sink (larger indeterminates are further from the sink).

show (**kwds)

Draw the underlying graph.

INPUT:

kwds - (optional) arguments passed to the show method for Graph or DiGraph

EXAMPLES:

```
sage: S = Sandpile({0:[], 1:[0,3,4], 2:[0,3,5], 3:[2,5], 4:[1,1], 5:[2,4]})
sage: S.show()
sage: S.show(graph_border=True, edge_labels=True)
```

show3d(**kwds)

Draw the underlying graph.

INPUT:

kwds - (optional) arguments passed to the show method for Graph or DiGraph

EXAMPLES:

```
sage: S = sandpiles.House()
sage: S.show3d() # long time
```

sink()

The sink vertex.

OUTPUT:

sink vertex

EXAMPLES:

```
sage: G = sandpiles.House()
sage: G.sink()
0
sage: H = sandpiles.Grid(2,2)
sage: H.sink()
(0, 0)
sage: type(H.sink())
<... 'tuple'>
```

smith_form()

The Smith normal form for the Laplacian. In detail: a list of integer matrices D, U, V such that ULV = D where L is the transpose of the Laplacian, D is diagonal, and U and V are invertible over the integers.

OUTPUT:

list of integer matrices

```
sage: s = sandpiles.Complete(4)
sage: D,U,V = s.smith_form()
sage: D
[1 0 0 0]
[0 4 0 0]
[0 0 4 0]
[0 0 0 0]
sage: U*s.laplacian()*V == D # Laplacian symmetric => transpose not necessary
True
```

solve()

Approximations of the complex affine zeros of the sandpile ideal.

OUTPUT:

list of complex numbers

EXAMPLES:

Note: The solutions form a multiplicative group isomorphic to the sandpile group. Generators for this group are given exactly by points().

stable configs(smax=None)

Generator for all stable configurations. If smax is provided, then the generator gives all stable configurations less than or equal to smax. If smax does not represent a stable configuration, then each component of smax is replaced by the corresponding component of the maximal stable configuration.

INPUT:

smax – (optional) SandpileConfig or list representing a SandpileConfig

OUTPUT:

generator for all stable configurations

```
sage: s = sandpiles.Complete(3)
sage: a = s.stable_configs()
sage: next(a)
{1: 0, 2: 0}
sage: [i.values() for i in a]
[[0, 1], [1, 0], [1, 1]]
sage: b = s.stable_configs([1,0])
sage: list(b)
[{1: 0, 2: 0}, {1: 1, 2: 0}]
```

stationary_density()

The stationary density of the sandpile.

OUTPUT:

rational number

EXAMPLES:

```
sage: s = sandpiles.Complete(3)
sage: s.stationary_density()
10/9
sage: s = Sandpile(digraphs.DeBruijn(2,2),'00')
sage: s.stationary_density()
9/8
```

Note: The stationary density of a sandpile is the sum $\sum_{c} (\deg(c) + \deg(s))$ where $\deg(s)$ is the degree of the sink and the sum is over all recurrent configurations.

REFERENCES:

• [Lev2014]

superstables (verbose=True)

The superstable configurations. If verbose is False, the configurations are converted to lists of integers. Superstables for undirected graphs are also known as G-parking functions.

INPUT:

```
verbose - (default: True) boolean
```

OUTPUT:

list of SandpileConfig

```
sage: sp = Sandpile(graphs.HouseXGraph(),0).superstables()
sage: sp[:3]
[\{1: 0, 2: 0, 3: 0, 4: 0\}, \{1: 1, 2: 0, 3: 0, 4: 1\}, \{1: 1, 2: 0, 3: 0, 4: 0\}]
sage: sandpiles.Complete(4).superstables(False)
[[0, 0, 0],
 [0, 0, 1],
 [0, 1, 0],
 [1, 0, 0],
 [0, 0, 2],
 [0, 2, 0],
 [2, 0, 0],
 [0, 1, 1],
 [1, 0, 1],
 [1, 1, 0],
 [0, 1, 2],
 [0, 2, 1],
 [1, 0, 2],
 [1, 2, 0],
 [2, 0, 1],
 [2, 1, 0]]
sage: sandpiles.Cycle(4).superstables(False)
[[0, 0, 0], [1, 0, 0], [0, 1, 0], [0, 0, 1]]
```

symmetric recurrents(orbits)

The symmetric recurrent configurations.

INPUT:

orbits - list of lists partitioning the vertices

OUTPUT:

list of recurrent configurations

EXAMPLES:

```
sage: S = Sandpile({0: {},
                   1: {0: 1, 2: 1, 3: 1},
...:
                   2: {1: 1, 3: 1, 4: 1},
. . . . :
                   3: {1: 1, 2: 1, 4: 1},
. . . . :
                   4: {2: 1, 3: 1}})
sage: S.symmetric_recurrents([[1],[2,3],[4]])
[\{1: 2, 2: 2, 3: 2, 4: 1\}, \{1: 2, 2: 2, 3: 2, 4: 0\}]
sage: S.recurrents()
[\{1: 2, 2: 2, 3: 2, 4: 1\},
{1: 2, 2: 2, 3: 2, 4: 0},
{1: 2, 2: 1, 3: 2, 4: 0},
{1: 2, 2: 2, 3: 0, 4: 1},
 \{1: 2, 2: 0, 3: 2, 4: 1\},\
 {1: 2, 2: 2, 3: 1, 4: 0},
{1: 2, 2: 1, 3: 2, 4: 1},
 {1: 2, 2: 2, 3: 1, 4: 1}]
```

Note: The user is responsible for ensuring that the list of orbits comes from a group of symmetries of the underlying graph.

tutte polynomial()

The Tutte polynomial of the underlying graph. Only defined for undirected sandpile graphs.

OUTPUT:

polynomial

EXAMPLES:

```
sage: s = sandpiles.Complete(4)
sage: s.tutte_polynomial()
x^3 + y^3 + 3*x^2 + 4*x*y + 3*y^2 + 2*x + 2*y
sage: s.tutte_polynomial().subs(x=1)
y^3 + 3*y^2 + 6*y + 6
sage: s.tutte_polynomial().subs(x=1).coefficients() == s.h_vector()
True
```

unsaturated_ideal()

The unsaturated, homogeneous toppling ideal.

OUTPUT:

ideal

```
sage: S = sandpiles.Diamond()
sage: S.unsaturated_ideal().gens()
[x1^3 - x3*x2*x0, x2^3 - x3*x1*x0, x3^2 - x2*x1]
sage: S.ideal().gens()
[x2*x1 - x0^2, x3^2 - x0^2, x1^3 - x3*x2*x0, x3*x1^2 - x2^2*x0, x2^3 - x3*x1*x0, x3*x2^2 - x1^2*x0]
```

static version()

The version number of Sage Sandpiles.

OUTPUT:

string

EXAMPLES:

```
sage: Sandpile.version()
Sage Sandpiles Version 2.4
sage: S = sandpiles.Complete(3)
sage: S.version()
Sage Sandpiles Version 2.4
```

zero_config()

The all-zero configuration.

OUTPUT:

SandpileConfig

EXAMPLES:

```
sage: s = sandpiles.Diamond()
sage: s.zero_config()
{1: 0, 2: 0, 3: 0}
```

zero_div()

The all-zero divisor.

OUTPUT:

SandpileDivisor

EXAMPLES:

```
sage: S = sandpiles.House()
sage: S.zero_div()
{0: 0, 1: 0, 2: 0, 3: 0, 4: 0}
```

class sage.sandpiles.sandpile.SandpileConfig (S, c)

Bases: dict

Class for configurations on a sandpile.

```
add_random(distrib=None)
```

Add one grain of sand to a random vertex. Optionally, a probability distribution, distrib, may be placed on the vertices or the nonsink vertices. See NOTE for details.

INPUT:

distrib – (optional) list of nonnegative numbers summing to 1 (representing a prob. dist.)

OUTPUT:

SandpileConfig

EXAMPLES:

```
sage: s = sandpiles.Complete(4)
sage: c = s.zero_config()
sage: c.add_random() # random
{1: 0, 2: 1, 3: 0}
sage: c
{1: 0, 2: 0, 3: 0}
sage: c.add_random([0.1,0.1,0.8]) # random
{1: 0, 2: 0, 3: 1}
sage: c.add_random([0.7,0.1,0.1]) # random
{1: 0, 2: 0, 3: 0}
```

We compute the "sizes" of the avalanches caused by adding random grains of sand to the maximal stable configuration on a grid graph. The function stabilize() returns the firing vector of the stabilization, a dictionary whose values say how many times each vertex fires in the stabilization.:

Note: If distrib is None, then the probability is the uniform probability on the nonsink vertices. Otherwise, there are two possibilities:

- (i) the length of distrib is equal to the number of vertices, and distrib represents a probability distribution on all of the vertices. In that case, the sink may be chosen at random, in which case, the configuration is unchanged.
- (ii) Otherwise, the length of distrib must be equal to the number of nonsink vertices, and distrib represents a probability distribution on the nonsink vertices.

Warning: If distrib != None, the user is responsible for assuring the sum of its entries is 1 and that its length is equal to the number of sink vertices or the number of nonsink vertices.

burst size(v)

The burst size of the configuration with respect to the given vertex.

INPUT:

v – vertex

OUTPUT:

integer

EXAMPLES:

```
sage: s = sandpiles.Diamond()
sage: [i.burst_size(0) for i in s.recurrents()]
[1, 1, 1, 1, 1, 1, 1]
sage: [i.burst_size(1) for i in s.recurrents()]
[0, 0, 1, 2, 1, 2, 0, 2]
```

Note: To define c.burst (v), if v is not the sink, let c' be the unique recurrent for which the stabilization of c' + v is c. The burst size is then the amount of sand that goes into the sink during this stabilization. If v is the sink, the burst size is defined to be 1.

REFERENCES:

• [Lev2014]

deg()

The degree of the configuration.

OUTPUT:

integer

EXAMPLES:

```
sage: S = sandpiles.Complete(3)
sage: c = SandpileConfig(S, [1,2])
sage: c.deg()
3
```

dualize()

The difference with the maximal stable configuration.

OUTPUT:

SandpileConfig

EXAMPLES:

```
sage: S = sandpiles.Cycle(3)
sage: c = SandpileConfig(S, [1,2])
sage: S.max_stable()
{1: 1, 2: 1}
sage: c.dualize()
{1: 0, 2: -1}
sage: S.max_stable() - c == c.dualize()
True
```

equivalent_recurrent (with_firing_vector=False)

The recurrent configuration equivalent to the given configuration. Optionally, return the corresponding firing vector.

INPUT:

```
with_firing_vector - (default: False) boolean
```

OUTPUT:

SandpileConfig or [SandpileConfig, firing_vector]

```
sage: S = sandpiles.Diamond()
sage: c = SandpileConfig(S, [0,0,0])
sage: c.equivalent_recurrent() == S.identity()
True
sage: x = c.equivalent_recurrent(True)
sage: r = vector([x[0][v] for v in S.nonsink_vertices()])
sage: f = vector([x[1][v] for v in S.nonsink_vertices()])
sage: cv = vector(c.values())
sage: r == cv - f*S.reduced_laplacian()
True
```

Note: Let L be the reduced Laplacian, c the initial configuration, r the returned configuration, and f the firing vector. Then $r = c - f \cdot L$.

equivalent_superstable (with_firing_vector=False)

The equivalent superstable configuration. Optionally, return the corresponding firing vector.

INPUT:

```
with_firing_vector - (default: False) boolean
```

OUTPUT:

SandpileConfig or [SandpileConfig, firing_vector]

EXAMPLES:

```
sage: S = sandpiles.Diamond()
sage: m = S.max_stable()
sage: m.equivalent_superstable().is_superstable()
True
sage: x = m.equivalent_superstable(True)
sage: s = vector(x[0].values())
sage: f = vector(x[1].values())
sage: mv = vector(m.values())
sage: s == mv - f*S.reduced_laplacian()
True
```

Note: Let L be the reduced Laplacian, c the initial configuration, s the returned configuration, and f the firing vector. Then $s = c - f \cdot L$.

fire_script (sigma)

Fire the given script. In other words, fire each vertex the number of times indicated by sigma.

INPUT:

sigma - SandpileConfig or (list or dict representing a SandpileConfig)

OUTPUT:

SandpileConfig

```
sage: S = sandpiles.Cycle(4)
sage: c = SandpileConfig(S, [1,2,3])
sage: c.unstable()
```

fire unstable()

Fire all unstable vertices.

OUTPUT:

SandpileConfig

EXAMPLES:

```
sage: S = sandpiles.Cycle(4)
sage: c = SandpileConfig(S, [1,2,3])
sage: c.fire_unstable()
{1: 2, 2: 1, 3: 2}
```

fire_vertex(v)

Fire the given vertex.

INPUT:

v - vertex

OUTPUT:

SandpileConfig

EXAMPLES:

```
sage: S = sandpiles.Cycle(3)
sage: c = SandpileConfig(S, [1,2])
sage: c.fire_vertex(2)
{1: 2, 2: 0}
```

static help(verbose=True)

List of SandpileConfig methods. If verbose, include short descriptions.

INPUT:

verbose - (default: True) boolean

OUTPUT:

printed string

```
-- Add one grain of sand to a random vertex.
add_random
                      -- The burst size of the configuration with respect to.
burst size

→the given vertex.

                     -- The degree of the configuration.
deg
                     -- The difference with the maximal stable.
dualize

→configuration.

equivalent_recurrent -- The recurrent configuration equivalent to the given_

→configuration.

equivalent_superstable -- The equivalent superstable configuration.
                     -- Fire the given script.
fire_script
                     -- Fire all unstable vertices.
fire_unstable
fire_vertex
                     -- Fire the given vertex.
help
                      -- List of SandpileConfig methods.
is_recurrent
                      -- Is the configuration recurrent?
is_stable
                      -- Is the configuration stable?
is_superstable
                    -- Is the configuration superstable?
is_symmetric
                     -- Is the configuration symmetric?
order
                     -- The order of the equivalent recurrent element.
sandpile
                     -- The configuration's underlying sandpile.
                     -- Show the configuration.
show
stabilize
                     -- The stabilized configuration.
support
                     -- The vertices containing sand.
                     -- The unstable vertices.
unstable
                     -- The values of the configuration as a list.
values
```

is_recurrent()

Is the configuration recurrent?

OUTPUT:

boolean

EXAMPLES:

```
sage: S = sandpiles.Diamond()
sage: S.identity().is_recurrent()
True
sage: S.zero_config().is_recurrent()
False
```

is_stable()

Is the configuration stable?

OUTPUT:

boolean

EXAMPLES:

```
sage: S = sandpiles.Diamond()
sage: S.max_stable().is_stable()
True
sage: (2*S.max_stable()).is_stable()
False
sage: (S.max_stable() & S.max_stable()).is_stable()
True
```

is_superstable()

Is the configuration superstable?

OUTPUT:

boolean

EXAMPLES:

```
sage: S = sandpiles.Diamond()
sage: S.zero_config().is_superstable()
True
```

is_symmetric(orbits)

Is the configuration symmetric? Return True if the values of the configuration are constant over the vertices in each sublist of orbits.

INPUT:

orbits - list of lists of vertices

OUTPUT:

boolean

EXAMPLES:

order()

The order of the equivalent recurrent element.

OUTPUT:

integer

EXAMPLES:

```
sage: S = sandpiles.Diamond()
sage: c = SandpileConfig(S,[2,0,1])
sage: c.order()
4
sage: ~(c + c + c + c) == S.identity()
True
sage: c = SandpileConfig(S,[1,1,0])
sage: c.order()
1
sage: c.is_recurrent()
False
sage: c.equivalent_recurrent() == S.identity()
True
```

sandpile()

The configuration's underlying sandpile.

OUTPUT:

Sandpile

EXAMPLES:

```
sage: S = sandpiles.Diamond()
sage: c = S.identity()
sage: c.sandpile()
Diamond sandpile graph: 4 vertices, sink = 0
sage: c.sandpile() == S
True
```

show (*sink=True*, *colors=True*, *heights=False*, *directed=None*, **kwds) Show the configuration.

show the configuration

INPUT:

- sink (default: True) whether to show the sink
- colors (default: True) whether to color-code the amount of sand on each vertex
- heights (default: False) whether to label each vertex with the amount of sand
- directed (optional) whether to draw directed edges
- kwds (optional) arguments passed to the show method for Graph

EXAMPLES:

```
sage: S = sandpiles.Diamond()
sage: c = S.identity()
sage: c.show()
sage: c.show(directed=False)
sage: c.show(sink=False,colors=False,heights=True)
```

stabilize (with_firing_vector=False)

The stabilized configuration. Optionally returns the corresponding firing vector.

INPUT:

```
with_firing_vector - (default: False) boolean
```

OUTPUT:

SandpileConfig or [SandpileConfig, firing_vector]

EXAMPLES:

```
sage: S = sandpiles.House()
sage: c = 2*S.max_stable()
sage: c._set_stabilize()
sage: '_stabilize' in c.__dict__
True
sage: S = sandpiles.House()
sage: c = S.max_stable() + S.identity()
sage: c.stabilize(True)
[{1: 1, 2: 2, 3: 2, 4: 1}, {1: 2, 2: 2, 3: 3, 4: 3}]
sage: S.max_stable() & S.identity() == c.stabilize()
True
sage: ~c == c.stabilize()
True
```

support()

The vertices containing sand.

OUTPUT:

list - support of the configuration

EXAMPLES:

```
sage: S = sandpiles.Diamond()
sage: c = S.identity()
sage: c
{1: 2, 2: 2, 3: 0}
sage: c.support()
[1, 2]
```

unstable()

The unstable vertices.

OUTPUT:

list of vertices

EXAMPLES:

```
sage: S = sandpiles.Cycle(4)
sage: c = SandpileConfig(S, [1,2,3])
sage: c.unstable()
[2, 3]
```

values()

The values of the configuration as a list. The list is sorted in the order of the vertices.

OUTPUT:

list of integers

boolean

EXAMPLES:

```
sage: S = Sandpile({'a':[1,'b'], 'b':[1,'a'], 1:['a']},'a')
sage: c = SandpileConfig(S, {'b':1, 1:2})
sage: c
{1: 2, 'b': 1}
sage: c.values()
[2, 1]
sage: S.nonsink_vertices()
[1, 'b']
```

```
class sage.sandpiles.sandpile.SandpileDivisor(S, D)
```

Bases: dict

Class for divisors on a sandpile.

Dcomplex()

The support-complex. (See NOTE.)

OUTPUT:

simplicial complex

```
sage: S = sandpiles.House()
sage: p = SandpileDivisor(S, [1,2,1,0,0]).Dcomplex()
sage: p.homology()
{0: 0, 1: Z x Z, 2: 0}
```

```
sage: p.f_vector()
[1, 5, 10, 4]
sage: p.betti()
{0: 1, 1: 2, 2: 0}
```

Note: The "support-complex" is the simplicial complex determined by the supports of the linearly equivalent effective divisors.

add random(distrib=None)

Add one grain of sand to a random vertex.

INPUT:

distrib – (optional) list of nonnegative numbers representing a probability distribution on the vertices

OUTPUT:

SandpileDivisor

EXAMPLES:

```
sage: s = sandpiles.Complete(4)
sage: D = s.zero_div()
sage: D.add_random() # random
{0: 0, 1: 0, 2: 1, 3: 0}
sage: D.add_random([0.1,0.1,0.7]) # random
{0: 0, 1: 0, 2: 0, 3: 1}
```

Warning: If distrib is not None, the user is responsible for assuring the sum of its entries is 1.

betti()

The Betti numbers for the support-complex. (See NOTE.)

OUTPUT:

dictionary of integers

EXAMPLES:

```
sage: S = sandpiles.Cycle(3)
sage: D = SandpileDivisor(S, [2,0,1])
sage: D.betti()
{0: 1, 1: 1}
```

Note: The "support-complex" is the simplicial complex determined by the supports of the linearly equivalent effective divisors.

deg()

The degree of the divisor.

OUTPUT:

integer

```
sage: S = sandpiles.Cycle(3)
sage: D = SandpileDivisor(S, [1,2,3])
sage: D.deg()
6
```

dualize()

The difference with the maximal stable divisor.

OUTPUT:

SandpileDivisor

EXAMPLES:

```
sage: S = sandpiles.Cycle(3)
sage: D = SandpileDivisor(S, [1,2,3])
sage: D.dualize()
{0: 0, 1: -1, 2: -2}
sage: S.max_stable_div() - D == D.dualize()
True
```

effective_div (*verbose=True*, *with_firing_vectors=False*)

All linearly equivalent effective divisors. If verbose is False, the divisors are converted to lists of integers. If with_firing_vectors is True then a list of firing vectors is also given, each of which prescribes the vertices to be fired in order to obtain an effective divisor.

INPUT:

- verbose (default: True) boolean
- with_firing_vectors (default: False) boolean

OUTPUT:

list (of divisors)

```
sage: s = sandpiles.Complete(4)
sage: D = SandpileDivisor(s, [4,2,0,0])
sage: sorted(D.effective_div(), key=str)
[{0: 0, 1: 2, 2: 0, 3: 4},
\{0: 0, 1: 2, 2: 4, 3: 0\},\
\{0: 0, 1: 6, 2: 0, 3: 0\},\
\{0: 1, 1: 3, 2: 1, 3: 1\},\
\{0: 2, 1: 0, 2: 2, 3: 2\},\
{0: 4, 1: 2, 2: 0, 3: 0}]
sage: sorted(D.effective_div(False))
[[0, 2, 0, 4],
 [0, 2, 4, 0],
 [0, 6, 0, 0],
 [1, 3, 1, 1],
 [2, 0, 2, 2],
 [4, 2, 0, 0]]
sage: sorted(D.effective_div(with_firing_vectors=True), key=str)
[(\{0: 0, 1: 2, 2: 0, 3: 4\}, (0, -1, -1, -2)),
 (\{0: 0, 1: 2, 2: 4, 3: 0\}, (0, -1, -2, -1)),
 (\{0: 0, 1: 6, 2: 0, 3: 0\}, (0, -2, -1, -1)),
 (\{0: 1, 1: 3, 2: 1, 3: 1\}, (0, -1, -1, -1)),
 (\{0: 2, 1: 0, 2: 2, 3: 2\}, (0, 0, -1, -1)),
 ({0: 4, 1: 2, 2: 0, 3: 0}, (0, 0, 0, 0))]
```

```
sage: a = _[2]
sage: a[0].values()
[0, 6, 0, 0]
sage: vector(D.values()) - s.laplacian()*a[1]
(0, 6, 0, 0)
sage: sorted(D.effective_div(False, True))
[([0, 2, 0, 4], (0, -1, -1, -2)),
([0, 2, 4, 0], (0, -1, -2, -1)),
([0, 6, 0, 0], (0, -2, -1, -1)),
([1, 3, 1, 1], (0, -1, -1, -1)),
([2, 0, 2, 2], (0, 0, -1, -1)),
([4, 2, 0, 0], (0, 0, 0, 0))]
sage: D = SandpileDivisor(s, [-1, 0, 0, 0])
sage: D.effective_div(False, True)
[]
```

fire_script (sigma)

Fire the given script. In other words, fire each vertex the number of times indicated by sigma.

INPUT

sigma – SandpileDivisor or (list or dict representing a SandpileDivisor)

OUTPUT:

SandpileDivisor

EXAMPLES:

fire_unstable()

Fire all unstable vertices.

OUTPUT:

SandpileDivisor

EXAMPLES:

```
sage: S = sandpiles.Cycle(3)
sage: D = SandpileDivisor(S, [1,2,3])
sage: D.fire_unstable()
{0: 3, 1: 1, 2: 2}
```

fire_vertex(v)

Fire the given vertex.

INPUT:

v - vertex

OUTPUT:

SandpileDivisor

EXAMPLES:

```
sage: S = sandpiles.Cycle(3)
sage: D = SandpileDivisor(S, [1,2,3])
sage: D.fire_vertex(1)
{0: 2, 1: 0, 2: 4}
```

static help(verbose=True)

List of SandpileDivisor methods. If verbose, include short descriptions.

INPUT:

```
verbose - (default: True) boolean
```

OUTPUT:

printed string

EXAMPLES:

```
sage: SandpileDivisor.help()
For detailed help with any method FOO listed below,
enter "SandpileDivisor.FOO?" or enter "D.FOO?" for any SandpileDivisor D.
Dcomplex
                     -- The support-complex.
add_random
                    -- Add one grain of sand to a random vertex.
betti
                     -- The Betti numbers for the support-complex.
                     -- The degree of the divisor.
deg
                     -- The difference with the maximal stable divisor.
dualize
effective_div
                    -- All linearly equivalent effective divisors.
                     -- Fire the given script.
fire_script
fire_unstable
                     -- Fire all unstable vertices.
fire_vertex
                     -- Fire the given vertex.
help
                     -- List of SandpileDivisor methods.
is_alive
                     -- Is the divisor stabilizable?
is_linearly_equivalent -- Is the given divisor linearly equivalent?
is_q_reduced -- Is the divisor q-reduced?
is_symmetric
                    -- Is the divisor symmetric?
is_weierstrass_pt -- Is the given vertex a Weierstrass point?
polytope
                    -- The polytope determining the complete linear system.
polytope_integer_pts -- The integer points inside divisor's polytope.
                    -- The linearly equivalent q-reduced divisor.
q_reduced
                     -- The rank of the divisor.
rank
                     -- The divisor's underlying sandpile.
sandpile
show
                     -- Show the divisor.
simulate_threshold
                      -- The first unstabilizable divisor in the closed.
→Markov chain.
stabilize
                      -- The stabilization of the divisor.
support
                     -- List of vertices at which the divisor is nonzero.
unstable
                     -- The unstable vertices.
values
                     -- The values of the divisor as a list.
weierstrass_div -- The Weierstrass divisor.
weierstrass_gap_seq -- The Weierstrass gap sequence at the given vertex.
weierstrass_pts -- The Weierstrass points (vertices).
weierstrass_rank_seq -- The Weierstrass rank sequence at the given vertex.
```

is_alive (cycle=False)

Is the divisor stabilizable? In other words, will the divisor stabilize under repeated firings of all unstable vertices? Optionally returns the resulting cycle.

INPUT:

cycle - (default: False) boolean

OUTPUT:

boolean or optionally, a list of SandpileDivisors

EXAMPLES:

```
sage: S = sandpiles.Complete(4)
sage: D = SandpileDivisor(S, {0: 4, 1: 3, 2: 3, 3: 2})
sage: D.is_alive()
True
sage: D.is_alive(True)
[{0: 4, 1: 3, 2: 3, 3: 2}, {0: 3, 1: 2, 2: 2, 3: 5}, {0: 1, 1: 4, 2: 4, 3: 3}]
```

is_linearly_equivalent (D, with_firing_vector=False)

Is the given divisor linearly equivalent? Optionally, returns the firing vector. (See NOTE.)

INPUT:

- D SandpileDivisor or list, tuple, etc. representing a divisor
- with_firing_vector (default: False) boolean

OUTPUT:

boolean or integer vector

EXAMPLES:

```
sage: s = sandpiles.Complete(3)
sage: D = SandpileDivisor(s,[2,0,0])
sage: D.is_linearly_equivalent([0,1,1])
True
sage: D.is_linearly_equivalent([0,1,1],True)
(1, 0, 0)
sage: v = vector(D.is_linearly_equivalent([0,1,1],True))
sage: vector(D.values()) - s.laplacian()*v
(0, 1, 1)
sage: D.is_linearly_equivalent([0,0,0])
False
sage: D.is_linearly_equivalent([0,0,0],True)
()
```

Note:

- If with_firing_vector is False, returns either True or False.
- If with_firing_vector is True then: (i) if self is linearly equivalent to D, returns a vector v such that self v*self.laplacian().transpose() = D. Otherwise, (ii) if self is not linearly equivalent to D, the output is the empty vector, ().

is_q_reduced()

Is the divisor q-reduced? This would mean that self = c + kq where c is superstable, k is an integer, and q is the sink vertex.

OUTPUT:

boolean

EXAMPLES:

```
sage: s = sandpiles.Complete(4)
sage: D = SandpileDivisor(s,[2,-3,2,0])
sage: D.is_q_reduced()
False
sage: SandpileDivisor(s,[10,0,1,2]).is_q_reduced()
True
```

For undirected or, more generally, Eulerian graphs, *q*-reduced divisors are linearly equivalent if and only if they are equal. The same does not hold for general directed graphs:

```
sage: s = Sandpile({0:[1],1:[1,1]})
sage: D = SandpileDivisor(s,[-1,1])
sage: Z = s.zero_div()
sage: D.is_q_reduced()
True
sage: Z.is_q_reduced()
True
sage: D == Z
False
sage: D.is_linearly_equivalent(Z)
True
```

is_symmetric(orbits)

Is the divisor symmetric? Return True if the values of the configuration are constant over the vertices in each sublist of orbits.

INPUT:

orbits - list of lists of vertices

OUTPUT:

boolean

EXAMPLES:

```
sage: S = sandpiles.House()
sage: S.dict()
{0: {1: 1, 2: 1},
    1: {0: 1, 3: 1},
    2: {0: 1, 3: 1, 4: 1},
    3: {1: 1, 2: 1, 4: 1},
    4: {2: 1, 3: 1}}
sage: D = SandpileDivisor(S, [0,0,1,1,3])
sage: D.is_symmetric([[2,3], [4]])
True
```

is weierstrass pt (v='sink')

Is the given vertex a Weierstrass point?

INPUT:

v - (default: sink) vertex

OUTPUT:

boolean

```
sage: s = sandpiles.House()
sage: K = s.canonical_divisor()
sage: K.weierstrass_rank_seq() # sequence at the sink vertex, 0
(1, 0, -1)
sage: K.is_weierstrass_pt()
False
sage: K.weierstrass_rank_seq(4)
(1, 0, 0, -1)
sage: K.is_weierstrass_pt(4)
True
```

Note: The vertex v is a (generalized) Weierstrass point for divisor D if the sequence of ranks r(D-nv) for $n=0,1,2,\ldots$ is not $r(D),r(D)-1,\ldots,0,-1,-1,\ldots$

polytope()

The polytope determining the complete linear system.

OUTPUT:

polytope

EXAMPLES:

```
sage: s = sandpiles.Complete(4)
sage: D = SandpileDivisor(s,[4,2,0,0])
sage: p = D.polytope()
sage: p.inequalities()
(An inequality (-3, 1, 1) x + 2 >= 0,
   An inequality (1, 1, 1) x + 4 >= 0,
   An inequality (1, -3, 1) x + 0 >= 0,
   An inequality (1, 1, -3) x + 0 >= 0)
sage: D = SandpileDivisor(s,[-1,0,0,0])
sage: D.polytope()
The empty polyhedron in QQ^3
```

Note: For a divisor D, this is the intersection of (i) the polyhedron determined by the system of inequalities $L^t x \leq D$ where L^t is the transpose of the Laplacian with (ii) the hyperplane $x_{\text{sink_vertex}} = 0$. The polytope is thought of as sitting in (n-1)-dimensional Euclidean space where n is the number of vertices.

polytope_integer_pts()

The integer points inside divisor's polytope. The polytope referred to here is the one determining the divisor's complete linear system (see the documentation for polytope).

OUTPUT:

tuple of integer vectors

```
(0, -1, -1),

(0, 0, 0)]

sage: D = SandpileDivisor(s,[-1,0,0,0])

sage: D.polytope_integer_pts()

()
```

q_reduced (verbose=True)

The linearly equivalent q-reduced divisor.

INPUT:

```
verbose - (default: True) boolean
```

OUTPUT:

SandpileDivisor or list representing SandpileDivisor

EXAMPLES:

```
sage: s = sandpiles.Complete(4)
sage: D = SandpileDivisor(s,[2,-3,2,0])
sage: D.q_reduced()
{0: -2, 1: 1, 2: 2, 3: 0}
sage: D.q_reduced(False)
[-2, 1, 2, 0]
```

Note: The divisor D is qreducedif'D = c + kq where c is superstable, k is an integer, and q is the sink.

rank (with_witness=False)

The rank of the divisor. Optionally returns an effective divisor E such that D-E is not winnable (has an empty complete linear system).

INPUT:

```
with_witness - (default: False) boolean
```

OUTPUT:

integer or (integer, SandpileDivisor)

```
sage: S = sandpiles.Complete(4)
sage: D = SandpileDivisor(S,[4,2,0,0])
sage: D.rank()
3
sage: D.rank(True)
(3, {0: 3, 1: 0, 2: 1, 3: 0})
sage: E = _[1]
sage: (D - E).rank()
-1

Riemann-Roch theorem::
sage: D.rank() - (S.canonical_divisor()-D).rank() == D.deg() + 1 - S.

--genus()
True

Riemann-Roch theorem::
```

```
sage: D.rank() - (S.canonical_divisor()-D).rank() == D.deg() + 1 - S.
⇒genus()
  True
  sage: S = Sandpile(\{0:[1,1,1,2],1:[0,0,0,1,1,1,2,2],2:[2,2,1,1,0]\},0) \#
→multigraph with loops
  sage: D = SandpileDivisor(S, [4, 2, 0])
  sage: D.rank (True)
  (2, \{0: 1, 1: 1, 2: 1\})
  sage: S = Sandpile({0:[1,2], 1:[0,2,2], 2: [0,1]},0) # directed graph
  sage: S.is_undirected()
  False
  sage: D = SandpileDivisor(S, [0, 2, 0])
  sage: D.effective_div()
  [\{0: 0, 1: 2, 2: 0\}, \{0: 2, 1: 0, 2: 0\}]
  sage: D.rank(True)
  (0, \{0: 0, 1: 0, 2: 1\})
  sage: E = D.rank(True)[1]
  sage: (D - E).effective_div()
```

Note: The rank of a divisor D is -1 if D is not linearly equivalent to an effective divisor (i.e., the dollar game represented by D is unwinnable). Otherwise, the rank of D is the largest integer r such that D - E is linearly equivalent to an effective divisor for all effective divisors E with $\deg(E) = r$.

sandpile()

The divisor's underlying sandpile.

OUTPUT:

Sandpile

EXAMPLES:

```
sage: S = sandpiles.Diamond()
sage: D = SandpileDivisor(S,[1,-2,0,3])
sage: D.sandpile()
Diamond sandpile graph: 4 vertices, sink = 0
sage: D.sandpile() == S
True
```

show (heights=True, directed=None, **kwds)

Show the divisor.

INPUT:

- heights (default: True) whether to label each vertex with the amount of sand
- directed (optional) whether to draw directed edges
- kwds (optional) arguments passed to the show method for Graph

```
sage: S = sandpiles.Diamond()
sage: D = SandpileDivisor(S,[1,-2,0,2])
sage: D.show(graph_border=True,vertex_size=700,directed=False)
```

simulate threshold(distrib=None)

The first unstabilizable divisor in the closed Markov chain. (See NOTE.)

INPUT:

distrib - (optional) list of nonnegative numbers representing a probability distribution on the vertices

OUTPUT:

SandpileDivisor

EXAMPLES:

```
sage: s = sandpiles.Complete(4)
sage: D = s.zero_div()
sage: D.simulate_threshold() # random
{0: 2, 1: 3, 2: 1, 3: 2}
sage: n(mean([D.simulate_threshold().deg() for _ in range(10)])) # random
7.10000000000000
sage: n(s.stationary_density()*s.num_verts())
6.93750000000000
```

Note: Starting at self, repeatedly choose a vertex and add a grain of sand to it. Return the first unstabilizable divisor that is reached. Also see the markov_chain method for the underlying sandpile.

stabilize (with_firing_vector=False)

The stabilization of the divisor. If not stabilizable, return an error.

INPUT:

with firing vector - (default: False) boolean

EXAMPLES:

```
sage: s = sandpiles.Complete(4)
sage: D = SandpileDivisor(s,[0,3,0,0])
sage: D.stabilize()
{0: 1, 1: 0, 2: 1, 3: 1}
sage: D.stabilize(with_firing_vector=True)
[{0: 1, 1: 0, 2: 1, 3: 1}, {0: 0, 1: 1, 2: 0, 3: 0}]
```

support()

List of vertices at which the divisor is nonzero.

OUTPUT:

list representing the support of the divisor

EXAMPLES:

```
sage: S = sandpiles.Cycle(4)
sage: D = SandpileDivisor(S, [0,0,1,1])
sage: D.support()
[2, 3]
sage: S.vertices()
[0, 1, 2, 3]
```

unstable()

The unstable vertices.

OUTPUT:

list of vertices

EXAMPLES:

```
sage: S = sandpiles.Cycle(3)
sage: D = SandpileDivisor(S, [1,2,3])
sage: D.unstable()
[1, 2]
```

values()

The values of the divisor as a list. The list is sorted in the order of the vertices.

OUTPUT:

list of integers

boolean

EXAMPLES:

```
sage: S = Sandpile({'a':[1,'b'], 'b':[1,'a'], 1:['a']},'a')
sage: D = SandpileDivisor(S, {'a':0, 'b':1, 1:2})
sage: D
{'a': 0, 1: 2, 'b': 1}
sage: D.values()
[2, 0, 1]
sage: S.vertices()
[1, 'a', 'b']
```

weierstrass_div (verbose=True)

The Weierstrass divisor. Its value at a vertex is the weight of that vertex as a Weierstrass point. (See SandpileDivisor.weierstrass_gap_seq.)

INPUT:

verbose - (default: True) boolean

OUTPUT:

SandpileDivisor

EXAMPLES:

```
sage: s = sandpiles.Diamond()
sage: D = SandpileDivisor(s, [4,2,1,0])
sage: [D.weierstrass_rank_seq(v) for v in s]
[(5, 4, 3, 2, 1, 0, 0, -1),
    (5, 4, 3, 2, 1, 0, 0, 0, -1),
    (5, 4, 3, 2, 1, 0, 0, 0, -1),
    (5, 4, 3, 2, 1, 0, 0, 0, -1)]
sage: D.weierstrass_div()
{0: 1, 1: 0, 2: 2, 3: 1}
sage: k5 = sandpiles.Complete(5)
sage: K = k5.canonical_divisor()
sage: K.weierstrass_div()
{0: 9, 1: 9, 2: 9, 3: 9, 4: 9}
```

weierstrass_gap_seq(v='sink', weight=True)

The Weierstrass gap sequence at the given vertex. If weight is True, then also compute the weight of each gap value.

INPUT:

- v (default: sink) vertex
- weight (default: True) boolean

OUTPUT:

list or (list of list) of integers

EXAMPLES:

```
sage: s = sandpiles.Cycle(4)
sage: D = SandpileDivisor(s,[2,0,0,0])
sage: [D.weierstrass_gap_seq(v,False) for v in s.vertices()]
[(1, 3), (1, 2), (1, 3), (1, 2)]
sage: [D.weierstrass_gap_seq(v) for v in s.vertices()]
[((1, 3), 1), ((1, 2), 0), ((1, 3), 1), ((1, 2), 0)]
sage: D.weierstrass_gap_seq() # gap sequence at sink vertex, 0
((1, 3), 1)
sage: D.weierstrass_rank_seq() # rank sequence at the sink vertex
(1, 0, 0, -1)
```

Note: The integer k is a Weierstrass gap for the divisor D at vertex v if the rank of D-(k-1)v does not equal the rank of D-kv. Let r be the rank of D and let k_i be the i-th gap at v. The Weierstrass weight of v for D is the sum of (k_i-i) as i ranges from 1 to v + 1. It measure the difference between the sequence v, v - 1, ..., 0, -1, -1, ... and the rank sequence v rank v - 1, rank v - 1, rank v - 2, rank v - 2, ...

weierstrass_pts (with_rank_seq=False)

The Weierstrass points (vertices). Optionally, return the corresponding rank sequences.

INPUT:

```
with_rank_seq-(default: False) boolean
```

OUTPUT:

tuple of vertices or list of (vertex, rank sequence)

EXAMPLES:

```
sage: s = sandpiles.House()
sage: K = s.canonical_divisor()
sage: K.weierstrass_pts()
(4,)
sage: K.weierstrass_pts(True)
[(4, (1, 0, 0, -1))]
```

Note: The vertex v is a (generalized) Weierstrass point for divisor D if the sequence of ranks r(D-nv) for $n=0,1,2,\ldots$ 'is not $r(D),r(D)-1,\ldots,0,-1,-1,\ldots$

weierstrass_rank_seq(v='sink')

The Weierstrass rank sequence at the given vertex. Computes the rank of the divisor D - nv starting with n = 0 and ending when the rank is -1.

INPUT:

```
v - (default: sink) vertex
```

OUTPUT:

tuple of int

EXAMPLES:

```
sage: s = sandpiles.House()
sage: K = s.canonical_divisor()
sage: [K.weierstrass_rank_seq(v) for v in s.vertices()]
[(1, 0, -1), (1, 0, -1), (1, 0, -1), (1, 0, -1), (1, 0, 0, -1)]
```

sage.sandpiles.sandpile.admissible_partitions (S, k)

The partitions of the vertices of S into k parts, each of which is connected.

INPUT:

S – Sandpile

k - integer

OUTPUT:

list of partitions

EXAMPLES:

```
sage: from sage.sandpiles.sandpile import admissible_partitions
sage: from sage.sandpiles.sandpile import partition_sandpile
sage: S = sandpiles.Cycle(4)
sage: P = [admissible_partitions(S, i) for i in [2,3,4]]
sage: P
[[{{0}, {1, 2, 3}},
  \{\{0, 2, 3\}, \{1\}\},\
  {{0, 1, 3}, {2}},
  \{\{0, 1, 2\}, \{3\}\},\
  \{\{0, 1\}, \{2, 3\}\},\
 \{\{0, 3\}, \{1, 2\}\}\}
 [{{0}, {1}, {2, 3}},
 {{0}, {1, 2}, {3}},
 {{0, 3}, {1}, {2}},
 \{\{0, 1\}, \{2\}, \{3\}\}\}
[{{0}, {1}, {2}, {3}}]]
sage: for p in P:
....: sum([partition_sandpile(S, i).betti(verbose=False)[-1] for i in p])
8
3
sage: S.betti()
     0
                1
    0:
         1
   1:
                       8
                              3
                 6
total: 1
               6
                      8
                            3
```

sage.sandpiles.sandpile.aztec_sandpile(n)

The aztec diamond graph.

INPUT:

n - integer

OUTPUT:

dictionary for the aztec diamond graph

EXAMPLES:

```
sage: from sage.sandpiles.sandpile import aztec_sandpile
sage: aztec_sandpile(2)
{ 'sink': { (-3/2, -1/2): 2, }
  (-3/2, 1/2): 2,
  (-1/2, -3/2): 2,
  (-1/2, 3/2): 2,
  (1/2, -3/2): 2,
  (1/2, 3/2): 2,
  (3/2, -1/2): 2,
  (3/2, 1/2): 2,
 (-3/2, -1/2): {'sink': 2, (-3/2, 1/2): 1, (-1/2, -1/2): 1},
 (-3/2, 1/2): {'sink': 2, (-3/2, -1/2): 1, (-1/2, 1/2): 1},
 (-1/2, -3/2): {'sink': 2, (-1/2, -1/2): 1, (1/2, -3/2): 1},
 (-1/2, -1/2): { (-3/2, -1/2): 1,
  (-1/2, -3/2): 1,
  (-1/2, 1/2): 1,
  (1/2, -1/2): 1,
 (-1/2, 1/2): {(-3/2, 1/2): 1, (-1/2, -1/2): 1, (-1/2, 3/2): 1, (1/2, 1/2): 1},
 (-1/2, 3/2): {'sink': 2, (-1/2, 1/2): 1, (1/2, 3/2): 1},
 (1/2, -3/2): {'sink': 2, (-1/2, -3/2): 1, (1/2, -1/2): 1},
 (1/2, -1/2): {(-1/2, -1/2): 1, (1/2, -3/2): 1, (1/2, 1/2): 1, (3/2, -1/2): 1},
 (1/2, 1/2): {(-1/2, 1/2): 1, (1/2, -1/2): 1, (1/2, 3/2): 1, (3/2, 1/2): 1},
 (1/2, 3/2): {'sink': 2, (-1/2, 3/2): 1, (1/2, 1/2): 1},
 (3/2, -1/2): {'sink': 2, (1/2, -1/2): 1, (3/2, 1/2): 1},
(3/2, 1/2): {'sink': 2, (1/2, 1/2): 1, (3/2, -1/2): 1}}
sage: Sandpile(aztec_sandpile(2),'sink').group_order()
4542720
```

Note: This is the aztec diamond graph with a sink vertex added. Boundary vertices have edges to the sink so that each vertex has degree 4.

```
sage.sandpiles.sandpile.firing_graph (S, eff)
```

Creates a digraph with divisors as vertices and edges between two divisors D and E if firing a single vertex in D gives E.

INPUT:

S – Sandpile

eff - list of divisors

OUTPUT:

DiGraph

EXAMPLES:

```
sage: S = sandpiles.Cycle(6)
sage: D = SandpileDivisor(S, [1,1,1,1,2,0])
sage: eff = D.effective_div()
sage: firing_graph(S,eff).show3d(edge_size=.005,vertex_size=0.01) # long time
```

```
sage.sandpiles.sandpile.glue_graphs (g, h, glue\_g, glue\_h)
```

Glue two graphs together.

INPUT:

- q, h dictionaries for directed multigraphs
- glue_h, glue_g dictionaries for a vertex

OUTPUT:

dictionary for a directed multigraph

EXAMPLES:

```
sage: from sage.sandpiles.sandpile import glue_graphs
sage: x = \{0: \{\}, 1: \{0: 1\}, 2: \{0: 1, 1: 1\}, 3: \{0: 1, 1: 1, 2: 1\}\}
sage: y = \{0: \{\}, 1: \{0: 2\}, 2: \{1: 2\}, 3: \{0: 1, 2: 1\}\}
sage: glue_x = \{1: 1, 3: 2\}
sage: glue_y = {0: 1, 1: 2, 3: 1}
sage: z = glue_graphs(x,y,glue_x,glue_y); z
{0: {},
 'x0': {0: 1, 'x1': 1, 'x3': 2, 'y1': 2, 'y3': 1},
 'x1': {'x0': 1},
 'x2': {'x0': 1, 'x1': 1},
 'x3': {'x0': 1, 'x1': 1, 'x2': 1},
 'y1': {0: 2},
 'y2': {'y1': 2},
 'y3': {0: 1, 'y2': 1}}
sage: S = Sandpile(z, 0)
sage: S.h_vector()
[1, 6, 17, 31, 41, 41, 31, 17, 6, 1]
sage: S.resolution()
'R^1 <-- R^7 <-- R^21 <-- R^35 <-- R^35 <-- R^21 <-- R^7 <-- R^1'
```

Note: This method makes a dictionary for a graph by combining those for g and h. The sink of g is replaced by a vertex that is connected to the vertices of g as specified by $glue_g$ the vertices of g as specified in $glue_h$. The sink of the glued graph is $glue_g$.

Both glue_g and glue_h are dictionaries with entries of the form v:w where v is the vertex to be connected to and w is the weight of the connecting edge.

```
sage.sandpiles.sandpile.min_cycles(G, v)
```

Minimal length cycles in the digraph G starting at vertex v.

INPUT:

- G DiGraph
- v vertex of G

OUTPUT:

list of lists of vertices

EXAMPLES:

```
sage: from sage.sandpiles.sandpile import min_cycles, sandlib
sage: T = sandlib('gor')
sage: [min_cycles(T, i) for i in T.vertices()]
[[], [[1, 3]], [[2, 3, 1], [2, 3]], [[3, 1], [3, 2]]]
```

```
sage.sandpiles.sandpile.parallel_firing_graph(S, eff)
```

Creates a digraph with divisors as vertices and edges between two divisors D and E if firing all unstable vertices in D gives E.

INPUT:

S – Sandpile

eff - list of divisors

OUTPUT:

DiGraph

EXAMPLES:

sage.sandpiles.sandpile.partition_sandpile(S, p)

Each set of vertices in p is regarded as a single vertex, with and edge between A and B if some element of A is connected by an edge to some element of B in S.

INPUT:

S – Sandpile

p - partition of the vertices of S

OUTPUT:

Sandpile

EXAMPLES:

```
sage: from sage.sandpiles.sandpile import admissible_partitions, partition_
sage: S = sandpiles.Cycle(4)
sage: P = [admissible_partitions(S, i) for i in [2,3,4]]
sage: for p in P:
....: sum([partition_sandpile(S, i).betti(verbose=False)[-1] for i in p])
8
3
sage: S.betti()
               1
          0
_____
   0:
        1
                          3
   1:
               6
                     8
total:
                          3
```

sage.sandpiles.sandpile.random_DAG(num_verts, p=0.5, weight_max=1)

A random directed acyclic graph with num_verts vertices. The method starts with the sink vertex and adds vertices one at a time. Each vertex is connected only to only previously defined vertices, and the probability of each possible connection is given by the argument p. The weight of an edge is a random integer between 1 and weight_max.

INPUT:

- num_verts positive integer
- p (default: 0,5) real number such that 0
- weight_max (default: 1) positive integer

OUTPUT:

a dictionary, encoding the edges of a directed acyclic graph with sink $\boldsymbol{0}$

EXAMPLES:

```
sage: d = DiGraph(random_DAG(5, .5)); d
Digraph on 5 vertices
```

```
sage.sandpiles.sandpile.sandlib(selector=None)
```

Returns the sandpile identified by selector. If no argument is given, a description of the sandpiles in the sandlib is printed.

INPUT:

selector - (optional) identifier or None

OUTPUT

sandpile or description

EXAMPLES:

$sage.sandpiles.sandpile.triangle_sandpile(n)$

A triangular sandpile. Each nonsink vertex has out-degree six. The vertices on the boundary of the triangle are connected to the sink.

INPUT:

n – integer

OUTPUT:

Sandpile

EXAMPLES:

```
sage: from sage.sandpiles.sandpile import triangle_sandpile
sage: T = triangle_sandpile(5)
sage: T.group_order()
135418115000
```

```
sage.sandpiles.sandpile.wilmes_algorithm(M)
```

Computes an integer matrix L with the same integer row span as M and such that L is the reduced Laplacian of a directed multigraph.

INPUT:

M – square integer matrix of full rank

OUTPUT:

integer matrix (L)

EXAMPLES:

REFERENCES:

• [PPW2013]

See also:

- sage.combinat.e_one_star
- sage.combinat.constellation

ARITHMETIC DYNAMICAL SYSTEMS

6.1 Generic dynamical systems on schemes

This is the generic class for dynamical systems and contains the exported constructor functions. The constructor functions can take either polynomials (or rational functions in the affine case) or morphisms from which to construct a dynamical system. If the domain is not specified, it is constructed. However, if you plan on working with points or subvarieties in the domain, it recommended to specify the domain. For products of projective spaces the domain must be specified.

The initialization checks are always performed by the constructor functions. It is possible, but not recommended, to skip these checks by calling the class initialization directly.

AUTHORS:

• Ben Hutz (July 2017): initial version

Bases: sage.schemes.generic.morphism.SchemeMorphism_polynomial

Base class for dynamical systems of schemes.

INPUT:

- polys_or_rat_fncts a list of polynomials or rational functions, all of which should have the same parent
- domain an affine or projective scheme, or product of projective schemes, on which polys defines an
 endomorphism. Subschemes are also ok
- names (default: ('X', 'Y')) tuple of strings to be used as coordinate names for a projective space that is constructed

The following combinations of morphism_or_polys and domain are meaningful:

- morphism_or_polys is a SchemeMorphism; domain is ignored in this case
- morphism_or_polys is a list of homogeneous polynomials that define a rational endomorphism of domain
- morphism_or_polys is a list of homogeneous polynomials and domain is unspecified; domain
 is then taken to be the projective space of appropriate dimension over the base ring of the first element
 of morphism_or_polys
- morphism_or_polys is a single polynomial or rational function; domain is ignored and taken to
 be a 1-dimensional projective space over the base ring of morphism_or_polys with coordinate
 names given by names

```
sage: A.<x> = AffineSpace(QQ,1)
sage: f = DynamicalSystem_affine([x^2+1])
sage: type(f)
<class 'sage.dynamics.arithmetic_dynamics.affine_ds.DynamicalSystem_affine_field'>
```

DynamicalSystem defaults to projective:

as scheme morphism()

Return this dynamical system as SchemeMorphism polynomial.

OUTPUT: SchemeMorphism_polynomial

change_ring(R, check=True)

Return a new dynamical system which is this map coerced to R.

If check is True, then the initialization checks are performed.

INPUT:

• R – ring or morphism

OUTPUT:

A new DynamicalSystem_projective that is this map coerced to R.

EXAMPLES:

specialization (D=None, phi=None, homset=None)

Specialization of this dynamical system.

Given a family of maps defined over a polynomial ring. A specialization is a particular member of that family. The specialization can be specified either by a dictionary or a SpecializationMorphism.

INPUT:

- D − (optional) dictionary
- phi (optional) SpecializationMorphism
- homset (optional) homset of specialized map

OUTPUT: DynamicalSystem

EXAMPLES:

6.2 Dynamical systems on affine schemes

An endomorphism of an affine scheme or subscheme determined by polynomials or rational functions.

The main constructor function is given by <code>DynamicalSystem_affine</code>. The constructor function can take polynomials, rational functions, or morphisms from which to construct a dynamical system. If the domain is not specified, it is constructed. However, if you plan on working with points or subvarieties in the domain, it recommended to specify the domain.

The initialization checks are always performed by the constructor functions. It is possible, but not recommended, to skip these checks by calling the class initialization directly.

AUTHORS:

- · David Kohel, William Stein
- Volker Braun (2011-08-08): Renamed classes, more documentation, misc cleanups.
- Ben Hutz (2017) relocate code and create new class

 ${\bf class} \ \, {\bf sage.dynamics.arithmetic_dynamics.affine_ds.DynamicalSystem_affine} \, (polys_or_rat_fncts, \\ do-$

 ${\it main}) \\ {\it Bases: sage.schemes.affine_morphism.SchemeMorphism_polynomial_affine_space,} \\$

sage.dynamics.arithmetic_dynamics.generic_ds.DynamicalSystem

An endomorphism of affine schemes determined by rational functions.

Warning: You should not create objects of this class directly because no type or consis-

tency checking is performed. The preferred method to construct such dynamical systems is to use DynamicalSystem_affine() function.

INPUT:

- morphism_or_polys a SchemeMorphism, a polynomial, a rational function, or a list or tuple of polynomials or rational functions
- domain optional affine space or subscheme of such; the following combinations of morphism_or_polys and domain are meaningful:
 - morphism_or_polys is a SchemeMorphism; domain is ignored in this case
 - morphism_or_polys is a list of polynomials or rational functions that define a rational endomorphism of domain

- morphism_or_polys is a list of polynomials or rational functions and domain is unspecified;
 domain is then taken to be the affine space of appropriate dimension over the base ring of the first element of morphism_or_polys
- morphism_or_polys is a single polynomial or rational function; domain is ignored and assumed to be the 1-dimensional affine space over the base ring of morphism_or_polys

OUTPUT: DynamicalSystem_affine

EXAMPLES:

Notice that ZZ becomes QQ since the function is rational:

```
sage: A.<x,y> = AffineSpace(QQ, 2)
sage: DynamicalSystem_affine([3/2*x^2, y^2])
Dynamical System of Affine Space of dimension 2 over Rational Field
```

```
Defn: Defined on coordinates by sending (x, y) to (3/2*x^2, y^2)
```

If you pass in quotient ring elements, they are reduced:

```
sage: x = var('x')
sage: DynamicalSystem_affine(x^2+1)
Traceback (most recent call last):
...
TypeError: Symbolic Ring cannot be the base ring
```

dynatomic polynomial(period)

Compute the (affine) dynatomic polynomial of a dynamical system $f: \mathbb{A}^1 \to \mathbb{A}^1$.

The dynatomic polynomial is the analog of the cyclotomic polynomial and its roots are the points of formal period n.

ALGORITHM:

Homogenize to a map $f: \mathbb{P}^1 \to \mathbb{P}^1$ and compute the dynatomic polynomial there. Then, dehomogenize.

INPUT:

• period – a positive integer or a list/tuple [m, n], where m is the preperiod and n is the period

OUTPUT:

If possible, a single variable polynomial in the coordinate ring of the polynomial. Otherwise a fraction field element of the coordinate ring of the polynomial.

```
sage: A.<x,y> = AffineSpace(QQ, 2)
sage: f = DynamicalSystem_affine([x^2+y^2, y^2])
sage: f.dynatomic_polynomial(2)
Traceback (most recent call last):
...
TypeError: does not make sense in dimension >1
```

```
sage: A.<x> = AffineSpace(ZZ, 1)
sage: f = DynamicalSystem_affine([(x^2+1)/x])
sage: f.dynatomic_polynomial(4)
2*x^12 + 18*x^10 + 57*x^8 + 79*x^6 + 48*x^4 + 12*x^2 + 1
```

```
sage: A.<x> = AffineSpace(CC, 1)
sage: f = DynamicalSystem_affine([(x^2+1)/(3*x)])
sage: f.dynatomic_polynomial(3)
13.0000000000000*x^6 + 117.00000000000*x^4 + 78.00000000000*x^2 +
1.00000000000000
```

```
sage: A.<x> = AffineSpace(QQ, 1)
sage: f = DynamicalSystem_affine([x^2-10/9])
sage: f.dynatomic_polynomial([2, 1])
531441*x^4 - 649539*x^2 - 524880
```

```
sage: A.<x> = AffineSpace(CC, 1)
sage: f = DynamicalSystem_affine([x^2+CC.0])
sage: f.dynatomic_polynomial(2)
x^2 + x + 1.000000000000000 + 1.0000000000000*I
```

```
sage: K.<c> = FunctionField(QQ)
sage: A.<x> = AffineSpace(K, 1)
sage: f = DynamicalSystem_affine([x^2 + c])
sage: f.dynatomic_polynomial(4)
x^12 + 6*c*x^10 + x^9 + (15*c^2 + 3*c)*x^8 + 4*c*x^7 + (20*c^3 + 12*c^2 + ...)*x^6
+ (6*c^2 + 2*c)*x^5 + (15*c^4 + 18*c^3 + 3*c^2 + 4*c)*x^4 + (4*c^3 + 4*c^2 + ...)*x^3
+ (6*c^5 + 12*c^4 + 6*c^3 + 5*c^2 + c)*x^2 + (c^4 + 2*c^3 + c^2 + 2*c)*x
+ c^6 + 3*c^5 + 3*c^4 + 3*c^3 + 2*c^2 + 1
```

```
sage: A.<z> = AffineSpace(QQ, 1)
sage: f = DynamicalSystem_affine([z^2+3/z+1/7])
sage: f.dynatomic_polynomial(1).parent()
Multivariate Polynomial Ring in z over Rational Field
```

```
sage: R.<c> = QQ[]
sage: A.<z> = AffineSpace(R,1)
sage: f = DynamicalSystem_affine([z^2 + c])
sage: f.dynatomic_polynomial([1,1])
z^2 + z + c
```

```
sage: A.<x> = AffineSpace(CC,1)
sage: F = DynamicalSystem_affine([1/2*x^2 + CC(sqrt(3))])
sage: F.dynatomic_polynomial([1,1])
(0.12500000000000*x^4 + 0.366025403784439*x^2 + 1.5000000000000)/(0.

→500000000000000*x^2 - x + 1.73205080756888)
```

homogenize(n)

Return the homogenization of this dynamical system.

If its domain is a subscheme, the domain of the homogenized map is the projective embedding of the domain. The domain and codomain can be homogenized at different coordinates: n[0] for the domain and n[1] for the codomain.

INPUT:

• n – a tuple of nonnegative integers. If n is an integer, then the two values of the tuple are assumed to be the same

OUTPUT: DynamicalSystem_projective

EXAMPLES:

```
sage: A.<x,y> = AffineSpace(ZZ, 2)
sage: f = DynamicalSystem_affine([(x^2-2)/x^5, y^2])
sage: f.homogenize(2)
Dynamical System of Projective Space of dimension 2 over Rational Field
   Defn: Defined on coordinates by sending (x0 : x1 : x2) to
        (x0^2*x2^5 - 2*x2^7 : x0^5*x1^2 : x0^5*x2^2)
```

```
sage: P.<x,y,z> = AffineSpace(QQ, 3)
sage: f = DynamicalSystem_affine([x^2 - 2*x*y + z*x, z^2 -y^2, 5*z*y])
sage: f.homogenize(2).dehomogenize(2) == f
True
```

multiplier (P, n, check=True)

Return the multiplier of the point P of period n by this dynamical system.

INPUT:

- P a point on domain of the map
- n a positive integer, the period of P
- check (default: True) boolean, verify that P has period n

OUTPUT:

A square matrix of size self.codomain().dimension_relative() in the base_ring of the map.

EXAMPLES:

```
sage: P.<x,y> = AffineSpace(QQ, 2)
sage: f = DynamicalSystem_affine([x^2, y^2])
sage: f.multiplier(P([1, 1]), 1)
[2 0]
[0 2]
```

```
sage: P.<x,y,z> = AffineSpace(QQ, 3)
sage: f = DynamicalSystem_affine([x, y^2, z^2 - y])
sage: f.multiplier(P([1/2, 1, 0]), 2)
[1 0 0]
[0 4 0]
[0 0 0]
```

```
sage: P.<x> = AffineSpace(CC, 1)
sage: f = DynamicalSystem_affine([x^2 + 1/2])
sage: f.multiplier(P([0.5 + 0.5*I]), 1)
[1.000000000000000 + 1.0000000000000*I]
```

```
sage: R.<t> = PolynomialRing(CC, 1)
sage: P.<x> = AffineSpace(R, 1)
sage: f = DynamicalSystem_affine([x^2 - t^2 + t])
sage: f.multiplier(P([-t + 1]), 1)
[(-2.000000000000000)*t + 2.000000000000]
```

```
sage: P.<x,y> = AffineSpace(QQ, 2)
sage: X = P.subscheme([x^2-y^2])
sage: f = DynamicalSystem_affine([x^2, y^2], domain=X)
sage: f.multiplier(X([1, 1]), 1)
[2 0]
[0 2]
```

${\tt nth_iterate}\,(P,n)$

Return the n-th iterate of the point P by this dynamical system.

INPUT:

- P a point in the map's domain
- n a positive integer

OUTPUT: a point in the map's codomain

```
sage: A.<x,y> = AffineSpace(QQ, 2)
sage: f = DynamicalSystem_affine([(x-2*y^2)/x, 3*x*y])
sage: f.nth_iterate(A(9, 3), 3)
(-104975/13123, -9566667)
```

```
sage: A.<x,y> = AffineSpace(ZZ, 2)
sage: X = A.subscheme([x-y^2])
sage: f = DynamicalSystem_affine([9*y^2, 3*y], domain=X)
sage: f.nth_iterate(X(9, 3), 4)
(59049, 243)
```

nth_iterate_map(n)

Return the n-th iterate of self.

ALGORITHM:

Uses a form of successive squaring to reducing computations.

Todo: This could be improved.

INPUT:

• n – a positive integer

OUTPUT: a dynamical system of affine space

```
sage: A.<x> = AffineSpace(QQ, 1)
sage: f = DynamicalSystem_affine([(3*x^2-2)/(x)])
sage: f.nth_iterate_map(3)
Dynamical System of Affine Space of dimension 1 over Rational Field
```

```
Defn: Defined on coordinates by sending (x) to  ((2187*x^8 - 6174*x^6 + 6300*x^4 - 2744*x^2 + 432)/(81*x^7 - 168*x^5 + 112*x^3 - 24*x))
```

orbit(P, n)

Return the orbit of P by the dynamical system.

Let F be this dynamical system. If n is an integer return $[P, F(P), \ldots, F^n(P)]$. If n is a list or tuple n = [m, k] return $[F^m(P), \ldots, F^k(P)]$.

INPUT:

- P a point in the map's domain
- n a non-negative integer or list or tuple of two non-negative integers

OUTPUT: a list of points in the map's codomain

```
sage: A.<x,y> = AffineSpace(QQ, 2)
sage: f = DynamicalSystem_affine([(x-2*y^2)/x, 3*x*y])
sage: f.orbit(A(9, 3), 3)
[(9, 3), (-1, 81), (13123, -243), (-104975/13123, -9566667)]
```

```
sage: A.<x> = AffineSpace(QQ, 1)
sage: f = DynamicalSystem_affine([(x-2)/x])
sage: f.orbit(A(1/2), [1, 3])
[(-3), (5/3), (-1/5)]
```

```
sage: A.<x,y> = AffineSpace(ZZ, 2)
sage: X = A.subscheme([x-y^2])
sage: f = DynamicalSystem_affine([9*y^2, 3*y], domain=X)
sage: f.orbit(X(9, 3), (0, 4))
[(9, 3), (81, 9), (729, 27), (6561, 81), (59049, 243)]
```

```
sage: R.<t> = PolynomialRing(QQ)
sage: A.<x,y> = AffineSpace(FractionField(R), 2)
sage: f = DynamicalSystem_affine([(x-t*y^2)/x, t*x*y])
sage: f.orbit(A(1, t), 3)
[(1, t), (-t^3 + 1, t^2), ((-t^5 - t^3 + 1)/(-t^3 + 1), -t^6 + t^3),
```

```
((-t^16 + 3*t^13 - 3*t^10 + t^7 + t^5 + t^3 - 1)/(t^5 + t^3 - 1), -t^9 - t^7 + t^4)]
```

main)

Bases: sage.dynamics.arithmetic_dynamics.affine_ds.DynamicalSystem_affine, sage.schemes.affine_affine_morphism.SchemeMorphism_polynomial_affine_space_field

weil_restriction()

Compute the Weil restriction of this morphism over some extension field.

If the field is a finite field, then this computes the Weil restriction to the prime subfield.

A Weil restriction of scalars - denoted $Res_{L/k}$ - is a functor which, for any finite extension of fields L/k and any algebraic variety X over L, produces another corresponding variety $Res_{L/k}(X)$, defined over k. It is useful for reducing questions about varieties over large fields to questions about more complicated varieties over smaller fields. Since it is a functor it also applied to morphisms. In particular, the functor applied to a morphism gives the equivalent morphism from the Weil restriction of the domain to the Weil restriction of the codomain.

OUTPUT:

Scheme morphism on the Weil restrictions of the domain and codomain of the map.

EXAMPLES:

```
sage: K.<v> = QuadraticField(5)
sage: PS.<x,y> = AffineSpace(K, 2)
sage: f = DynamicalSystem_affine([x, y])
sage: F = f.weil_restriction()
sage: P = PS(2, 1)
sage: Q = P.weil_restriction()
sage: f(P).weil_restriction() == F(Q)
True
```

main)

Bases: sage.dynamics.arithmetic_dynamics.affine_ds.DynamicalSystem_affine_field, sage.schemes.affine_morphism.SchemeMorphism_polynomial_affine_space_finite_fiel

cyclegraph()

Return the digraph of all orbits of this morphism mod p.

For subschemes, only points on the subscheme whose image are also on the subscheme are in the digraph.

OUTPUT: a digraph

EXAMPLES:

```
sage: P.<x,y> = AffineSpace(GF(5), 2)
sage: f = DynamicalSystem_affine([x^2-y, x*y+1])
sage: f.cyclegraph()
Looped digraph on 25 vertices
```

```
sage: P.<x> = AffineSpace(GF(3^3, 't'), 1)
sage: f = DynamicalSystem_affine([x^2-1])
sage: f.cyclegraph()
Looped digraph on 27 vertices
```

```
sage: P.<x,y> = AffineSpace(GF(7), 2)
sage: X = P.subscheme(x-y)
sage: f = DynamicalSystem_affine([x^2, y^2], domain=X)
sage: f.cyclegraph()
Looped digraph on 7 vertices
```

orbit structure (P)

Every point is preperiodic over a finite field.

This function returns the pair [m, n] where m is the preperiod and n is the period of the point P by this map.

INPUT:

• P - a point in the map's domain

OUTPUT: a list [m, n] of integers

EXAMPLES:

```
sage: A.<x,y> = AffineSpace(GF(13), 2)
sage: f = DynamicalSystem_affine([x^2 - 1, y^2])
sage: f.orbit_structure(A(2, 3))
[1, 6]
```

```
sage: A.<x,y,z> = AffineSpace(GF(49, 't'), 3)
sage: f = DynamicalSystem_affine([x^2 - z, x - y + z, y^2 - x^2])
sage: f.orbit_structure(A(1, 1, 2))
[7, 6]
```

6.3 Dynamical systems on projective schemes

A dynamical system of projective schemes determined by homogeneous polynomials functions that define what the morphism does on points in the ambient projective space.

The main constructor functions are given by <code>DynamicalSystem</code> and <code>DynamicalSystem_projective</code>. The constructors function can take either polynomials or a morphism from which to construct a dynamical system. If the domain is not specified, it is constructed. However, if you plan on working with points or subvarieties in the domain, it recommended to specify the domain.

The initialization checks are always performed by the constructor functions. It is possible, but not recommended, to skip these checks by calling the class initialization directly.

AUTHORS:

• David Kohel, William Stein

- William Stein (2006-02-11): fixed bug where P(0,0,0) was allowed as a projective point.
- Volker Braun (2011-08-08): Renamed classes, more documentation, misc cleanups.
- Ben Hutz (2013-03) iteration functionality and new directory structure for affine/projective, height functionality
- Brian Stout, Ben Hutz (Nov 2013) added minimal model functionality
- Dillon Rose (2014-01): Speed enhancements
- Ben Hutz (2015-11): iteration of subschemes
- Ben Hutz (2017-7): relocate code and create class

Bases: sage.schemes.projective.projective_morphism.SchemeMorphism_polynomial_projective_sage.dynamics.arithmetic_dynamics.generic_ds.DynamicalSystem

A dynamical system of projective schemes determined by homogeneous polynomials that define what the morphism does on points in the ambient projective space.

Warning: You should not create objects of this class directly because no type or consistency checking is performed. The preferred method to construct such dynamical systems is to use <code>DynamicalSystem_projective()</code> function

INPUT:

- morphism_or_polys a SchemeMorphism, a polynomial, a rational function, or a list or tuple of homogeneous polynomials.
- domain optional projective space or projective subscheme.
- names optional tuple of strings to be used as coordinate names for a projective space that is constructed; defaults to 'X', 'Y'.

The following combinations of morphism_or_polys and domain are meaningful:

- morphism_or_polys is a SchemeMorphism; domain is ignored in this case.
- morphism_or_polys is a list of homogeneous polynomials that define a rational endomorphism of domain.
- morphism_or_polys is a list of homogeneous polynomials and domain is unspecified; domain is then taken to be the projective space of appropriate dimension over the base ring of the first element of morphism or polys.
- morphism_or_polys is a single polynomial or rational function; domain is ignored and taken to
 be a 1-dimensional projective space over the base ring of morphism_or_polys with coordinate
 names given by names.

OUTPUT: DynamicalSystem_projectve.

We can define dynamical systems on P^1 by giving a polynomial or rational function:

Symbolic Ring elements are not allows:

```
sage: x,y = var('x,y')
sage: DynamicalSystem_projective([x^2,y^2])
Traceback (most recent call last):
...
ValueError: [x^2, y^2] must be elements of a polynomial ring
```

When elements of the quotient ring are used, they are reduced:

We can also compute the forward image of subschemes through elimination. In particular, let $X = V(h_1, \ldots, h_t)$ and define the ideal $I = (h_1, \ldots, h_t, y_0 - f_0(\bar{x}), \ldots, y_n - f_n(\bar{x}))$. Then the elimination ideal $I_{n+1} = I \cap K[y_0, \ldots, y_n]$ is a homogeneous ideal and $f(X) = V(I_{n+1})$:

```
sage: P.<x,y,z> = ProjectiveSpace(QQ, 2)
sage: f = DynamicalSystem_projective([(x-2*y)^2, (x-2*z)^2, x^2])
sage: X = P.subscheme(y-z)
sage: f(f(f(X)))
Closed subscheme of Projective Space of dimension 2 over Rational Field
defined by:
    y - z
```

```
sage: P.<x,y,z,w> = ProjectiveSpace(QQ, 3)
sage: f = DynamicalSystem_projective([(x-2*y)^2, (x-2*z)^2, (x-2*w)^2, x^2])
sage: f(P.subscheme([x,y,z]))
Closed subscheme of Projective Space of dimension 3 over Rational Field
defined by:
    w,
    y,
    x
```

automorphism_group(**kwds)

Calculates the subgroup of PGL2 that is the automorphism group of this dynamical system.

The automorphism group is the set of PGL(2) elements that fixes this map under conjugation.

INPUT:

keywords:

- starting_prime (default: 5) the first prime to use for CRT
- algorithm—(optional) can be one of the following:
 - 'CRT' Chinese Remainder Theorem
 - 'fixed_points' fixed points algorithm
- return_functions-(default: False) boolean; True returns elements as linear fractional transformations and False returns elements as PGL2 matrices
- iso_type (default: False) boolean; True returns the isomorphism type of the automorphism group

OUTPUT: a list of elements in the automorphism group

AUTHORS:

- Original algorithm written by Xander Faber, Michelle Manes, Bianca Viray
- Modified by Joao Alberto de Faria, Ben Hutz, Bianca Thompson

REFERENCES:

• [FMV2014]

```
sage: R.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2-y^2, x*y])
```

```
sage: f.automorphism_group(return_functions=True)
[x, -x]
```

```
sage: R.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2-2*x*y-2*y^2, -2*x^2-2*x*y+y^2])
sage: f.automorphism_group(return_functions=True)
[x, 2/(2*x), -x - 1, -2*x/(2*x + 2), (-x - 1)/x, -1/(x + 1)]
```

```
sage: A.<z> = AffineSpace(QQ,1)
sage: f = DynamicalSystem_affine([1/z^3])
sage: F = f.homogenize(1)
sage: F.automorphism_group()
[
[1 0] [0 2] [-1 0] [ 0 -2]
[0 1], [2 0], [ 0 1], [ 2 0]
]
```

```
sage: P.<x,y,z> = ProjectiveSpace(QQ,2)
sage: f = DynamicalSystem_projective([x**2 + x*z, y**2, z**2])
sage: f.automorphism_group() # long time
[
[1 0 0]
[0 1 0]
[0 0 1]
]
```

canonical_height (P, **kwds)

Evaluate the (absolute) canonical height of P with respect to this dynamical system.

Must be over number field or order of a number field. Specify either the number of terms of the series to evaluate or the error bound required.

ALGORITHM:

The sum of the Green's function at the archimedean places and the places of bad reduction.

If function is defined over \mathbf{Q} uses Wells' Algorithm, which allows us to not have to factor the resultant.

INPUT:

• P - a projective point

kwds:

- badprimes (optional) a list of primes of bad reduction
- N (default: 10) positive integer. number of terms of the series to use in the local green functions
- prec (default: 100) positive integer, float point or p-adic precision
- error_bound (optional) a positive real number

OUTPUT: a real number

EXAMPLES:

```
sage: P.<x,y> = ProjectiveSpace(ZZ,1)
sage: f = DynamicalSystem_projective([x^2+y^2, 2*x*y]);
sage: f.canonical_height(P.point([5,4]), error_bound=0.001)
2.1970553519503404898926835324
sage: f.canonical_height(P.point([2,1]), error_bound=0.001)
1.0984430632822307984974382955
```

Notice that preperiodic points may not return exactly 0:

```
sage: P.<x,y,z> = ProjectiveSpace(QQ,2)
sage: X = P.subscheme(x^2-y^2);
sage: f = DynamicalSystem_projective([x^2,y^2, 4*z^2], domain=X);
sage: Q = X([4,4,1])
sage: f.canonical_height(Q, badprimes=[2])
0.0013538030870311431824555314882
```

```
sage: P.<x,y,z> = ProjectiveSpace(QQ,2)
sage: X = P.subscheme(x^2-y^2);
sage: f = DynamicalSystem_projective([x^2,y^2, 30*z^2], domain=X)
sage: Q = X([4, 4, 1])
sage: f.canonical_height(Q, badprimes=[2,3,5], prec=200)
2.7054056208276961889784303469356774912979228770208655455481
```

```
sage: P.<x,y> = ProjectiveSpace(QQ, 1)
sage: f = DynamicalSystem_projective([1000*x^2-29*y^2, 1000*y^2])
sage: Q = P(-1/4, 1)
```

```
sage: f.canonical_height(Q, error_bound=0.01)
3.7996079979254623065837411853
```

conjugate(M)

Conjugate this dynamical system by M, i.e. $M^{-1} \circ f \circ M$.

If possible the new map will be defined over the same space. Otherwise, will try to coerce to the base ring of M.

INPUT:

• M − a square invertible matrix

OUTPUT: a dynamical system

```
sage: P.<x,y> = ProjectiveSpace(ZZ,1)
sage: f = DynamicalSystem_projective([x^2+y^2, y^2])
sage: f.conjugate(matrix([[2,0], [0,1/2]]))
Dynamical System of Projective Space of dimension 1 over Rational Field
```

```
Defn: Defined on coordinates by sending (x : y) to (2*x^2 + 1/8*y^2 : 1/2*y^2)
```

critical height(**kwds)

Compute the critical height of this dynamical system.

The critical height is defined by J. Silverman as the sum of the canonical heights of the critical points. This must be dimension 1 and defined over a number field or number field order.

INPUT:

kwds:

- badprimes (optional) a list of primes of bad reduction
- N (default: 10) positive integer; number of terms of the series to use in the local green functions
- prec (default: 100) positive integer, float point or p-adic precision
- error_bound (optional) a positive real number
- ullet embedding (optional) the embedding of the base field to $\overline{\mathbf{Q}}$

OUTPUT: real number

EXAMPLES:

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^3+7*y^3, 11*y^3])
sage: f.critical_height()
1.1989273321156851418802151128
```

```
sage: K.<w> = QuadraticField(2)
sage: P.<x,y> = ProjectiveSpace(K,1)
sage: f = DynamicalSystem_projective([x^2+w*y^2, y^2])
sage: f.critical_height()
0.16090842452312941163719755472
```

Postcritically finite maps have critical height 0:

critical_point_portrait (check=True, embedding=None)

If this dynamical system is post-critically finite, return its critical point portrait.

This is the directed graph of iterates starting with the critical points. Must be dimension 1. If check is True, then the map is first checked to see if it is postcritically finite.

INPUT:

- check boolean
- ullet embedding embedding of base ring into $\overline{\mathbf{Q}}$

OUTPUT: a digraph

EXAMPLES:

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^5 + 5/4*x*y^4, y^5])
sage: f.critical_point_portrait(check=False)
Looped digraph on 5 vertices
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2 + 2*y^2, y^2])
sage: f.critical_point_portrait()
Traceback (most recent call last):
...
TypeError: map must be post-critically finite
```

critical_points(R=None)

Return the critical points of this dynamical system defined over the ring R or the base ring of this map.

Must be dimension 1.

INPUT:

• R – (optional) a ring

OUTPUT: a list of projective space points defined over R

EXAMPLES:

```
sage: set_verbose(None)
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^3-2*x*y^2 + 2*y^3, y^3])
sage: f.critical_points()
[(1 : 0)]
sage: K.<w> = QuadraticField(6)
sage: f.critical_points(K)
[(-1/3*w : 1), (1/3*w : 1), (1 : 0)]
```

```
sage: set_verbose(None)
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([2*x^2-y^2, x*y])
sage: f.critical_points(QQbar)
[(-0.7071067811865475?*I : 1), (0.7071067811865475?*I : 1)]
```

critical_subscheme()

Return the critical subscheme of this dynamical system.

OUTPUT: projective subscheme

EXAMPLES:

```
sage: set_verbose(None)
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^3-2*x*y^2 + 2*y^3, y^3])
sage: f.critical_subscheme()
Closed subscheme of Projective Space of dimension 1 over Rational Field
defined by:
9*x^2*y^2 - 6*y^4
```

```
sage: set_verbose(None)
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([2*x^2-y^2, x*y])
sage: f.critical_subscheme()
Closed subscheme of Projective Space of dimension 1 over Rational Field
defined by:
4*x^2 + 2*y^2
```

```
sage: P.<x,y,z> = ProjectiveSpace(QQ,2)
sage: f = DynamicalSystem_projective([2*x^2-y^2, x*y, z^2])
sage: f.critical_subscheme()
Closed subscheme of Projective Space of dimension 2 over Rational Field
defined by:
8*x^2*z + 4*y^2*z
```

```
sage: P.<x,y,z,w> = ProjectiveSpace(GF(81),3)
sage: g = DynamicalSystem_projective([x^3+y^3, y^3+z^3, z^3+x^3, w^3])
sage: g.critical_subscheme()
Closed subscheme of Projective Space of dimension 3 over Finite Field in
z4 of size 3^4 defined by:
0
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2,x*y])
sage: f.critical_subscheme()
Traceback (most recent call last):
...
TypeError: the function is not a morphism
```

degree_sequence (iterates=2)

Return sequence of degrees of normalized iterates starting with the degree of this dynamical system.

INPUT: iterates - (default: 2) positive integer

OUTPUT: list of integers

```
sage: P2.<X,Y,Z> = ProjectiveSpace(QQ, 2)
sage: f = DynamicalSystem_projective([Z^2, X*Y, Y^2])
sage: f.degree_sequence(15)
[2, 3, 5, 8, 11, 17, 24, 31, 45, 56, 68, 91, 93, 184, 275]
```

```
sage: F.<t> = PolynomialRing(QQ)
sage: P2.<X,Y,Z> = ProjectiveSpace(F, 2)
sage: f = DynamicalSystem_projective([Y*Z, X*Y, Y^2 + t*X*Z])
```

```
sage: f.degree_sequence(5)
[2, 3, 5, 8, 13]
```

```
sage: P2.<X,Y,Z> = ProjectiveSpace(QQ, 2)
sage: f = DynamicalSystem_projective([X^2, Y^2, Z^2])
sage: f.degree_sequence(10)
[2, 4, 8, 16, 32, 64, 128, 256, 512, 1024]
```

```
sage: P2.<X,Y,Z> = ProjectiveSpace(ZZ, 2)
sage: f = DynamicalSystem_projective([X*Y, Y*Z+Z^2, Z^2])
sage: f.degree_sequence(10)
[2, 3, 4, 5, 6, 7, 8, 9, 10, 11]
```

dehomogenize(n)

Return the standard dehomogenization at the n[0] coordinate for the domain and the n[1] coordinate for the codomain.

Note that the new function is defined over the fraction field of the base ring of this map.

INPUT:

• n – a tuple of nonnegative integers; if n is an integer, then the two values of the tuple are assumed to be the same

OUTPUT:

 $DynamicalSystem_affine$ given by dehomogenizing the source and target of self with respect to the given indices.

EXAMPLES:

$dynamical_degree(N=3, prec=53)$

Return an approximation to the dynamical degree of this dynamical system. The dynamical degree is defined as $\lim_{n\to\infty} \sqrt[n]{\deg(f^n)}$.

INPUT:

- N (default: 3) positive integer, iterate to use for approximation
- prec (default: 53) positive integer, real precision to use when computing root

OUTPUT: real number

```
sage: P.<x,y> = ProjectiveSpace(QQ, 1)
sage: f = DynamicalSystem_projective([x^2 + (x*y), y^2])
sage: f.dynamical_degree()
2.000000000000000
```

```
sage: P2.<X,Y,Z> = ProjectiveSpace(ZZ, 2)
sage: f = DynamicalSystem_projective([X*Y, Y*Z+Z^2, Z^2])
```

```
sage: f.dynamical_degree(N=5, prec=100)
1.4309690811052555010452244131
```

dynatomic_polynomial (period)

For a dynamical system of \mathbb{P}^1 compute the dynatomic polynomial.

The dynatomic polynomial is the analog of the cyclotomic polynomial and its roots are the points of formal period *period*. If possible the division is done in the coordinate ring of this map and a polynomial is returned. In rings where that is not possible, a FractionField element will be returned. In certain cases, when the conversion back to a polynomial fails, a SymbolRing element will be returned.

ALGORITHM:

For a positive integer n, let $[F_n, G_n]$ be the coordinates of the nth iterate of f. Then construct

$$\Phi_n^*(f)(x,y) = \sum_{d|n} (yF_d(x,y) - xG_d(x,y))^{\mu(n/d)},$$

where μ is the Möbius function.

For a pair [m, n], let $f^m = [F_m, G_m]$. Compute

$$\Phi_{m,n}^*(f)(x,y) = \Phi_n^*(f)(F_m, G_m)/\Phi_n^*(f)(F_{m-1}, G_{m-1})$$

REFERENCES:

- [Hutz2015]
- [MoPa1994]

INPUT:

• period – a positive integer or a list/tuple [m, n] where m is the preperiod and n is the period

OUTPUT:

If possible, a two variable polynomial in the coordinate ring of this map. Otherwise a fraction field element of the coordinate ring of this map. Or, a SymbolicRing element.

Todo:

- Do the division when the base ring is p-adic so that the output is a polynomial.
- Convert back to a polynomial when the base ring is a function field (not over \mathbf{Q} or F_p).

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2 + y^2, y^2])
sage: f.dynatomic_polynomial(2)
x^2 + x*y + 2*y^2
```

```
sage: P.<x,y> = ProjectiveSpace(ZZ,1)
sage: f = DynamicalSystem_projective([x^2 + y^2, x*y])
sage: f.dynatomic_polynomial(4)
2*x^12 + 18*x^10*y^2 + 57*x^8*y^4 + 79*x^6*y^6 + 48*x^4*y^8 + 12*x^2*y^10 + y^4
→12
```

```
sage: P.<x,y> = ProjectiveSpace(CC,1)
sage: f = DynamicalSystem_projective([x^2 + y^2, 3*x*y])
sage: f.dynatomic_polynomial(3)
13.000000000000*x^6 + 117.00000000000*x^4*y^2 +
78.000000000000*x^2*y^4 + y^6
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2 - 10/9*y^2, y^2])
sage: f.dynatomic_polynomial([2,1])
x^4*y^2 - 11/9*x^2*y^4 - 80/81*y^6
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2 - 29/16*y^2, y^2])
sage: f.dynatomic_polynomial([2,3])
x^12 - 95/8*x^10*y^2 + 13799/256*x^8*y^4 - 119953/1024*x^6*y^6 +
8198847/65536*x^4*y^8 - 31492431/524288*x^2*y^10 +
172692729/16777216*y^12
```

```
sage: P.<x,y> = ProjectiveSpace(ZZ,1)
sage: f = DynamicalSystem_projective([x^2 - y^2, y^2])
sage: f.dynatomic_polynomial([1,2])
x^2 - x*y
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^3 - y^3, 3*x*y^2])
sage: f.dynatomic_polynomial([0,4]) == f.dynatomic_polynomial(4)
True
```

```
sage: P.<x,y,z> = ProjectiveSpace(QQ,2)
sage: f = DynamicalSystem_projective([x^2 + y^2, x*y, z^2])
sage: f.dynatomic_polynomial(2)
Traceback (most recent call last):
...
TypeError: does not make sense in dimension >1
```

```
sage: P.<x,y> = ProjectiveSpace(Qp(5),1)
sage: f = DynamicalSystem_projective([x^2 + y^2, y^2])
sage: f.dynatomic_polynomial(2)
(x^4*y + (2 + O(5^20))*x^2*y^3 - x*y^4 + (2 + O(5^20))*y^5)/(x^2*y - x*y^2 + y^3)
```

```
sage: L.<t> = PolynomialRing(QQ)
sage: P.<x,y> = ProjectiveSpace(L,1)
sage: f = DynamicalSystem_projective([x^2 + t*y^2, y^2])
sage: f.dynatomic_polynomial(2)
x^2 + x*y + (t + 1)*y^2
```

```
sage: K.<c> = PolynomialRing(ZZ)
sage: P.<x,y> = ProjectiveSpace(K,1)
sage: f = DynamicalSystem_projective([x^2 + c*y^2, y^2])
sage: f.dynatomic_polynomial([1, 2])
x^2 - x*y + (c + 1)*y^2
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2 + y^2, y^2])
sage: f.dynatomic_polynomial(2)
x^2 + x*y + 2*y^2
sage: R.<X> = PolynomialRing(QQ)
sage: K.<c> = NumberField(X^2 + X + 2)
sage: PP = P.change_ring(K)
sage: ff = f.change_ring(K)
sage: ff = f.change_ring(K)
sage: ff = f.change_ring(K)
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2 + y^2, x*y])
sage: f.dynatomic_polynomial([2, 2])
x^4 + 4*x^2*y^2 + y^4
sage: R.<X> = PolynomialRing(QQ)
sage: K.<c> = NumberField(X^4 + 4*X^2 + 1)
sage: PP = P.change_ring(K)
sage: ff = f.change_ring(K)
sage: p = PP((c, 1))
sage: ff.nth_iterate(p, 4) == ff.nth_iterate(p, 2)
True
```

```
sage: P.<x,y> = ProjectiveSpace(CC, 1)
sage: f = DynamicalSystem_projective([x^2-CC.0/5*y^2, y^2])
sage: f.dynatomic_polynomial(2)
x^2 + x*y + (1.000000000000000 - 0.2000000000000*I)*y^2
```

```
sage: L.<t> = PolynomialRing(QuadraticField(2).maximal_order())
sage: P.<x, y> = ProjectiveSpace(L.fraction_field() , 1)
sage: f = DynamicalSystem_projective([x^2 + (t^2 + 1)*y^2 , y^2])
sage: f.dynatomic_polynomial(2)
x^2 + x*y + (t^2 + 2)*y^2
```

```
sage: P.<x,y> = ProjectiveSpace(ZZ, 1)
sage: f = DynamicalSystem_projective([x^2 - 5*y^2, y^2])
sage: f.dynatomic_polynomial([3,0])
0
```

```
sage: P.<x, y> = ProjectiveSpace(QQ, 1)
sage: f = DynamicalSystem_projective([x^3 - y^3*2, y^3])
sage: f.dynatomic_polynomial(1).parent()
Multivariate Polynomial Ring in x, y over Rational Field
```

```
sage: R.<c> = QQ[]
sage: P.<x,y> = ProjectiveSpace(R,1)
sage: f = DynamicalSystem_projective([x^2 + c*y^2, y^2])
sage: f.dynatomic_polynomial([1,2]).parent()
Multivariate Polynomial Ring in x, y over Univariate
Polynomial Ring in c over Rational Field
```

```
sage: R.<c> = QQ[]
sage: P.<x,y> = ProjectiveSpace(ZZ,1)
sage: f = DynamicalSystem_projective([x^2 + y^2, (1)*y^2 + (1)*x*y])
sage: f.dynatomic_polynomial([1,2]).parent()
Multivariate Polynomial Ring in x, y over Integer Ring
```

```
sage: P.<x, y> = ProjectiveSpace(QQ, 1)
sage: f = DynamicalSystem_projective([x^2 + y^2, y^2])
sage: f.dynatomic_polynomial(0)
0
sage: f.dynatomic_polynomial([0,0])
0
sage: f.dynatomic_polynomial(-1)
Traceback (most recent call last):
...
TypeError: period must be a positive integer
```

```
sage: R.<c> = QQ[]
sage: P.<x,y> = ProjectiveSpace(R,1)
sage: f = DynamicalSystem_projective([x^2 + c*y^2,y^2])
sage: f.dynatomic_polynomial([1,2]).parent()
Multivariate Polynomial Ring in x, y over Univariate Polynomial Ring in c over Rational Field
```

Some rings still return SymoblicRing elements:

```
sage: S.<t> = FunctionField(CC)
sage: P.<x,y> = ProjectiveSpace(S,1)
sage: f = DynamicalSystem_projective([t*x^2-1*y^2, t*y^2])
sage: f.dynatomic_polynomial([1, 2]).parent()
Symbolic Ring
```

```
sage: R.<x,y> = PolynomialRing(QQ)
sage: S = R.quo(R.ideal(y^2-x+1))
sage: P.<u,v> = ProjectiveSpace(FractionField(S),1)
sage: f = DynamicalSystem_projective([u^2 + S(x^2)*v^2, v^2])
sage: dyn = f.dynatomic_polynomial([1,1]); dyn
v^3*xbar^2 + u^2*v + u*v^2
sage: dyn.parent()
Symbolic Ring
```

green_function (P, v, **kwds)

Evaluate the local Green's function at the place v for P with N terms of the series or to within a given error bound.

Must be over a number field or order of a number field. Note that this is the absolute local Green's function so is scaled by the degree of the base field.

Use v=0 for the archimedean place over \mathbf{Q} or field embedding. Non-archimedean places are prime ideals for number fields or primes over \mathbf{Q} .

ALGORITHM:

See Exercise 5.29 and Figure 5.6 of [Sil2007].

INPUT:

- P a projective point
- v non-negative integer. a place, use 0 for the archimedean place

kwds:

- N (optional default: 10) positive integer. number of terms of the series to use
- prec (default: 100) positive integer, float point or p-adic precision
- error_bound (optional) a positive real number

OUTPUT: a real number

EXAMPLES:

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2+y^2, x*y]);
sage: Q = P(5, 1)
sage: f.green_function(Q, 0, N=30)
1.6460930159932946233759277576
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2+y^2, x*y]);
sage: Q = P(5, 1)
sage: f.green_function(Q, 0, N=200, prec=200)
1.6460930160038721802875250367738355497198064992657997569827
```

```
sage: K.<w> = QuadraticField(3)
sage: P.<x,y> = ProjectiveSpace(K,1)
sage: f = DynamicalSystem_projective([17*x^2+1/7*y^2, 17*w*x*y])
sage: f.green_function(P.point([w, 2], False), K.places()[1])
1.7236334013785676107373093775
sage: f.green_function(P([2, 1]), K.ideal(7), N=7)
0.48647753726382832627633818586
sage: f.green_function(P([w, 1]), K.ideal(17), error_bound=0.001)
-0.70813041039490996737374178059
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2+y^2, x*y])
sage: f.green_function(P.point([5,2], False), 0, N=30)
1.7315451844777407992085512000
sage: f.green_function(P.point([2,1], False), 0, N=30)
0.86577259223181088325226209926
sage: f.green_function(P.point([1,1], False), 0, N=30)
0.43288629610862338612700146098
```

height_difference_bound(prec=None)

Return an upper bound on the different between the canonical height of a point with respect to this dynamical system and the absolute height of the point.

This map must be a morphism.

ALGORITHM:

Uses a Nullstellensatz argument to compute the constant. For details: see [Hutz2015].

INPUT:

• prec – (default: RealField default) positive integer, float point precision

OUTPUT: a real number

EXAMPLES:

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2+y^2, x*y])
sage: f.height_difference_bound()
1.38629436111989
```

This function does not automatically normalize.

A number field example:

is_PGL_minimal(prime_list=None)

Check if this dynamical system is a minimal model in its conjugacy class.

See [BM2012] and [Mol2015] for a description of the algorithm.

INPUT:

• prime_list - (optional) list of primes to check minimality

OUTPUT: boolean

```
sage: PS.<X,Y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([X^2+3*Y^2, X*Y])
sage: f.is_PGL_minimal()
True
```

```
sage: PS.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([6*x^2+12*x*y+7*y^2, 12*x*y])
sage: f.is_PGL_minimal()
False
```

```
sage: PS.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([6*x^2+12*x*y+7*y^2, y^2])
sage: f.is_PGL_minimal()
Traceback (most recent call last):
...
TypeError: affine minimality is only considered for maps not of the form
f or 1/f for a polynomial f
```

is_postcritically_finite(err=0.01, embedding=None)

Determine if this dynamical system is post-critically finite.

Only for endomorphisms of \mathbb{P}^1 . It checks if each critical point is preperiodic. The optional parameter err is passed into is_preperiodic() as part of the preperiodic check.

INPUT:

- err (default: 0.01) positive real number
- embedding embedding of base ring into $\overline{\mathbf{Q}}$

OUTPUT: boolean

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2 - y^2, y^2])
sage: f.is_postcritically_finite()
True
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^3- y^3, y^3])
sage: f.is_postcritically_finite()
False
```

```
sage: R.<z> = QQ[]
sage: K.<v> = NumberField(z^8 + 3*z^6 + 3*z^4 + z^2 + 1)
sage: PS.<x,y> = ProjectiveSpace(K,1)
sage: f = DynamicalSystem_projective([x^3+v*y^3, y^3])
sage: f.is_postcritically_finite(embedding=K.embeddings(QQbar)[0]) # long time
True
```

minimal_model (return_transformation=False, prime_list=None)

Determine if this dynamical system is minimal.

This dynamical system must be defined over the projective line over the rationals. In particular, determine if this map is affine minimal, which is enough to decide if it is minimal or not. See Proposition 2.10 in [BM2012].

REFERENCES:

- [BM2012]
- [Mol2015]

INPUT:

- return_transformation (default: False) boolean; this signals a return of the PGL_2 transformation to conjugate this map to the calculated minimal model
- prime_list (optional) a list of primes, in case one only wants to determine minimality at those specific primes

OUTPUT:

- a scheme morphism on the projective line which is a minimal model of this map
- a $PGL(2, \mathbf{Q})$ element which conjugates this map to a minimal model

```
sage: PS.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([6*x^2+12*x*y+7*y^2, 12*x*y])
```

```
sage: PS.<x,y> = ProjectiveSpace(ZZ,1)
sage: f = DynamicalSystem_projective([6*x^2+12*x*y+7*y^2, 12*x*y + 42*y^2])
sage: g,M=f.minimal_model(return_transformation=True)
sage: f.conjugate(M) == g
True
```

```
sage: PS.<X,Y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([X+Y, X-3*Y])
sage: f.minimal_model()
Traceback (most recent call last):
...
NotImplementedError: minimality is only for degree 2 or higher
```

```
sage: PS.<X,Y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([X^2-Y^2, X^2+X*Y])
sage: f.minimal_model()
Traceback (most recent call last):
...
TypeError: the function is not a morphism
```

multiplier (P, n, check=True)

Return the multiplier of the point P of period n with respect to this dynamical system.

INPUT:

- P a point on domain of this map
- n a positive integer, the period of P
- check (default: True) boolean; verify that P has period n

OUTPUT:

A square matrix of size self.codomain().dimension_relative() in the base_ring of this dynamical system.

```
sage: P.<x,y,z> = ProjectiveSpace(QQ,2)
sage: f = DynamicalSystem_projective([x^2,y^2, 4*z^2]);
sage: Q = P.point([4,4,1], False);
sage: f.multiplier(Q,1)
[2 0]
[0 2]
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([7*x^2 - 28*y^2, 24*x*y])
sage: f.multiplier(P(2,5), 4)
[231361/20736]
```

```
sage: P.<x,y> = ProjectiveSpace(CC,1)
sage: f = DynamicalSystem_projective([x^3 - 25*x*y^2 + 12*y^3, 12*y^3])
```

```
sage: f.multiplier(P(1,1), 5)
[0.389017489711935]
```

```
sage: P.<x,y> = ProjectiveSpace(RR,1)
sage: f = DynamicalSystem_projective([x^2-2*y^2, y^2])
sage: f.multiplier(P(2,1), 1)
[4.00000000000000]
```

```
sage: P.<x,y> = ProjectiveSpace(Qp(13),1)
sage: f = DynamicalSystem_projective([x^2-29/16*y^2, y^2])
sage: f.multiplier(P(5,4), 3)
[6 + 8*13 + 13^2 + 8*13^3 + 13^4 + 8*13^5 + 13^6 + 8*13^7 + 13^8 +
8*13^9 + 13^10 + 8*13^11 + 13^12 + 8*13^13 + 13^14 + 8*13^15 + 13^16 +
8*13^17 + 13^18 + 8*13^19 + O(13^20)]
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2-y^2, y^2])
sage: f.multiplier(P(0,1), 1)
Traceback (most recent call last):
...
ValueError: (0 : 1) is not periodic of period 1
```

multiplier_spectra (n, formal=False, embedding=None, type='point')

Computes the n multiplier spectra of this dynamical system.

This is the set of multipliers of the periodic points of formal period n included with the appropriate multiplicity. User can also specify to compute the n multiplier spectra instead which includes the multipliers of all periodic points of period n. The map must be defined over projective space of dimension 1 over a number field.

INPUT:

- n a positive integer, the period
- formal (default: False) boolean; True specifies to find the formal n multiplier spectra of this map and False specifies to find the n multiplier spectra
- embedding embedding of the base field into $\overline{\mathbf{Q}}$
- type (default: 'point') string; either 'point' or 'cycle' depending on whether you compute one multiplier per point or one per cycle

OUTPUT: a list of $\overline{\mathbf{Q}}$ elements

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([4608*x^10 - 2910096*x^9*y + 325988068*x^
→8*y^2 + 31825198932*x^7*y^3 - 4139806626613*x^6*y^4\
- 44439736715486*x^5*y^5 + 2317935971590902*x^4*y^6 - 15344764859590852*x^3*y^
→7 + 2561851642765275*x^2*y^8\
+ 113578270285012470*x*y^9 - 150049940203963800*y^10, 4608*y^10])
sage: f.multiplier_spectra(1)
[0, -7198147681176255644585/256, 848446157556848459363/19683, -
→3323781962860268721722583135/35184372088832,
529278480109921/256, -4290991994944936653/2097152, 1061953534167447403/19683, →
-3086380435599991/9,
82911372672808161930567/8192, -119820502365680843999, →3553497751559301575157261317/8192]
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2 - 3/4*y^2, y^2])
sage: f.multiplier_spectra(2, formal=False, type='cycle')
[0, 1, 1, 9]
sage: f.multiplier_spectra(2, formal=False, type='point')
[0, 1, 1, 1, 9]
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2 - 7/4*y^2, y^2])
sage: f.multiplier_spectra(3, formal=True, type='cycle')
[1, 1]
sage: f.multiplier_spectra(3, formal=True, type='point')
[1, 1, 1, 1, 1, 1]
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2 + y^2, x*y])
sage: f.multiplier_spectra(1)
[1, 1, 1]
```

nth_iterate(P, n, **kwds)

Return the n-th iterate of the point P by this dynamical system.

If normalize is True, then the coordinates are automatically normalized.

Todo: Is there a more efficient way to do this?

INPUT:

- P a point in this map's domain
- n a positive integer

kwds:

• normalize - (default: False) boolean

OUTPUT: a point in this map's codomain

```
sage: P.<x,y> = ProjectiveSpace(ZZ,1)
sage: f = DynamicalSystem_projective([x^2+y^2, 2*y^2])
sage: Q = P(1,1)
sage: f.nth_iterate(Q,4)
(32768 : 32768)
```

```
sage: P.<x,y> = ProjectiveSpace(ZZ,1)
sage: f = DynamicalSystem_projective([x^2+y^2, 2*y^2])
sage: Q = P(1,1)
sage: f.nth_iterate(Q, 4, normalize=True)
(1 : 1)
```

```
sage: P.<x,y,z> = ProjectiveSpace(QQ,2)
sage: f = DynamicalSystem_projective([x^2, 2*y^2, z^2-x^2])
sage: Q = P(2,7,1)
sage: f.nth_iterate(Q,2)
(-16/7 : -2744 : 1)
```

```
sage: R.<t> = PolynomialRing(QQ)
sage: P.<x,y,z> = ProjectiveSpace(R,2)
sage: f = DynamicalSystem_projective([x^2+t*y^2, (2-t)*y^2, z^2])
sage: Q = P(2+t,7,t)
sage: f.nth_iterate(Q,2)
(t^4 + 2507*t^3 - 6787*t^2 + 10028*t + 16 : -2401*t^3 + 14406*t^2 -
28812*t + 19208 : t^4)
```

```
sage: P.<x,y,z> = ProjectiveSpace(ZZ,2)
sage: X = P.subscheme(x^2-y^2)
sage: f = DynamicalSystem_projective([x^2, y^2, z^2], domain=X)
sage: f.nth_iterate(X(2,2,3), 3)
(256: 256: 6561)
```

```
sage: K.<c> = FunctionField(QQ)
sage: P.<x,y> = ProjectiveSpace(K,1)
sage: f = DynamicalSystem_projective([x^3 - 2*x*y^2 - c*y^3, x*y^2])
sage: f.nth_iterate(P(c,1), 2)
((c^6 - 9*c^4 + 25*c^2 - c - 21)/(c^2 - 3) : 1)

sage: P.<x,y,z> = ProjectiveSpace(QQ,2)
sage: H = Hom(P,P)
sage: f = H([x^2+3*y^2, 2*y^2,z^2])
sage: P(2, 7, 1).nth_iterate(f, -2)
Traceback (most recent call last):
...
TypeError: must be a forward orbit
```

```
sage: P.<x,y> = ProjectiveSpace(QQ, 1)
sage: f = DynamicalSystem_projective([x^3, x*y^2], domain=P)
sage: f.nth_iterate(P(0, 1), 3, check=False)
(0 : 0)
sage: f.nth_iterate(P(0, 1), 3)
Traceback (most recent call last):
...
ValueError: [0, 0] does not define a valid point since all entries are 0
```

```
sage: P.<x,y> = ProjectiveSpace(ZZ, 1)
sage: f = DynamicalSystem_projective([x^3, x*y^2], domain=P)
sage: f.nth_iterate(P(2,1), 3, normalize=False)
(134217728 : 524288)
sage: f.nth_iterate(P(2,1), 3, normalize=True)
(256 : 1)
```

nth iterate map(n, normalize=False)

Return the n-th iterate of this dynamical system.

ALGORITHM:

Uses a form of successive squaring to reducing computations.

Todo: This could be improved.

INPUT:

- n positive integer
- normalize boolean; remove gcd's during iteration

OUTPUT: a projective dynamical system

```
sage: P.<x,y,z> = ProjectiveSpace(QQ, 2)
sage: f = DynamicalSystem_projective([y^2 * z^3, y^3 * z^2, x^5])
sage: f.nth_iterate_map( 5, normalize=True)
```

```
Dynamical System of Projective Space of dimension 2 over Rational Field Defn: Defined on coordinates by sending (x:y:z) to (y^202*z^443:x^140*y^163*z^342:x^645)
```

orbit(P, N, **kwds)

Return the orbit of the point P by this dynamical system.

Let F be this dynamical system. If N is an integer return $[P, F(P), \ldots, F^N(P)]$. If N is a list or tuple N = [m, k] return $[F^m(P), \ldots, F^k(P)]$. Automatically normalize the points if normalize=True. Perform the checks on point initialization if check=True.

INPUT:

- P a point in this dynamical system's domain
- n a non-negative integer or list or tuple of two non-negative integers

kwds:

- check (default: True) boolean
- normalize (default: False) boolean

OUTPUT: a list of points in this dynamical system's codomain

```
sage: P.<x,y,z> = ProjectiveSpace(ZZ,2)
sage: f = DynamicalSystem_projective([x^2+y^2, y^2-z^2, 2*z^2])
sage: f.orbit(P(1,2,1), 3)
[(1 : 2 : 1), (5 : 3 : 2), (34 : 5 : 8), (1181 : -39 : 128)]
```

```
sage: P.<x,y,z> = ProjectiveSpace(ZZ,2)
sage: f = DynamicalSystem_projective([x^2+y^2, y^2-z^2, 2*z^2])
sage: f.orbit(P(1,2,1), [2,4])
[(34 : 5 : 8), (1181 : -39 : 128), (1396282 : -14863 : 32768)]
```

```
sage: P.<x,y,z> = ProjectiveSpace(ZZ,2)
sage: X = P.subscheme(x^2-y^2)
sage: f = DynamicalSystem_projective([x^2, y^2, x*z], domain=X)
sage: f.orbit(X(2,2,3), 3, normalize=True)
[(2 : 2 : 3), (2 : 2 : 3), (2 : 2 : 3)]
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2+y^2, y^2])
sage: f.orbit(P.point([1,2],False), 4, check=False)
[(1 : 2), (5 : 4), (41 : 16), (1937 : 256), (3817505 : 65536)]
```

```
sage: K.<c> = FunctionField(QQ)
sage: P.<x,y> = ProjectiveSpace(K,1)
sage: f = DynamicalSystem_projective([x^2+c*y^2, y^2])
sage: f.orbit(P(0,1), 3)
[(0:1), (c:1), (c^2 + c:1), (c^4 + 2*c^3 + c^2 + c:1)]
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2+y^2,y^2], domain=P)
sage: f.orbit(P.point([1, 2], False), 4, check=False)
[(1 : 2), (5 : 4), (41 : 16), (1937 : 256), (3817505 : 65536)]
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2, 2*y^2], domain=P)
sage: P(2, 1).orbit(f,[-1, 4])
Traceback (most recent call last):
...
TypeError: orbit bounds must be non-negative
sage: P(2, 1).orbit(f, 0.1)
Traceback (most recent call last):
...
TypeError: Attempt to coerce non-integral RealNumber to Integer
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^3, x*y^2], domain=P)
sage: P(0, 1).orbit(f, 3)
Traceback (most recent call last):
...
ValueError: [0, 0] does not define a valid point since all entries are 0
sage: P(0, 1).orbit(f, 3, check=False)
[(0 : 1), (0 : 0), (0 : 0), (0 : 0)]
```

```
sage: P.<x,y> = ProjectiveSpace(ZZ, 1)
sage: f = DynamicalSystem_projective([x^3, x*y^2], domain=P)
sage: P(2,1).orbit(f, 3, normalize=False)
[(2:1), (8:2), (512:32), (134217728:524288)]
sage: P(2, 1).orbit(f, 3, normalize=True)
[(2:1), (4:1), (16:1), (256:1)]
```

periodic_points (*n*, *minimal=True*, *R=None*, *algorithm='variety'*, *return_scheme=False*)

Computes the periodic points of period n of this dynamical system defined over the ring R or the base ring of the map.

This can be done either by finding the rational points on the variety defining the points of period n, or, for finite fields, finding the cycle of appropriate length in the cyclegraph. For small cardinality fields, the cyclegraph algorithm is effective for any map and length cycle, but is slow when the cyclegraph is large. The variety algorithm is good for small period, degree, and dimension, but is slow as the defining equations of the variety get more complicated.

For rational map, where there are potentially infinitely many peiodic points of a given period, you must use the return_scheme option. Note that this scheme will include the indeterminacy locus.

INPUT:

- n a positive integer
- minimal (default: True) boolean; True specifies to find only the periodic points of minimal period n and False specifies to find all periodic points of period n
- R a commutative ring
- \bullet algorithm (default: 'variety') must be one of the following:
 - 'variety' find the rational points on the appropriate variety
 - 'cyclegraph' find the cycles from the cycle graph
- return_scheme return a subscheme of the ambient space that defines the n th periodic points

OUTPUT:

A list of periodic points of this map or the subscheme defining the periodic points.

```
sage: set_verbose(None)
sage: P.<x,y> = ProjectiveSpace(QQbar,1)
sage: f = DynamicalSystem_projective([x^2-x*y+y^2, x^2-y^2+x*y])
sage: f.periodic_points(1)
[(-0.5000000000000000 - 0.866025403784439?*I : 1),
    (-0.5000000000000000 + 0.866025403784439?*I : 1),
    (1 : 1)]
```

```
sage: P.<x,y,z> = ProjectiveSpace(QuadraticField(5,'t'),2)
sage: f = DynamicalSystem_projective([x^2 - 21/16*z^2, y^2-z^2, z^2])
sage: f.periodic_points(2)
[(-5/4 : -1 : 1), (-5/4 : -1/2*t + 1/2 : 1), (-5/4 : 0 : 1),
    (-5/4 : 1/2*t + 1/2 : 1), (-3/4 : -1 : 1), (-3/4 : 0 : 1),
    (1/4 : -1 : 1), (1/4 : -1/2*t + 1/2 : 1), (1/4 : 0 : 1),
    (1/4 : 1/2*t + 1/2 : 1), (7/4 : -1 : 1), (7/4 : 0 : 1)]
```

```
sage: w = QQ['w'].0
sage: K = NumberField(w^6 - 3*w^5 + 5*w^4 - 5*w^3 + 5*w^2 - 3*w + 1,'s')
sage: P.\langle x,y,z\rangle = ProjectiveSpace(K,2)
sage: f = DynamicalSystem_projective([x^2+z^2, y^2+x^2, z^2+y^2])
sage: f.periodic_points(1)
[(-s^5 + 3*s^4 - 5*s^3 + 4*s^2 - 3*s + 1 : s^5 - 2*s^4 + 3*s^3 - 3*s^2 + 4*s -
\rightarrow 1 : 1),
 (-2*s^5 + 4*s^4 - 5*s^3 + 3*s^2 - 4*s : -2*s^5 + 5*s^4 - 7*s^3 + 6*s^2 - 7*s
\rightarrow+ 3 : 1),
 (-s^5 + 3*s^4 - 4*s^3 + 4*s^2 - 4*s + 2 : -s^5 + 2*s^4 - 2*s^3 + s^2 - s : ...
\hookrightarrow 1),
(s^5 - 2*s^4 + 3*s^3 - 3*s^2 + 3*s - 1 : -s^5 + 3*s^4 - 5*s^3 + 4*s^2 - 4*s_0
\hookrightarrow+ 2 : 1),
(2*s^5 - 6*s^4 + 9*s^3 - 8*s^2 + 7*s - 4 : 2*s^5 - 5*s^4 + 7*s^3 - 5*s^2 + ...
\hookrightarrow6*s - 2 : 1),
(1:1:1),
 (s^5 - 2*s^4 + 2*s^3 + s : s^5 - 3*s^4 + 4*s^3 - 3*s^2 + 2*s - 1 : 1)
```

```
sage: P.<x,y,z> = ProjectiveSpace(QQ,2)
sage: f = DynamicalSystem_projective([x^2 - 21/16*z^2, y^2-2*z^2, z^2])
sage: f.periodic_points(2)
[(-5/4 : -1 : 1), (-5/4 : 2 : 1), (1/4 : -1 : 1), (1/4 : 2 : 1)]
```

```
sage: set_verbose(None)
sage: P.<x,y> = ProjectiveSpace(ZZ, 1)
sage: f = DynamicalSystem_projective([x^2+y^2,y^2])
sage: f.periodic_points(2, R=QQbar, minimal=False)
[(-0.50000000000000000 - 1.322875655532296?*I : 1),
  (-0.5000000000000000 + 1.322875655532296?*I : 1),
  (0.5000000000000000 - 0.866025403784439?*I : 1),
  (0.5000000000000000 + 0.866025403784439?*I : 1),
  (1 : 0)]
```

```
sage: P.<x,y> = ProjectiveSpace(GF(13^2,'t'),1)
sage: f = DynamicalSystem_projective([x^3 + 3*y^3, x^2*y])
sage: f.periodic_points(30, minimal=True, algorithm='cyclegraph')
[(t + 3 : 1), (6*t + 6 : 1), (7*t + 1 : 1), (2*t + 8 : 1),
    (3*t + 4 : 1), (10*t + 12 : 1), (8*t + 10 : 1), (5*t + 11 : 1),
    (7*t + 4 : 1), (4*t + 8 : 1), (9*t + 1 : 1), (2*t + 2 : 1),
    (11*t + 9 : 1), (5*t + 7 : 1), (t + 10 : 1), (12*t + 4 : 1),
    (7*t + 12 : 1), (6*t + 8 : 1), (11*t + 10 : 1), (10*t + 7 : 1),
    (3*t + 9 : 1), (5*t + 5 : 1), (8*t + 3 : 1), (6*t + 11 : 1),
    (9*t + 12 : 1), (4*t + 10 : 1), (11*t + 4 : 1), (2*t + 7 : 1),
    (8*t + 12 : 1), (12*t + 11 : 1)]
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([3*x^2+5*y^2,y^2])
sage: f.periodic_points(2, R=GF(3), minimal=False)
[(2 : 1)]
```

```
sage: P.<x,y,z> = ProjectiveSpace(QQ, 2)
sage: f = DynamicalSystem_projective([x^2, x*y, z^2])
sage: f.periodic_points(1)
Traceback (most recent call last):
...
TypeError: use return_scheme=True
```

```
sage: P.<x,y,z> = ProjectiveSpace(QQ,2)
sage: f = DynamicalSystem_projective([x^2-y^2, x^2-z^2, y^2-z^2])
sage: f.periodic_points(1)
[(-1 : 0 : 1)]
sage: f.periodic_points(1, return_scheme=True)
Closed subscheme of Projective Space of dimension 2 over Rational Field
defined by:
   -x^3 + x^2*y - y^3 + x*z^2,
   -x*y^2 + x^2*z - y^2*z + x*z^2,
   -y^3 + x^2*z + y*z^2 - z^3
sage: f.periodic_points(2, minimal=True, return_scheme=True)
Traceback (most recent call last):
...
NotImplementedError: return_subscheme only implemented for minimal=False
```

possible_periods(**kwds)

Return the set of possible periods for rational periodic points of this dynamical system.

Must be defined over **Z** or **Q**.

ALGORITHM:

Calls self.possible_periods() modulo all primes of good reduction in range prime_bound. Return the intersection of those lists.

INPUT:

kwds:

- prime_bound (default: [1, 20]) a list or tuple of two positive integers or an integer for the upper bound
- bad_primes (optional) a list or tuple of integer primes, the primes of bad reduction
- ncpus (default: all cpus) number of cpus to use in parallel

OUTPUT: a list of positive integers

EXAMPLES:

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2-29/16*y^2, y^2])
sage: f.possible_periods(ncpus=1)
[1, 3]
```

```
sage: PS.<x,y> = ProjectiveSpace(1,QQ)
sage: f = DynamicalSystem_projective([5*x^3 - 53*x*y^2 + 24*y^3, 24*y^3])
sage: f.possible_periods(prime_bound=[1,5])
Traceback (most recent call last):
...
ValueError: no primes of good reduction in that range
sage: f.possible_periods(prime_bound=[1,10])
[1, 4, 12]
sage: f.possible_periods(prime_bound=[1,20])
[1, 4]
```

primes_of_bad_reduction (check=True)

Determine the primes of bad reduction for this dynamical system.

Must be defined over a number field.

If check is True, each prime is verified to be of bad reduction.

ALGORITHM:

p is a prime of bad reduction if and only if the defining polynomials of self have a common zero. Or stated another way, p is a prime of bad reduction if and only if the radical of the ideal defined by the defining polynomials of self is not (x_0, x_1, \ldots, x_N) . This happens if and only if some power of each x_i is not in the ideal defined by the defining polynomials of self. This last condition is what is checked. The lcm of the coefficients of the monomials x_i in a Groebner basis is computed. This may return extra primes.

INPUT:

• check - (default: True) boolean

OUTPUT: a list of primes

EXAMPLES:

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([1/3*x^2+1/2*y^2, y^2])
sage: f.primes_of_bad_reduction()
[2, 3]
```

```
sage: P.<x,y,z,w> = ProjectiveSpace(QQ,3)
sage: f = DynamicalSystem_projective([12*x*z-7*y^2, 31*x^2-y^2, 26*z^2, 3*w^2-y^2)
sage: f.primes_of_bad_reduction()
[2, 3, 7, 13, 31]
```

A number field example:

```
sage: R.<z> = QQ[]
sage: K.<a> = NumberField(z^2 - 2)
sage: P.<x,y> = ProjectiveSpace(K,1)
sage: f = DynamicalSystem_projective([1/3*x^2+1/a*y^2, y^2])
sage: f.primes_of_bad_reduction()
[Fractional ideal (a), Fractional ideal (3)]
```

This is an example where check = False returns extra primes:

reduced_form (prec=300, return_conjugation=True, error_limit=1e-06)

Return reduced form of this dynamical system.

The reduced form is the $SL(2, \mathbf{Z})$ equivalent morphism obtained by applying the binary form reduction algorithm from Stoll and Cremona [SC] to the homogeneous polynomial defining the periodic points (the dynatomic polynomial). The smallest period n with enough periodic points is used.

This should also minimize the sum of the squares of the coefficients, but this is not always the case.

See sage.rings.polynomial.multi_polynomial.reduced_form() for the information on binary form reduction.

Implemented by Rebecca Lauren Miller as part of GSOC 2016.

INPUT:

- prec (default: 300) integer, desired precision
- return_conjuagtion (default: True) boolean; return an element of $SL(2,{f Z})$
- \bullet error_limit (default: 0.000001) a real number, sets the error tolerance

OUTPUT:

- · a projective morphism
- a matrix

```
sage: PS.<x,y> = ProjectiveSpace(QQ, 1)
sage: f = DynamicalSystem_projective([x^3 + x*y^2, y^3])
sage: m = matrix(QQ, 2, 2, [-221, -1, 1, 0])
sage: f = f.conjugate(m)
sage: f.reduced_form(prec=100) #needs 2 periodic
Traceback (most recent call last):
...
ValueError: not enough precision
sage: f.reduced_form()
(
Dynamical System of Projective Space of dimension 1 over Rational Field
Defn: Defined on coordinates by sending (x : y) to
(x^3 + x*y^2 : y^3)
,
[ 0 -1]
[ 1 221]
)
```

```
sage: PS.<x,y> = ProjectiveSpace(ZZ, 1)
sage: f = DynamicalSystem_projective([x^2+ x*y, y^2]) #needs 3 periodic
sage: m = matrix(QQ, 2, 2, [-221, -1, 1, 0])
sage: f = f.conjugate(m)
sage: f.reduced_form(prec=200)
(
Dynamical System of Projective Space of dimension 1 over Integer Ring
Defn: Defined on coordinates by sending (x : y) to
(-x^2 + x*y - y^2 : -y^2)
,
[ 0 -1]
[ 1 220]
)
```

```
(-1 12]
[ 0 -1]
)
```

resultant (normalize=False)

Computes the resultant of the defining polynomials of this dynamical system.

If normalize is True, then first normalize the coordinate functions with normalize_coordinates().

INPUT:

• normalize - (default: False) boolean

OUTPUT: an element of the base ring of this map

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2+y^2, 6*y^2])
sage: f.resultant()
36
```

```
sage: R.<t> = PolynomialRing(GF(17))
sage: P.<x,y> = ProjectiveSpace(R,1)
sage: f = DynamicalSystem_projective([t*x^2+t*y^2, 6*y^2])
sage: f.resultant()
2*t^2
```

```
sage: R.<t> = PolynomialRing(GF(17))
sage: P.<x,y,z> = ProjectiveSpace(R,2)
sage: f = DynamicalSystem_projective([t*x^2+t*y^2, 6*y^2, 2*t*z^2])
sage: f.resultant()
13*t^8
```

```
sage: P.<x,y,z> = ProjectiveSpace(QQ,2)
sage: F = DynamicalSystem_projective([x^2+y^2,6*y^2,10*x*z+z^2+y^2])
sage: F.resultant()
1296
```

```
sage: R.<t>=PolynomialRing(QQ)
sage: s = (t^3+t+1).roots(QQbar)[0][0]
sage: P.<x,y>=ProjectiveSpace(QQbar,1)
sage: f = DynamicalSystem_projective([s*x^3-13*y^3, y^3-15*y^3])
sage: f.resultant()
871.6925062959149?
```

sigma_invariants (n, formal=False, embedding=None, type='point')

Computes the values of the elementary symmetric polynomials of the n multiplier spectra of this dynamical system.

Can specify to instead compute the values corresponding to the elementary symmetric polynomials of the formal n multiplier spectra. The map must be defined over projective space of dimension 1. The base ring should be a number field, number field order, or a finite field or a polynomial ring or function field over a number field, number field order, or finite field.

The parameter type determines if the sigma are computed from the multipliers calculated at one per cycle (with multiplicity) or one per point (with multiplicity). Note that in the cycle case, a map with a cycle which collapses into multiple smaller cycles, this is still considered one cycle. In other words, if a 4-cycle collapses into a 2-cycle with multiplicity 2, there is only one multiplier used for the doubled 2-cycle when computing n=4.

ALGORITHM:

We use the Poisson product of the resultant of two polynomials:

$$res(f,g) = \prod_{f(a)=0} g(a).$$

Letting f be the polynomial defining the periodic or formal periodic points and g the polynomial w - f' for an auxiliary variable w. Note that if f is a rational function, we clear denominators for g.

INPUT:

- n a positive integer, the period
- formal (default: False) boolean; True specifies to find the values of the elementary symmetric polynomials corresponding to the formal n multiplier spectra and False specifies to instead find the values corresponding to the n multiplier spectra, which includes the multipliers of all periodic points of period n
- embedding deprecated in trac ticket #23333
- type (default: 'point') string; either 'point' or 'cycle' depending on whether you compute with one multiplier per point or one per cycle

OUTPUT: a list of elements in the base ring

```
sage: set_verbose(None)
sage: z = QQ['z'].0
sage: K = NumberField(z^4 - 4*z^2 + 1, 'z')
sage: P.<x,y> = ProjectiveSpace(K, 1)
sage: f = DynamicalSystem_projective([x^2 - 5/4*y^2, y^2])
sage: f.sigma_invariants(2, formal=False, type='cycle')
[13, 11, -25, 0]
sage: f.sigma_invariants(2, formal=False, type='point')
[12, -2, -36, 25, 0]
```

check that infinity as part of a longer cycle is handled correctly:

```
sage: P.<x,y> = ProjectiveSpace(QQ, 1)
sage: f = DynamicalSystem_projective([y^2, x^2])
sage: f.sigma_invariants(2, type='cycle')
[12, 48, 64, 0]
sage: f.sigma_invariants(2, type='point')
[12, 48, 64, 0, 0]
sage: f.sigma_invariants(2, type='cycle', formal=True)
[0]
sage: f.sigma_invariants(2, type='point', formal=True)
[0, 0]
```

```
sage: P.<x,y,z> = ProjectiveSpace(QQ, 2)
sage: f = DynamicalSystem_projective([x^2, y^2, z^2])
sage: f.sigma_invariants(1)
Traceback (most recent call last):
...
NotImplementedError: only implemented for dimension 1
```

```
sage: K.<w> = QuadraticField(3)
sage: P.<x,y> = ProjectiveSpace(K, 1)
sage: f = DynamicalSystem_projective([x^2 - w*y^2, (1-w)*x*y])
sage: f.sigma_invariants(2, formal=False, type='cycle')
[6*w + 21, 78*w + 159, 210*w + 367, 90*w + 156]
sage: f.sigma_invariants(2, formal=False, type='point')
[6*w + 24, 96*w + 222, 444*w + 844, 720*w + 1257, 270*w + 468]
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2 + x*y + y^2, y^2 + x*y])
sage: f.sigma_invariants(1)
[3, 3, 1]
```

```
sage: R. <c> = QQ[]
sage: Pc. <x, y> = ProjectiveSpace(R, 1)
sage: f = DynamicalSystem_projective([x^2 + c*y^2, y^2])
sage: f.sigma_invariants(1)
[2, 4*c, 0]
sage: f.sigma_invariants(2, formal=True, type='point')
[8*c + 8, 16*c^2 + 32*c + 16]
sage: f.sigma_invariants(2, formal=True, type='cycle')
[4*c + 4]
```

doubled fixed point:

```
sage: P.\langle x,y \rangle = ProjectiveSpace(QQ, 1)
sage: f = DynamicalSystem_projective([x^2 - 3/4*y^2, y^2])
sage: f.sigma_invariants(2, formal=True)
[2, 1]
```

doubled 2 cycle:

```
sage: P.<x,y> = ProjectiveSpace(QQ, 1)
sage: f = DynamicalSystem_projective([x^2 - 5/4*y^2, y^2])
sage: f.sigma_invariants(4, formal=False, type='cycle')
[170, 5195, 172700, 968615, 1439066, 638125, 0]
```

class sage.dynamics.arithmetic_dynamics.projective_ds.DynamicalSystem_projective_field (polystem_projective_field)

Bases:

```
sage.dynamics.arithmetic_dynamics.projective_ds.
DynamicalSystem_projective,
                              sage.schemes.projective.projective_morphism.
SchemeMorphism_polynomial_projective_space_field
```

main

all rational preimages (points)

Given a set of rational points in the domain of this dynamical system, return all the rational preimages of those points.

In others words, all the rational points which have some iterate in the set points. This function repeatedly calls rational_preimages. If the degree is at least two, by Northocott, this is always a finite set. The map must be defined over number fields and be an endomorphism.

INPUT:

• points – a list of rational points in the domain of this map

OUTPUT: a list of rational points in the domain of this map

EXAMPLES:

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([16*x^2 - 29*y^2, 16*y^2])
sage: sorted(f.all_rational_preimages([P(-1,4)]))
[(-7/4:1), (-5/4:1), (-3/4:1), (-1/4:1), (1/4:1), (3/4:1),
(5/4 : 1), (7/4 : 1)]
```

```
sage: P.<x,y,z> = ProjectiveSpace(QQ,2)
sage: f = DynamicalSystem_projective([76*x^2 - 180*x*y + 45*y^2 + 14*x*z + ...)
45 \times y \times z - 90 \times z^2, 67 \times x^2 - 180 \times x \times y - 157 \times x \times z + 90 \times y \times z, -90 \times z^2])
sage: sorted(f.all_rational_preimages([P(-9,-4,1)]))
[(-9:-4:1), (0:-1:1), (0:0:1), (0:1:1), (0:4:1),
(1:0:1), (1:1:1), (1:2:1), (1:3:1)]
```

A non-periodic example

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2 + y^2, 2*x*y])
sage: sorted(f.all_rational_preimages([P(17,15)]))
[(1/3:1), (3/5:1), (5/3:1), (3:1)]
```

A number field example:

```
sage: z = QQ['z'].0
sage: K. < w > = NumberField(z^3 + (z^2)/4 - (41/16)*z + 23/64);
```

```
sage: P.<x,y> = ProjectiveSpace(K,1)
sage: f = DynamicalSystem_projective([16*x^2 - 29*y^2, 16*y^2])
sage: f.all_rational_preimages([P(16*w^2 - 29,16)])
[(-w^2 + 21/16 : 1),
  (w : 1),
  (w : 1),
  (w + 1/2 : 1),
  (w^2 + w - 33/16 : 1),
  (-w^2 - w + 25/16 : 1),
  (w^2 - 21/16 : 1),
  (-w^2 - w + 33/16 : 1),
  (-w = 1/2 : 1),
  (-w = 1/2 : 1),
  (-w - 1/2 : 1),
  (w^2 + 29/16 : 1),
  (w^2 - 29/16 : 1),
  (w^2 + w - 25/16 : 1)]
```

```
sage: K.<w> = QuadraticField(3)
sage: P.<u,v> = ProjectiveSpace(K,1)
sage: f = DynamicalSystem_projective([u^2+v^2, v^2])
sage: f.all_rational_preimages(P(4))
[(-w : 1), (w : 1)]
```

conjugating_set (other)

Return the set of elements in PGL that conjugates one dynamical system to the other.

Given two nonconstant rational functions of equal degree determine to see if there is an element of PGL that conjugates one rational function to another. It does this by taking the fixed points of one map and mapping them to all unique permutations of the fixed points of the other map. If there are not enough fixed points the function compares the mapping between rational preimages of fixed points and the rational preimages of the preimages of fixed points until there are enough points; such that there are n+2 points with all n+1 subsets linearly independent.

ALGORITHM:

Implementing invariant set algorithm from the paper [FMV2014]. Given that the set of n th preimages of fixed points is invariant under conjugation find all elements of PGL that take one set to another.

INPUT:

• other - a nonconstant rational function of same degree as self

OUTPUT:

Set of conjugating n + 1 by n + 1 matrices.

AUTHORS:

- Original algorithm written by Xander Faber, Michelle Manes, Bianca Viray [FMV2014].
- Implimented by Rebecca Lauren Miller, as part of GSOC 2016.

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2 - 2*y^2, y^2])
sage: m = matrix(QQbar, 2, 2, [-1, 3, 2, 1])
sage: g = f.conjugate(m)
sage: f.conjugating_set(g)
[
[-1 3]
```

```
[ 2 1]
```

```
sage: K.<w> = QuadraticField(-1)
sage: P.<x,y> = ProjectiveSpace(K,1)
sage: f = DynamicalSystem_projective([x^2 + y^2, x*y])
sage: m = matrix(K, 2, 2, [1, 1, 2, 1])
sage: g = f.conjugate(m)
sage: f.conjugating_set(g) # long time
[
[1 1] [-1 -1]
[2 1], [2 1]
]
```

```
sage: K.<i> = QuadraticField(-1)
sage: P.<x,y> = ProjectiveSpace(K,1)
sage: D8 = DynamicalSystem_projective([y^3, x^3])
sage: D8.conjugating_set(D8) # long time
[
[1 0] [0 1] [ 0 -i] [i 0] [ 0 -1] [-1 0] [-i 0] [0 i]
[0 1], [1 0], [ 1 0], [ 0 1], [ 1 0], [ 0 1], [ 1 0]
]
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: D8 = DynamicalSystem_projective([y^2, x^2])
sage: D8.conjugating_set(D8)
Traceback (most recent call last):
...
ValueError: not enough rational preimages
```

```
sage: P.<x,y> = ProjectiveSpace(GF(7),1)
sage: D6 = DynamicalSystem_projective([y^2, x^2])
sage: D6.conjugating_set(D6)
[
[1 0] [0 1] [0 2] [4 0] [2 0] [0 4]
[0 1], [1 0], [1 0], [0 1], [1 0]
]
```

```
sage: P.<x,y,z> = ProjectiveSpace(QQ,2)
sage: f = DynamicalSystem_projective([x^2 + x*z, y^2, z^2])
sage: f.conjugating_set(f) # long time
[
[1 0 0]
[0 1 0]
[0 0 1]
]
```

connected_rational_component (P, n=0)

Computes the connected component of a rational preperiodic point $\ensuremath{\mathbb{P}}$ by this dynamical system.

Will work for non-preperiodic points if n is positive. Otherwise this will not terminate.

INPUT:

- P a rational preperiodic point of this map
- n (default: 0) integer; maximum distance from P to branch out; a value of 0 indicates no bound

OUTPUT:

A list of points connected to P up to the specified distance.

EXAMPLES:

```
sage: R.<x> = PolynomialRing(QQ)
sage: K. < w > = NumberField(x^3+1/4*x^2-41/16*x+23/64)
sage: PS.<x,y> = ProjectiveSpace(1,K)
sage: f = DynamicalSystem_projective([x^2 - 29/16*y^2, y^2])
sage: P = PS([w,1])
sage: f.connected_rational_component(P)
[(w : 1),
 (w^2 - 29/16 : 1),
 (-w^2 - w + 25/16 : 1),
 (w^2 + w - 25/16 : 1),
 (-w : 1),
 (-w^2 + 29/16 : 1),
 (w + 1/2 : 1),
 (-w - 1/2 : 1),
 (-w^2 + 21/16 : 1)
 (w^2 - 21/16 : 1),
 (w^2 + w - 33/16 : 1),
 (-w^2 - w + 33/16 : 1)
```

```
sage: PS.\langle x, y, z \rangle = ProjectiveSpace(2,QQ)
sage: f = DynamicalSystem_projective([x^2 - 21/16*z^2, y^2-2*z^2, z^2])
sage: P = PS([17/16,7/4,1])
sage: f.connected_rational_component(P,3)
[(17/16:7/4:1),
 (-47/256 : 17/16 : 1),
 (-83807/65536 : -223/256 : 1),
 (-17/16 : -7/4 : 1),
 (-17/16 : 7/4 : 1),
 (17/16 : -7/4 : 1),
 (1386468673/4294967296 : -81343/65536 : 1),
 (-47/256 : -17/16 : 1),
 (47/256 : -17/16 : 1),
 (47/256 : 17/16 : 1),
 (-1/2 : -1/2 : 1),
 (-1/2 : 1/2 : 1),
 (1/2 : -1/2 : 1),
 (1/2 : 1/2 : 1)]
```

is_conjugate(other)

Return whether or not two dynamical systems are conjugate.

ALGORITHM:

Implementing invariant set algorithm from the paper [FMV2014]. Given that the set of n th preimages is invariant under conjugation this function finds whether two maps are conjugate.

INPUT:

• other - a nonconstant rational function of same degree as self

OUTPUT: boolean

AUTHORS:

• Original algorithm written by Xander Faber, Michelle Manes, Bianca Viray [FMV2014].

• Implimented by Rebecca Lauren Miller as part of GSOC 2016.

EXAMPLES:

```
sage: K.<w> = CyclotomicField(3)
sage: P.<x,y> = ProjectiveSpace(K,1)
sage: D8 = DynamicalSystem_projective([y^2, x^2])
sage: D8.is_conjugate(D8)
True
```

```
sage: set_verbose(None)
sage: P.<x,y> = ProjectiveSpace(QQbar,1)
sage: f = DynamicalSystem_projective([x^2 + x*y,y^2])
sage: m = matrix(QQbar, 2, 2, [1, 1, 2, 1])
sage: g = f.conjugate(m)
sage: f.is_conjugate(g) # long time
True
```

```
sage: P.<x,y> = ProjectiveSpace(GF(5),1)
sage: f = DynamicalSystem_projective([x^3 + x*y^2,y^3])
sage: m = matrix(GF(5), 2, 2, [1, 3, 2, 9])
sage: g = f.conjugate(m)
sage: f.is_conjugate(g)
True
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2 + x*y,y^2])
sage: g = DynamicalSystem_projective([x^3 + x^2*y, y^3])
sage: f.is_conjugate(g)
False
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2 + x*y, y^2])
sage: g = DynamicalSystem_projective([x^2 - 2*y^2, y^2])
sage: f.is_conjugate(g)
False
```

is polynomial()

Check to see if the dynamical system has a totally ramified fixed point.

The function must be defined over an absolute number field or a finite field.

OUTPUT: boolean

```
sage: R.<x> = QQ[]
sage: K.<w> = QuadraticField(7)
sage: P.<x,y> = ProjectiveSpace(K, 1)
sage: f = DynamicalSystem_projective([x**2 + 2*x*y - 5*y**2, 2*x*y])
sage: f.is_polynomial()
False
```

```
sage: R.<x> = QQ[]
sage: K.<w> = QuadraticField(7)
sage: P.<x,y> = ProjectiveSpace(K, 1)
sage: f = DynamicalSystem_projective([x**2 - 7*x*y, 2*y**2])
sage: m = matrix(K, 2, 2, [w, 1, 0, 1])
```

```
sage: f = f.conjugate(m)
sage: f.is_polynomial()
True
```

```
sage: K.<w> = QuadraticField(4/27)
sage: P.<x,y> = ProjectiveSpace(K,1)
sage: f = DynamicalSystem_projective([x**3 + w*y^3,x*y**2])
sage: f.is_polynomial()
False
```

```
sage: K = GF(3**2, prefix='w')
sage: P.<x,y> = ProjectiveSpace(K,1)
sage: f = DynamicalSystem_projective([x**2 + K.gen()*y**2, x*y])
sage: f.is_polynomial()
False
```

```
sage: PS.<x,y> = ProjectiveSpace(QQ, 1)
sage: f = DynamicalSystem_projective([6*x^2+12*x*y+7*y^2, 12*x*y + 42*y^2])
sage: f.is_polynomial()
False
```

lift_to_rational_periodic (points_modp, B=None)

Given a list of points in projective space over \mathbf{F}_p , determine if they lift to \mathbf{Q} -rational periodic points.

The map must be an endomorphism of projective space defined over Q.

ALGORITHM:

Use Hensel lifting to find a p-adic approximation for that rational point. The accuracy needed is determined by the height bound B. Then apply the LLL algorithm to determine if the lift corresponds to a rational point.

If the point is a point of high multiplicity (multiplier 1), the procedure can be very slow.

INPUT:

- points_modp a list or tuple of pairs containing a point in projective space over \mathbf{F}_p and the possible period
- B (optional) a positive integer; the height bound for a rational preperiodic point

OUTPUT: a list of projective points

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([x^2 - y^2, y^2])
sage: f.lift_to_rational_periodic([[P(0,1).change_ring(GF(7)), 4]])
[[(0:1), 2]]
```

```
There may be multiple points in the lift. sage: P.\langle x,y\rangle = ProjectiveSpace(QQ,1) sage: f = DynamicalSystem\_projective([-5*x^2 + 4*y^2, 4*x*y]) sage: f.lift\_to\_rational\_periodic([[P(1,0).change\_ring(GF(3)), 1]]) # long\_<math>\rightarrow time [[(1 : 0), 1], [(2/3 : 1), 1], [(-2/3 : 1), 1]]
```

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([16*x^2 - 29*y^2, 16*y^2])
```

```
sage: f.lift_to_rational_periodic([[P(3,1).change_ring(GF(13)), 3]])
[[(-1/4 : 1), 3]]
```

```
sage: P.<x,y,z> = ProjectiveSpace(QQ, 2)
sage: f = DynamicalSystem_projective([76*x^2 - 180*x*y + 45*y^2 + 14*x*z + 45*y*z - 90*z^2, 67*x^2 - 180*x*y - 157*x*z + 90*y*z, -90*z^2])
sage: f.lift_to_rational_periodic([[P(14,19,1).change_ring(GF(23)), 9]]) #_\delta long time
[[(-9: -4: 1), 9]]
```

normal form(return conjugation=False)

Return a normal form in the moduli space of dynamical systems.

Currently implemented only for polynomials. The totally ramified fixed point is moved to infinity and the map is conjugated to the form $x^n + a_{n-2}x^{n-2} + \cdots + a_0$. Note that for finite fields we can only remove the (n-1)-st term when the characteristic does not divide n.

INPUT:

• return_conjugation - (default: False) boolean; if True, then return the conjugation element of PGL along with the embedding into the new field

OUTPUT:

- SchemeMorphism_polynomial
- (optional) an element of PGL as a matrix
- (optional) the field embedding

```
sage: P.<x,y> = ProjectiveSpace(QQ, 1)
sage: f = DynamicalSystem_projective([x^2 + 2*x*y - 5*x^2, 2*x*y])
sage: f.normal_form()
Traceback (most recent call last):
...
NotImplementedError: map is not a polynomial
```

rational_periodic_points(**kwds)

Determine the set of rational periodic points for this dynamical system.

The map must be defined over \mathbf{Q} and be an endomorphism of projective space. If the map is a polynomial endomorphism of \mathbb{P}^1 , i.e. has a totally ramified fixed point, then the base ring can be an absolute number field. This is done by passing to the Weil restriction.

The default parameter values are typically good choices for \mathbb{P}^1 . If you are having trouble getting a particular map to finish, try first computing the possible periods, then try various different lifting_prime values.

ALGORITHM:

Modulo each prime of good reduction p determine the set of periodic points modulo p. For each cycle modulo p compute the set of possible periods (mrp^e) . Take the intersection of the list of possible periods modulo several primes of good reduction to get a possible list of minimal periods of rational periodic points. Take each point modulo p associated to each of these possible periods and try to lift it to a rational point with a combination of p-adic approximation and the LLL basis reduction algorithm.

See [Hutz2015].

INPUT:

kwds:

- prime_bound (default: [1, 20]) a pair (list or tuple) of positive integers that represent the limits of primes to use in the reduction step or an integer that represents the upper bound
- lifting_prime (default: 23) a prime integer; argument that specifies modulo which prime to try and perform the lifting
- periods (optional) a list of positive integers that is the list of possible periods
- bad_primes (optional) a list or tuple of integer primes; the primes of bad reduction
- ncpus (default: all cpus) number of cpus to use in parallel

OUTPUT: a list of rational points in projective space

```
sage: P.<x,y> = ProjectiveSpace(QQ,1)
sage: f = DynamicalSystem_projective([-5*x^2 + 4*y^2, 4*x*y])
sage: sorted(f.rational_periodic_points()) # long time
[(-2:1), (-2/3:1), (2/3:1), (1:0), (2:1)]
```

```
sage: R.<x> = QQ[]
sage: K.<w> = NumberField(x^2-x+1)
sage: P.<u,v> = ProjectiveSpace(K,1)
sage: f = DynamicalSystem_projective([u^2 + v^2,v^2])
sage: f.rational_periodic_points()
[(w: 1), (1: 0), (-w + 1: 1)]
```

```
sage: R.<x> = QQ[]
sage: K.<w> = NumberField(x^2-x+1)
sage: P.<u,v> = ProjectiveSpace(K,1)
sage: f = DynamicalSystem_projective([u^2+v^2,u*v])
sage: f.rational_periodic_points()
Traceback (most recent call last):
...
NotImplementedError: rational periodic points for number fields only_
→implemented for polynomials
```

rational preperiodic graph (**kwds)

Determine the directed graph of the rational preperiodic points for this dynamical system.

The map must be defined over \mathbf{Q} and be an endomorphism of projective space. If this map is a polynomial endomorphism of \mathbb{P}^1 , i.e. has a totally ramified fixed point, then the base ring can be an absolute number field. This is done by passing to the Weil restriction.

ALGORITHM:

- Determines the list of possible periods.
- Determines the rational periodic points from the possible periods.
- Determines the rational preperiodic points from the rational periodic points by determining rational preimages.

INPUT:

kwds:

- prime_bound (default: [1, 20]) a pair (list or tuple) of positive integers that represent the limits of primes to use in the reduction step or an integer that represents the upper bound
- lifting_prime (default: 23) a prime integer; specifies modulo which prime to try and perform the lifting
- periods (optional) a list of positive integers that is the list of possible periods
- bad_primes (optional) a list or tuple of integer primes; the primes of bad reduction
- ncpus (default: all cpus) number of cpus to use in parallel

OUTPUT:

A digraph representing the orbits of the rational preperiodic points in projective space.

```
sage: PS.<x,y> = ProjectiveSpace(1,QQ)
sage: f = DynamicalSystem_projective([7*x^2 - 28*y^2, 24*x*y])
sage: f.rational_preperiodic_graph()
Looped digraph on 12 vertices
```

```
sage: PS.<x,y> = ProjectiveSpace(1,QQ)
sage: f = DynamicalSystem_projective([-3/2*x^3 +19/6*x*y^2, y^3])
sage: f.rational_preperiodic_graph(prime_bound=[1,8])
Looped digraph on 12 vertices
```

```
sage: K.<w> = QuadraticField(-3)
sage: P.<x,y> = ProjectiveSpace(K,1)
sage: f = DynamicalSystem_projective([x^2+y^2, y^2])
sage: f.rational_preperiodic_graph() # long time
Looped digraph on 5 vertices
```

rational_preperiodic_points(**kwds)

Determine the set of rational preperiodic points for this dynamical system.

The map must be defined over \mathbf{Q} and be an endomorphism of projective space. If the map is a polynomial endomorphism of \mathbb{P}^1 , i.e. has a totally ramified fixed point, then the base ring can be an absolute number field. This is done by passing to the Weil restriction.

The default parameter values are typically good choices for \mathbb{P}^1 . If you are having trouble getting a particular map to finish, try first computing the possible periods, then try various different values for lifting_prime.

ALGORITHM:

- Determines the list of possible periods.
- Determines the rational periodic points from the possible periods.
- Determines the rational preperiodic points from the rational periodic points by determining rational preimages.

INPUT:

kwds:

- prime_bound (default: [1, 20]) a pair (list or tuple) of positive integers that represent the limits of primes to use in the reduction step or an integer that represents the upper bound
- lifting_prime (default: 23) a prime integer; specifies modulo which prime to try and perform the lifting
- periods (optional) a list of positive integers that is the list of possible periods
- bad_primes (optional) a list or tuple of integer primes; the primes of bad reduction
- ncpus (default: all cpus) number of cpus to use in parallel

OUTPUT: a list of rational points in projective space

```
sage: PS.<x,y> = ProjectiveSpace(1,QQ)
sage: f = DynamicalSystem_projective([x^2 -y^2, 3*x*y])
sage: sorted(f.rational_preperiodic_points())
[(-2 : 1), (-1 : 1), (-1/2 : 1), (0 : 1), (1/2 : 1), (1 : 0), (1 : 1),
(2 : 1)]
```

```
sage: PS.<x,y> = ProjectiveSpace(1,QQ)
sage: f = DynamicalSystem_projective([5*x^3 - 53*x*y^2 + 24*y^3, 24*y^3])
sage: sorted(f.rational_preperiodic_points(prime_bound=10))
[(-1 : 1), (0 : 1), (1 : 0), (1 : 1), (3 : 1)]
```

```
sage: K.<w> = QuadraticField(33)
sage: PS.<x,y> = ProjectiveSpace(K,1)
sage: f = DynamicalSystem_projective([x^2-71/48*y^2, y^2])
sage: sorted(f.rational_preperiodic_points()) # long time
[(-1/12*w - 1 : 1),
 (-1/6*w - 1/4 : 1),
 (-1/12*w - 1/2 : 1),
 (-1/6*w + 1/4 : 1),
 (1/12*w - 1 : 1),
 (1/12*w - 1/2 : 1),
 (-1/12*w + 1/2 : 1),
 (-1/12*w + 1 : 1),
 (1/6*w - 1/4 : 1),
 (1/12*w + 1/2 : 1),
 (1:0),
 (1/6*w + 1/4 : 1),
 (1/12*w + 1 : 1)
```

class sage.dynamics.arithmetic_dynamics.projective_ds.DynamicalSystem_projective_finite_fic

```
Bases: sage.dynamics.arithmetic_dynamics.projective_ds.

DynamicalSystem_projective_field, sage.schemes.projective.

projective_morphism.SchemeMorphism_polynomial_projective_space_finite_field
```

automorphism_group (absolute=False, iso_type=False, return_functions=False)

Return the subgroup of PGL2 that is the automorphism group of this dynamical system.

Only for dimension 1. The automorphism group is the set of PGL2 elements that fixed the map under conjugation. See [FMV2014] for the algorithm.

INPUT:

• absolute—(default: False) boolean; if True, then return the absolute automorphism group and a field of definition

- iso_type (default: False) boolean; if True, then return the isomorphism type of the automorphism group
- return_functions (default: False) boolean; True returns elements as linear fractional transformations and False returns elements as PGL2 matrices

OUTPUT: a list of elements of the automorphism group

AUTHORS:

- Original algorithm written by Xander Faber, Michelle Manes, Bianca Viray
- Modified by Joao Alberto de Faria, Ben Hutz, Bianca Thompson

EXAMPLES:

```
sage: R.<x,y> = ProjectiveSpace(GF(7^3,'t'),1)
sage: f = DynamicalSystem_projective([x^2-y^2, x*y])
sage: f.automorphism_group()
[
[1 0] [6 0]
[0 1], [0 1]
]
```

```
sage: R.<x,y> = ProjectiveSpace(GF(3^2,'t'),1)
sage: f = DynamicalSystem_projective([x^3,y^3])
sage: f.automorphism_group(return_functions=True, iso_type=True) # long time
([x, x/(x + 1), x/(2*x + 1), 2/(x + 2), (2*x + 1)/(2*x), (2*x + 2)/x,
1/(2*x + 2), x + 1, x + 2, x/(x + 2), 2*x/(x + 1), 2*x, 1/x, 2*x + 1,
2*x + 2, ((t + 2)*x + t + 2)/((2*t + 1)*x + t + 2), (t*x + 2*t)/(t*x + t), 2/x, (x + 1)/(x + 2), (2*t*x + t)/(t*x), (2*t + 1)/((2*t + 1)*x + 2*t + 1), ((2*t + 1)*x + 2*t + 1)/((2*t + 1)*x), t/(t*x + 2*t), (2*x + 1)/(x + 1)], 'PGL(2,3)')
```

```
sage: R.<x,y> = ProjectiveSpace(GF(2^5,'t'),1)
sage: f = DynamicalSystem_projective([x^5,y^5])
sage: f.automorphism_group(return_functions=True, iso_type=True)
([x, 1/x], 'Cyclic of order 2')
```

```
sage: R.<x,y> = ProjectiveSpace(GF(3^4,'t'),1)
sage: f = DynamicalSystem_projective([x^2+25*x*y+y^2, x*y+3*y^2])
sage: f.automorphism_group(absolute=True)
[Univariate Polynomial Ring in w over Finite Field in b of size 3^4,
    [
    [1 0]
    [0 1]
    ]]
```

cyclegraph()

Return the digraph of all orbits of this dyanmical system.

Over a finite field this is a finite graph. For subscheme domains, only points on the subscheme whose image are also on the subscheme are in the digraph.

OUTPUT: a digraph

```
sage: P.<x,y> = ProjectiveSpace(GF(13),1)
sage: f = DynamicalSystem_projective([x^2-y^2, y^2])
```

```
sage: f.cyclegraph()
Looped digraph on 14 vertices
```

```
sage: P.<x,y,z> = ProjectiveSpace(GF(3^2,'t'),2)
sage: f = DynamicalSystem_projective([x^2+y^2, y^2, z^2+y*z])
sage: f.cyclegraph()
Looped digraph on 91 vertices
```

```
sage: P.<x,y,z> = ProjectiveSpace(GF(7),2)
sage: X = P.subscheme(x^2-y^2)
sage: f = DynamicalSystem_projective([x^2, y^2, z^2], domain=X)
sage: f.cyclegraph()
Looped digraph on 15 vertices
```

```
sage: P.<x,y,z> = ProjectiveSpace(GF(3),2)
sage: f = DynamicalSystem_projective([x*z-y^2, x^2-y^2, y^2-z^2])
sage: f.cyclegraph()
Looped digraph on 13 vertices
```

```
sage: P.<x,y,z> = ProjectiveSpace(GF(3),2)
sage: X = P.subscheme([x-y])
sage: f = DynamicalSystem_projective([x^2-y^2, x^2-y^2, y^2-z^2], domain=X)
sage: f.cyclegraph()
Looped digraph on 4 vertices
```

orbit structure (P)

Return the pair [m, n], where m is the preperiod and n is the period of the point P by this dynamical system.

Every point is preperiodic over a finite field so every point will be preperiodic.

INPUT:

• P - a point in the domain of this map

OUTPUT: a list [m, n] of integers

```
sage: P.<x,y,z> = ProjectiveSpace(GF(5),2)
sage: f = DynamicalSystem_projective([x^2 + y^2,y^2, z^2 + y*z], domain=P)
sage: f.orbit_structure(P(2,1,2))
[0, 6]
```

```
sage: P.<x,y,z> = ProjectiveSpace(GF(7),2)
sage: X = P.subscheme(x^2-y^2)
sage: f = DynamicalSystem_projective([x^2, y^2, z^2], domain=X)
sage: f.orbit_structure(X(1,1,2))
[0, 2]
```

```
sage: P.<x,y> = ProjectiveSpace(GF(13),1)
sage: f = DynamicalSystem_projective([x^2 - y^2, y^2], domain=P)
sage: f.orbit_structure(P(3,4))
[2, 3]
```

```
sage: R.<t> = GF(13^3)
sage: P.<x,y> = ProjectiveSpace(R,1)
```

```
sage: f = DynamicalSystem_projective([x^2 - y^2, y^2], domain=P)
sage: f.orbit_structure(P(t, 4))
[11, 6]
```

possible_periods (return_points=False)

Return the list of possible minimal periods of a periodic point over \mathbf{Q} and (optionally) a point in each cycle.

ALGORITHM:

See [Hutz2009].

INPUT:

• return_points - (default: False) boolean; if True, then return the points as well as the possible periods

OUTPUT:

A list of positive integers, or a list of pairs of projective points and periods if return_points is True.

EXAMPLES:

```
sage: P.<x,y> = ProjectiveSpace(GF(23),1)
sage: f = DynamicalSystem_projective([x^2-2*y^2, y^2])
sage: f.possible_periods()
[1, 5, 11, 22, 110]
```

```
sage: P.<x,y> = ProjectiveSpace(GF(13),1)
sage: f = DynamicalSystem_projective([x^2-y^2, y^2])
sage: sorted(f.possible_periods(True))
[[(0 : 1), 2], [(1 : 0), 1], [(3 : 1), 3], [(3 : 1), 36]]
```

Todo:

- do not return duplicate points
- improve hash to reduce memory of point-table

6.4 Dynamical systems for products of projective spaces

This class builds on the prouct projective space class. The main constructor functions are given by DynamicalSystem and DynamicalSystem_projective. The constructors function can take either polynomials or a morphism from which to construct a dynamical system.

The must be specified.

class sage.dynamics.arithmetic_dynamics.product_projective_ds.DynamicalSystem_product_projective

Bases: sage.dynamics.arithmetic_dynamics.generic_ds.DynamicalSystem, sage.schemes.product_projective.morphism.ProductProjectiveSpaces_morphism_ring

The class of dynamical systems on products of projective spaces.

Warning: You should not create objects of this class directly because no type or consistency checking is performed. The preferred method to construct such dynamical systems is to use DynamicalSystem_projective() function.

INPUT:

- polys a list of n_1 + \cdots + n_r multi-homogeneous polynomials, all of which should have the same parent
- domain a projective scheme embedded in P^{n_1-1} \times \cdots \times P^{n_r-1}

EXAMPLES:

nth_iterate(P, n, normalize=False)

Return the n-th iterate of P by this dynamical system.

If normalize is True, then the coordinates are automatically normalized.

Todo: Is there a more efficient way to do this?

INPUT:

- P a point in self.domain()
- n a positive integer
- normalize (default: False) boolean

OUTPUT: A point in self.codomain()

```
sage: Z.<a,b,x,y> = ProductProjectiveSpaces([1, 1], ZZ)
sage: f = DynamicalSystem_projective([a*b, b^2, x^3 - y^3, y^2*x], domain=Z)
sage: P = Z([2, 6, 2, 4])
sage: f.nth_iterate(P, 2, normalize = True)
(1 : 3 , 407 : 112)
```

$nth_iterate_map(n)$

Return the nth iterate of this dynamical system.

ALGORITHM:

Uses a form of successive squaring to reduce computations.

Todo: This could be improved.

INPUT:

• n − a positive integer

OUTPUT: A dynamical system of products of projective spaces

EXAMPLES:

orbit(P, N, **kwds)

Return the orbit of P by this dynamical system.

Let F be this dynamical system. If N is an integer return $[P, F(P), \dots, F^N(P)]$.

If N is a list or tuple N = [m, k] return $[F^m(P), \dots, F^k(P)]$. Automatically normalize the points if normalize == True. Perform the checks on point initialize if check==True.

INPUT:

- P a point in self.domain()
- N a non-negative integer or list or tuple of two non-negative integers

kwds:

- check (default: True) boolean
- normalize (default: False) boolean

OUTPUT: a list of points in self.codomain()

6.5 Wehler K3 Surfaces

AUTHORS:

- Ben Hutz (11-2012)
- Joao Alberto de Faria (10-2013)

Todo: Hasse-Weil Zeta Function

Picard Number

Number Fields

REFERENCES: [FH2015], [CS1996], [Weh1998], [Hutz2007]

```
sage.dynamics.arithmetic_dynamics.wehlerK3.WehlerK3Surface(polys)
```

Defines a K3 Surface over $\mathbb{P}^2 \times \mathbb{P}^2$ defined as the intersection of a bilinear and biquadratic form. [Weh1998]

INPUT: Bilinear and biquadratic polynomials as a tuple or list

OUTPUT: WehlerK3Surface_ring

EXAMPLES:

```
sage: PP.<x0,x1, x2, y0, y1, y2> = ProductProjectiveSpaces([2, 2],QQ)
sage: L = x0*y0 + x1*y1 - x2*y2
sage: Q = x0*x1*y1^2 + x2^2*y0*y2
sage: WehlerK3Surface([L, Q])
Closed subscheme of Product of projective spaces P^2 x P^2 over Rational
Field defined by:
  x0*y0 + x1*y1 - x2*y2,
  x0*x1*y1^2 + x2^2*y0*y2
```

```
class sage.dynamics.arithmetic_dynamics.wehlerK3.WehlerK3Surface_finite_field(polys)
    Bases: sage.dynamics.arithmetic dynamics.wehlerK3.WehlerK3Surface field
```

cardinality()

Counts the total number of points on the K3 surface.

ALGORITHM:

Enumerate points over \mathbb{P}^2 , and then count the points on the fiber of each of those points.

OUTPUT: Integer - total number of points on the surface

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], GF(7))
sage: Z = x0^2*y0^2 + 3*x0*x1*y0^2 + x1^2*y0^2 + 4*x0^2*y0*y1 + \
3*x0*x1*y0*y1 -2*x2^2*y0*y1 - x0^2*y1^2 + 2*x1^2*y1^2 - x0*x2*y1^2 \
- 4*x1*x2*y1^2 + 5*x0*x2*y0*y2 -4*x1*x2*y0*y2 + 7*x0^2*y1*y2 + 4*x1^2*y1*y2 \
+ x0*x1*y2^2 + 3*x2^2*y2^2
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: X = WehlerK3Surface([Z, Y])
sage: X.cardinality()
```

class sage.dynamics.arithmetic_dynamics.wehlerK3.WehlerK3Surface_ring(polys)

Bases: sage.schemes.product_projective.subscheme.AlgebraicScheme_subscheme_product_projective.subscheme.algebraicScheme_subscheme_product_projective.subscheme.algebraicScheme_subscheme.algebraicScheme_subscheme.algebraicScheme_subscheme.algebraicScheme_subscheme.algebraicScheme_subscheme.algebraicScheme_subscheme.algebraicScheme_subscheme.algebraicScheme_subscheme.algebraicScheme_subscheme.algebraicScheme_subscheme.algebraicScheme_subscheme.algebraicScheme_subscheme.algebraicScheme_subscheme.algebraicSche

A K3 surface in $\mathbb{P}^2 \times \mathbb{P}^2$ defined as the intersection of a bilinear and biquadratic form. [Weh1998]

EXAMPLES:

```
sage: R.<x,y,z,u,v,w> = PolynomialRing(QQ, 6)
sage: L = x*u - y*v
sage: Q = x*y*v^2 + z^2*u*w
sage: WehlerK3Surface([L, Q])
Closed subscheme of Product of projective spaces P^2 x P^2 over Rational
Field defined by:
    x*u - y*v,
    x*y*v^2 + z^2*u*w
```

Gpoly (*component*, *k*)

Returns the G polynomials G_k^* .

They are defined as: $G_k^* = (L_j^*)^2 Q_{ii}^* - L_i^* L_j^* Q_{ij}^* + (L_i^*)^2 Q_{jj}^*$ where $\{i, j, k\}$ is some permutation of (0, 1, 2) and * is either x (Component = 1) or y (Component = 0).

INPUT:

- component Integer: 0 or 1
- k Integer: 0, 1 or 2

OUTPUT: polynomial in terms of either y (Component = 0) or x (Component = 1)

EXAMPLES:

```
sage: R.<x0,x1,x2,y0,y1,y2> = PolynomialRing(ZZ, 6)
sage: Y = x0*y0 + x1*y1 - x2*y2
sage: Z = x0^2*y0*y1 + x0^2*y2^2 - x0*x1*y1*y2 + x1^2*y2*y1 \
+ x2^2*y2^2 + x2^2*y1^2 + x1^2*y2^2
sage: X = WehlerK3Surface([Z, Y])
sage: X.Gpoly(1, 0)
x0^2*x1^2 + x1^4 - x0*x1^2*x2 + x1^3*x2 + x1^2*x2^2 + x2^4
```

Hpoly (component, i, j)

Returns the H polynomials defined as H_{ii}^* .

This polynomial is defined by:

 $H_{ij}^* = 2L_i^*L_j^*Q_{kk}^* - L_i^*L_k^*Q_{jk}^* - L_j^*L_k^*Q_{ik}^* + (L_k^*)^2Q_{ij}^*$ where {i, j, k} is some permutation of (0, 1, 2) and * is either y (Component = 0) or x (Component = 1).

INPUT:

• component - Integer: 0 or 1

- i Integer: 0, 1 or 2
- \(\dagger\) Integer: 0, 1 or 2

OUTPUT: polynomial in terms of either y (Component = 0) or x (Component = 1)

EXAMPLES:

```
sage: R.<x0,x1,x2,y0,y1,y2> = PolynomialRing(ZZ, 6)
sage: Y = x0*y0 + x1*y1 - x2*y2
sage: Z = x0^2*y0*y1 + x0^2*y2^2 - x0*x1*y1*y2 + x1^2*y2*y1 \
+ x2^2*y2^2 + x2^2*y1^2 + x1^2*y2^2
sage: X = WehlerK3Surface([Z, Y])
sage: X.Hpoly(0, 1, 0)
2*y0*y1^3 + 2*y0*y1*y2^2 - y1*y2^3
```

Lxa(a)

Function will return the L polynomial defining the fiber, given by L_a^x .

This polynomial is defined as:

```
L_a^x = \{(a, y) \in \mathbb{P}^2 \times \mathbb{P}^2 : L(a, y) = 0\}.
```

Notation and definition from: [CS1996]

INPUT: a - Point in \mathbb{P}^2

OUTPUT: A polynomial representing the fiber

EXAMPLES:

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: Z = x0^2*y0^2 + 3*x0*x1*y0^2 + x1^2*y0^2 + 4*x0^2*y0*y1 \
+ 3*x0*x1*y0*y1 - 2*x2^2*y0*y1 - \
x0^2*y1^2 + 2*x1^2*y1^2 - x0*x2*y1^2 - 4*x1*x2*y1^2 \
+ 5*x0*x2*y0*y2 - 4*x1*x2*y0*y2 + 7*x0^2*y1*y2 \
+ 4*x1^2*y1*y2 + x0*x1*y2^2 + 3*x2^2*y2^2
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: X = WehlerK3Surface([Z, Y])
sage: T = PP(1, 1, 0, 1, 0, 0);
sage: X.Lxa(T[0])
y0 + y1
```

Lyb (b)

Function will return a fiber by L_b^y .

This polynomial is defined as:

```
L_b^y = \{(x, b) \in \mathbb{P}^2 \times \mathbb{P}^2 : L(x, b) = 0\}.
```

Notation and definition from: [CS1996]

INPUT: b - Point in projective space

OUTPUT: A polynomial representing the fiber

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: Z =x0^2*y0^2 + 3*x0*x1*y0^2 + x1^2*y0^2 + 4*x0^2*y0*y1 \
+ 3*x0*x1*y0*y1 \
- 2*x2^2*y0*y1 - x0^2*y1^2 + 2*x1^2*y1^2 - x0*x2*y1^2 - 4*x1*x2*y1^2 \
+ 5*x0*x2*y0*y2 \
- 4*x1*x2*y0*y2 + 7*x0^2*y1*y2 + 4*x1^2*y1*y2 + x0*x1*y2^2 + 3*x2^2*y2^2
```

```
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: X = WehlerK3Surface([Z, Y])
sage: T = PP(1, 1, 0, 1, 0, 0);
sage: X.Lyb(T[1])
x0
```

Qxa(a)

Function will return the Q polynomial defining a fiber given by Q_a^x .

This polynomial is defined as:

$$Q_a^x = \{(a, y) \in \mathbb{P}^2 \times \mathbb{P}^2 : Q(a, y) = 0\}.$$

Notation and definition from: [CS1996]

INPUT: a - Point in \mathbb{P}^2

OUTPUT: A polynomial representing the fiber

EXAMPLES:

Qyb(b)

Function will return a fiber by Q_b^y .

This polynomial is defined as:

$$Q_b^y = \{(x, b) \in \mathbb{P}^2 \times \mathbb{P}^2 : Q(x, b) = 0\}.$$

Notation and definition from: [CS1996]

INPUT: b - Point in projective space

OUTPUT: A polynomial representing the fiber

EXAMPLES:

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: Z = x0^2*y0^2 + 3*x0*x1*y0^2 + x1^2*y0^2 + 4*x0^2*y0*y1 \
+ 3*x0*x1*y0*y1 - 2*x2^2*y0*y1 - x0^2*y1^2 + 2*x1^2*y1^2 - x0*x2*y1^2 \
- 4*x1*x2*y1^2 + 5*x0*x2*y0*y2 - 4*x1*x2*y0*y2 + 7*x0^2*y1*y2 \
+ 4*x1^2*y1*y2 + x0*x1*y2^2 + 3*x2^2*y2^2
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: X = WehlerK3Surface([Z, Y])
sage: T = PP(1, 1, 0, 1, 0, 0);
sage: X.Qyb(T[1])
x0^2 + 3*x0*x1 + x1^2
```

Ramification_poly(i)

Function will return the Ramification polynomial g^* .

This polynomial is defined by:

$$g^* = \frac{(H_{ij}^*)^2 - 4G_i^* G_j^*}{(L_k^*)^2}.$$

The roots of this polynomial will either be degenerate fibers or fixed points of the involutions σ_x or σ_y for more information, see [CS1996].

INPUT: i - Integer, either 0 (polynomial in y) or 1 (polynomial in x)

OUTPUT: Polynomial in the coordinate ring of the ambient space

EXAMPLES:

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: Z = x0^2*y0^2 + 3*x0*x1*y0^2 + x1^2*y0^2 + 4*x0^2*y0*y1 + 3*x0*x1*y0*y1\
- 2*x2^2*y0*y1 - x0^2*y1^2 + 2*x1^2*y1^2 - x0*x2*y1^2 - 4*x1*x2*y1^2 + \_
\[ \times 5*x0*x2*y0*y2\\
- 4*x1*x2*y0*y2 + 7*x0^2*y1*y2 + 4*x1^2*y1*y2 + x0*x1*y2^2 + 3*x2^2*y2^2
\]
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: X = WehlerK3Surface([Z, Y])
sage: X.Ramification_poly(0)
8*y0^5*y1 - 24*y0^4*y1^2 + 48*y0^2*y1^4 - 16*y0*y1^5 + y1^6 + 84*y0^3*y1^2*y2
+ 46*y0^2*y1^3*y2 - 20*y0*y1^4*y2 + 16*y1^5*y2 + 53*y0^4*y2^2 + 56*y0^3*y1*y2^4
\[ \times 2
- 32*y0^2*y1^2*y2^2 - 80*y0*y1^3*y2^2 - 92*y1^4*y2^2 - 12*y0^2*y1*y2^3
- 168*y0*y1^2*y2^3 - 122*y1^3*y2^3 + 14*y0^2*y2^4 + 8*y0*y1*y2^4 - 112*y1^4
\[ \times 2*y2^4 + y2^6
\]
```

Sxa(a)

Function will return fiber by S_a^x .

This function is defined as:

$$S_a^x = L_a^x \cap Q_a^x$$
.

Notation and definition from: [CS1996]

INPUT: a - Point in \mathbb{P}^2

OUTPUT: A subscheme representing the fiber

EXAMPLES:

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: Z = x0^2*y0^2 + 3*x0*x1*y0^2 + x1^2*y0^2 + 4*x0^2*y0*y1 \
+ 3*x0*x1*y0*y1 \
- 2*x2^2*y0*y1 - x0^2*y1^2 + 2*x1^2*y1^2 - x0*x2*y1^2 - 4*x1*x2*y1^2 \
+ 5*x0*x2*y0*y2 \
- 4*x1*x2*y0*y2 + 7*x0^2*y1*y2 + 4*x1^2*y1*y2 + x0*x1*y2^2 + 3*x2^2*y2^2
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: X = WehlerK3Surface([Z, Y])
sage: T = PP(1, 1, 0, 1, 0, 0);
sage: X.Sxa(T[0])
Closed subscheme of Projective Space of dimension 2 over Rational Field.

defined by:
    y0 + y1,
    5*y0^2 + 7*y0*y1 + y1^2 + 11*y1*y2 + y2^2
```

$\mathbf{Syb}(b)$

Function will return fiber by S_b^y .

This function is defined by:

```
S_h^y = L_h^y \cap Q_h^y.
```

Notation and definition from: [CS1996]

INPUT: b - Point in \mathbb{P}^2

OUTPUT: A subscheme representing the fiber

EXAMPLES:

canonical_height (P, N, badprimes=None, prec=100)

Evaluates the canonical height for P with N terms of the series of the local heights.

ALGORITHM:

The sum of the canonical height minus and canonical height plus, for more info see section 4 of [CS1996].

INPUT:

- P a surface point
- N positive integer (number of terms of the series to use)
- badprimes (optional) list of integer primes (where the surface is degenerate)
- prec (default: 100) float point or p-adic precision

OUTPUT: A real number

EXAMPLES:

Call-Silverman example:

```
sage: set_verbose(None)
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: Z = x0^2*y0^2 + 3*x0*x1*y0^2 + x1^2*y0^2 + 4*x0^2*y0*y1 + 3*x0*x1*y0*y1__
```

```
2*x2^2*y0*y1 - x0^2*y1^2 + 2*x1^2*y1^2 - x0*x2*y1^2 - 4*x1*x2*y1^2 + ...

-5*x0*x2*y0*y2 \

-4*x1*x2*y0*y2 + 7*x0^2*y1*y2 + 4*x1^2*y1*y2 + x0*x1*y2^2 + 3*x2^2*y2^2

sage: Y = x0*y0 + x1*y1 + x2*y2

sage: X = WehlerK3Surface([Z, Y])

sage: P = X(0, 1, 0, 0, 0, 1)

sage: X.canonical_height(P, 4)

0.69826458668659859569990618895
```

canonical_height_minus (P, N, badprimes=None, prec=100)

Evaluates the canonical height minus function of Call-Silverman for P with N terms of the series of the local heights.

Must be over **Z** or **Q**.

ALGORITHM:

Sum over the lambda minus heights (local heights) in a convergent series, for more detail see section 7 of [CS1996].

INPUT:

- P a surface point
- N positive integer (number of terms of the series to use)
- badprimes (optional) list of integer primes (where the surface is degenerate)
- prec (default: 100) float point or p-adic precision

OUTPUT: A real number

EXAMPLES:

Call-Silverman example:

```
sage: set_verbose(None)
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: Z = x0^2*y0^2 + 3*x0*x1*y0^2 + x1^2*y0^2 + 4*x0^2*y0*y1 +\
    3*x0*x1*y0*y1 - 2*x2^2*y0*y1 - x0^2*y1^2 + 2*x1^2*y1^2 - x0*x2*y1^2 - \
    4*x1*x2*y1^2 + 5*x0*x2*y0*y2 - 4*x1*x2*y0*y2 + 7*x0^2*y1*y2 + 4*x1^2*y1*y2 + \
    \times\
    x0*x1*y2^2 + 3*x2^2*y2^2
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: X = WehlerK3Surface([Z, Y])
sage: P = X([0, 1, 0, 0, 0, 1])
sage: X.canonical_height_minus(P, 4) # long time
0.55073705369676788175590206734
```

canonical_height_plus (P, N, badprimes=None, prec=100)

Evaluates the canonical height plus function of Call-Silverman for P with N terms of the series of the local heights.

Must be over **Z** or **Q**.

ALGORITHM:

Sum over the lambda plus heights (local heights) in a convergent series, for more detail see section 7 of [CS1996].

INPUT:

- P a surface point
- N positive integer. Number of terms of the series to use
- badprimes (optional) list of integer primes (where the surface is degenerate)
- prec (default: 100) float point or p-adic precision

OUTPUT: A real number

EXAMPLES:

Call-Silverman Example:

```
sage: set_verbose(None)
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: Z = x0^2*y0^2 + 3*x0*x1*y0^2 + x1^2*y0^2 + 4*x0^2*y0*y1 + \
3*x0*x1*y0*y1 - 2*x2^2*y0*y1 - x0^2*y1^2 + 2*x1^2*y1^2 - x0*x2*y1^2 \
- 4*x1*x2*y1^2 + 5*x0*x2*y0*y2 - 4*x1*x2*y0*y2 + 7*x0^2*y1*y2 + 4*x1^2*y1*y2 \
+ x0*x1*y2^2 + 3*x2^2*y2^2
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: X = WehlerK3Surface([Z, Y])
sage: P = X([0, 1, 0, 0, 0, 1])
sage: X.canonical_height_plus(P, 4) # long time
0.14752753298983071394400412161
```

$change_ring(R)$

Changes the base ring on which the Wehler K3 Surface is defined.

INPUT: R - ring

OUTPUT: K3 Surface defined over input ring

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], GF(3))
sage: L = x0*y0 + x1*y1 - x2*y2
sage: Q = x0*x1*y1^2 + x2^2*y0*y2
```

```
sage: W = WehlerK3Surface([L, Q])
sage: W.base_ring()
Finite Field of size 3
sage: T = W.change_ring(GF(7))
sage: T.base_ring()
Finite Field of size 7
```

degenerate fibers()

Function will return the (rational) degenerate fibers of the surface defined over the base ring, or the fraction field of the base ring if it is not a field.

ALGORITHM:

The criteria for degeneracy by the common vanishing of the polynomials self.Gpoly(1, 0), self.Gpoly(1, 1), self.Gpoly(1, 2), self.Hpoly(1, 0, 1), "self.Hpoly(1, 0, 2)", self.Hpoly(1, 1, 2) (for the first component), is from Proposition 1.4 in the following article: [CS1996].

This function finds the common solution through elimination via Groebner bases by using the .variety() function on the three affine charts in each component.

OUTPUT: The output is a list of lists where the elements of lists are points in the appropriate projective space. The first list is the points whose pullback by the projection to the first component (projective space) is dimension greater than 0. The second list is points in the second component

EXAMPLES:

```
sage: R.<x0,x1,x2,y0,y1,y2> = PolynomialRing(ZZ, 6)
sage: Y = x0*y0 + x1*y1 - x2*y2
sage: Z = x0^2*y0*y1 + x0^2*y2^2 - x0*x1*y1*y2 + x1^2*y2*y1 + x2^2*y2^2\
+ x2^2*y1^2 + x1^2*y2^2
sage: X = WehlerK3Surface([Z, Y])
sage: X.degenerate_fibers()
[[], [(1 : 0 : 0)]]
```

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: Z = x0^2*y0^2 + 3*x0*x1*y0^2 + x1^2*y0^2 + 4*x0^2*y0*y1 + 3*x0*x1*y0*y1\
- 2*x2^2*y0*y1 - x0^2*y1^2 + 2*x1^2*y1^2 - x0*x2*y1^2 - 4*x1*x2*y1^2 + \
- 5*x0*x2*y0*y2\
- 4*x1*x2*y0*y2 + 7*x0^2*y1*y2 + 4*x1^2*y1*y2 + x0*x1*y2^2 + 3*x2^2*y2^2
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: X = WehlerK3Surface([Z, Y])
sage: X.degenerate_fibers()
[[], []]
```

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: R = PP.coordinate_ring()
sage: l = y0*x0 + y1*x1 + (y0 - y1)*x2
sage: q = (y1*y0 + y2^2)*x0^2 + ((y0^2 - y2*y1)*x1 + (y0^2 + (y1^2 - y2^2 + (y2*y0 + y1^2)*x1)*x2) + (y0^2 + (y1^2 + y2^2))*x2)*x0 \
+ (y2*y0 + y1^2)*x1^2 + (y0^2 + (-y1^2 + y2^2))*x2*x1
sage: X = WehlerK3Surface([1,q])
sage: X.degenerate_fibers()
[[(-1 : 1 : 1), (0 : 0 : 1)], [(-1 : -1 : 1), (0 : 0 : 1)]]
```

degenerate_primes (check=True)

Determine which primes p self has degenerate fibers over GF(p).

If check is False, then may return primes that do not have degenerate fibers. Raises an error if the surface is degenerate. Works only for $\mathbb{Z}\mathbb{Z}$ or $\mathbb{Q}\mathbb{Q}$.

INPUT: check – (default: True) boolean, whether the primes are verified

ALGORITHM:

p is a prime of bad reduction if and only if the defining polynomials of self plus the G and H polynomials have a common zero. Or stated another way, p is a prime of bad reduction if and only if the radical of the ideal defined by the defining polynomials of self plus the G and H polynomials is not (x_0, x_1, \ldots, x_N) . This happens if and only if some power of each x_i is not in the ideal defined by the defining polynomials of self (with G and H). This last condition is what is checked. The lcm of the coefficients of the monomials x_i in a groebner basis is computed. This may return extra primes.

OUTPUT: List of primes.

EXAMPLES:

```
sage: R.<x0,x1,x2,y0,y1,y2> = PolynomialRing(QQ, 6)
sage: L = y0*x0 + (y1*x1 + y2*x2)
sage: Q = (2*y0^2 + y2*y0 + (2*y1^2 + y2^2))*x0^2 + ((y0^2 + y1*y0 + \
    (y1^2 + 2*y2*y1 + y2^2))*x1 + (2*y1^2 + y2*y1 + y2^2)*x2)*x0 + ((2*y0^2\
    + (y1 + 2*y2)*y0 + (2*y1^2 + y2*y1))*x1^2 + ((2*y1 + 2*y2)*y0 + (y1^2 + \
    y2*y1 + 2*y2^2))*x2*x1 + (2*y0^2 + y1*y0 + (2*y1^2 + y2^2))*x2^2)
sage: X = WehlerK3Surface([L, Q])
sage: X.degenerate_primes()
[2, 3, 5, 11, 23, 47, 48747691, 111301831]
```

fiber (p, component)

Returns the fibers [y (component = 1) or x (Component = 0)] of a point on a K3 Surface, will work for nondegenerate fibers only.

For algorithm, see [Hutz2007].

INPUT:

-p - a point in \mathbb{P}^2

OUTPUT: The corresponding fiber (as a list)

```
sage: R.<x0,x1,x2,y0,y1,y2> = PolynomialRing(ZZ, 6)
sage: Y = x0*y0 + x1*y1 - x2*y2
sage: Z = y0^2*x0*x1 + y0^2*x2^2 - y0*y1*x1*x2 + y1^2*x2*x1 + y2^2*x2^2 +\
y2^2*x1^2 + y1^2*x2^2
sage: X = WehlerK3Surface([Z, Y])
sage: Proj = ProjectiveSpace(QQ, 2)
sage: P = Proj([1, 0, 0])
sage: X.fiber(P, 1)
Traceback (most recent call last):
...
TypeError: fiber is degenerate
```

```
sage: X.fiber(T, 1)
[(0:0:1,0:1:0), (0:0:1,2:0:0)]
```

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], GF(7))
sage: L = x0*y0 + x1*y1 - 1*x2*y2
sage: Q=(2*x0^2 + x2*x0 + (2*x1^2 + x2^2))*y0^2 + ((x0^2 + x1*x0 + (x1^2 + x2*x1 + x2^2))*y1 + \
    (2*x1^2 + x2*x1 + x2^2)*y1 + \
    (2*x1^2 + x2*x1 + x2^2)*y2)*y0 + ((2*x0^2 + (x1 + 2*x2)*x0 + (2*x1^2 + x2^2))*y1^2 + ((2*x1 + 2*x2)*x0 + \
    (x1^2 + x2*x1 + 2*x2^2))*y2*y1 + (2*x0^2 + x1*x0 + (2*x1^2 + x2^2))*y2^2)
sage: W = WehlerK3Surface([L, Q])
sage: W.fiber([4, 0, 1], 0)
[(0 : 1 : 0 , 4 : 0 : 1), (4 : 0 : 2 , 4 : 0 : 1)]
```

is_degenerate()

Function will return True if there is a fiber (over the algebraic closure of the base ring) of dimension greater than 0 and False otherwise.

OUTPUT: boolean

EXAMPLES:

```
sage: R.<x0,x1,x2,y0,y1,y2> = PolynomialRing(ZZ, 6)
sage: Y = x0*y0 + x1*y1 - x2*y2
sage: Z = x0^2*y0*y1 + x0^2*y2^2 - x0*x1*y1*y2 + x1^2*y2*y1 + x2^2*y2^2 + \
x2^2*y1^2 + x1^2*y2^2
sage: X = WehlerK3Surface([Z, Y])
sage: X.is_degenerate()
True
```

is_isomorphic(right)

Checks to see if two K3 surfaces have the same defining ideal.

INPUT:

• right - the K3 surface to compare to the original

OUTPUT: Boolean

EXAMPLES:

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: Z = x0^2*y0^2 + 3*x0*x1*y0^2 + x1^2*y0^2 + 4*x0^2*y0*y1 + \
3*x0*x1*y0*y1 -2*x2^2*y0*y1 - x0^2*y1^2 + 2*x1^2*y1^2 - x0*x2*y1^2 \
-4*x1*x2*y1^2 + 5*x0*x2*y0*y2 - 4*x1*x2*y0*y2 + 7*x0^2*y1*y2 + 4*x1^2*y1*y2 \
+ x0*x1*y2^2 + 3*x2^2*y2^2
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: X = WehlerK3Surface([Z, Y])
sage: W = WehlerK3Surface([Z + Y^2, Y])
sage: X.is_isomorphic(W)
True
```

```
sage: R.<x,y,z,u,v,w> = PolynomialRing(QQ, 6)
sage: L = x*u-y*v
sage: Q = x*y*v^2 + z^2*u*w
sage: W1 = WehlerK3Surface([L, Q])
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: L = x0*y0 + x1*y1 + x2*y2
sage: Q = x1^2*y0^2 + 2*x2^2*y0*y1 + x0^2*y1^2 -x0*x1*y2^2
sage: W2 = WehlerK3Surface([L, Q])
sage: W1.is_isomorphic(W2)
False
```

is smooth()

Function will return the status of the smoothness of the surface.

ALGORITHM:

Checks to confirm that all of the 2x2 minors of the Jacobian generated from the biquadratic and bilinear forms have no common vanishing points.

OUTPUT: Boolean

EXAMPLES:

```
sage: R.<x0,x1,x2,y0,y1,y2> = PolynomialRing(ZZ, 6)
sage: Y = x0*y0 + x1*y1 - x2*y2
sage: Z = x0^2*y0*y1 + x0^2*y2^2 - x0*x1*y1*y2 + x1^2*y2*y1 +\
    x2^2*y2^2 + x2^2*y1^2 + x1^2*y2^2
sage: X = WehlerK3Surface([Z, Y])
sage: X.is_smooth()
False
```

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: Z = x0^2*y0^2 + 3*x0*x1*y0^2 + x1^2*y0^2 + 4*x0^2*y0*y1 + \
3*x0*x1*y0*y1 - 2*x2^2*y0*y1 - x0^2*y1^2 + 2*x1^2*y1^2 - x0*x2*y1^2 \
- 4*x1*x2*y1^2 + 5*x0*x2*y0*y2 - 4*x1*x2*y0*y2 + 7*x0^2*y1*y2 + 4*x1^2*y1*y2 \
+ x0*x1*y2^2 + 3*x2^2*y2^2
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: X = WehlerK3Surface([Z, Y])
sage: X.is_smooth()
True
```

is_symmetric_orbit (orbit)

Checks to see if the orbit is symmetric (i.e. if one of the points on the orbit is fixed by 'sigma_x' or 'sigma_y').

INPUT:

orbit- a periodic cycle of either psi or phi

OUTPUT: Boolean

EXAMPLES:

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], GF(7))
sage: Z = x0^2*y0^2 + 3*x0*x1*y0^2 + x1^2*y0^2 + 4*x0^2*y0*y1 + 3*x0*x1*y0*y1

-\( \)
-2*x2^2*y0*y1 - x0^2*y1^2 + 2*x1^2*y1^2 - x0*x2*y1^2 - 4*x1*x2*y1^2 + \)
-5*x0*x2*y0*y2 \
-4*x1*x2*y0*y2 + 7*x0^2*y1*y2 + 4*x1^2*y1*y2 + x0*x1*y2^2 + 3*x2^2*y2^2
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: X = WehlerK3Surface([Z, Y])
sage: T = PP([0, 0, 1, 1, 0, 0])
sage: orbit = X.orbit_psi(T, 4)
sage: X.is_symmetric_orbit(orbit)
True
```

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: L = x0*y0 + x1*y1 + x2*y2
sage: Q = x1^2*y0^2 + 2*x2^2*y0*y1 + x0^2*y1^2 - x0*x1*y2^2
sage: W = WehlerK3Surface([L, Q])
sage: T = W([-1, -1, 1, 1, 0, 1])
sage: Orb = W.orbit_phi(T, 7)
sage: W.is_symmetric_orbit(Orb)
False
```

$lambda_minus(P, v, N, m, n, prec=100)$

Evaluates the local canonical height minus function of Call-Silverman at the place v for P with N terms of the series.

Use $\nabla = 0$ for the Archimedean place. Must be over **Z** or **Q**.

ALGORITHM:

Sum over local heights using convergent series, for more details, see section 4 of [CS1996].

INPUT:

- P a projective point
- N positive integer. number of terms of the series to use
- v non-negative integer. a place, use v = 0 for the Archimedean place
- m, n positive integers, We compute the local height for the divisor E_{mn}^+ . These must be indices of non-zero coordinates of the point P.
- prec (default: 100) float point or p-adic precision

OUTPUT: A real number

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: Z = x0^2*y0^2 + 3*x0*x1*y0^2 + x1^2*y0^2 + 4*x0^2*y0*y1 + 3*x0*x1*y0*y1

- 2*x2^2*y0*y1 - x0^2*y1^2 + 2*x1^2*y1^2 - x0*x2*y1^2 - 4*x1*x2*y1^2 +

- 5*x0*x2*y0*y2\

- 4*x1*x2*y0*y2 + 7*x0^2*y1*y2 + 4*x1^2*y1*y2 + x0*x1*y2^2 + 3*x2^2*y2^2
```

```
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: X = WehlerK3Surface([Z, Y])
sage: P = X([0, 0, 1, 1, 0, 0])
sage: X.lambda_minus(P, 2, 20, 2, 0, 200)
-0.18573351672047135037172805779671791488351056677474271893705
```

lambda_plus (P, v, N, m, n, prec=100)

Evaluates the local canonical height plus function of Call-Silverman at the place v for P with v terms of the series.

Use v = 0 for the archimedean place. Must be over **Z** or **Q**.

ALGORITHM:

Sum over local heights using convergent series, for more details, see section 4 of [CS1996].

INPUT:

- P a surface point
- N positive integer. number of terms of the series to use
- \forall non-negative integer. a place, use \forall = 0 for the Archimedean place
- m, n positive integers, We compute the local height for the divisor E_{mn}^+ . These must be indices of non-zero coordinates of the point P.
- prec (default: 100) float point or p-adic precision

OUTPUT: A real number

EXAMPLES:

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: Z = x0^2*y0^2 + 3*x0*x1*y0^2 + x1^2*y0^2 + 4*x0^2*y0*y1 + 3*x0*x1*y0*y1\
- 2*x2^2*y0*y1 - x0^2*y1^2 + 2*x1^2*y1^2 - x0*x2*y1^2 - 4*x1*x2*y1^2 + __

$\to 5*x0*x2*y0*y2\\
- 4*x1*x2*y0*y2 + 7*x0^2*y1*y2 + 4*x1^2*y1*y2 + x0*x1*y2^2 + 3*x2^2*y2^2
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: X = WehlerK3Surface([Z, Y])
sage: P = X([0, 0, 1, 1, 0, 0])
sage: X.lambda_plus(P, 0, 10, 2, 0)
0.89230705169161608922595928129
```

nth_iterate_phi(P, n, **kwds)

Computes the nth iterate for the phi function.

INPUT:

- \mathbb{P} - a point in $\mathbb{P}^2 \times \mathbb{P}^2$
- n an integer

kwds:

- check (default: True) boolean checks to see if point is on the surface
- normalize (default: False) boolean normalizes the point

OUTPUT: The nth iterate of the point given the phi function (if n is positive), or the psi function (if n is negative)

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: L = x0*y0 + x1*y1 + x2*y2
sage: Q = x1^2*y0^2 + 2*x2^2*y0*y1 + x0^2*y1^2 - x0*x1*y2^2
sage: W = WehlerK3Surface([L,Q])
sage: T = W([-1, -1, 1, 1, 0, 1])
sage: W.nth_iterate_phi(T, 7)
(-1: 0: 1, 1: -2: 1)
```

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: L = x0*y0 + x1*y1 + x2*y2
sage: Q = x1^2*y0^2 + 2*x2^2*y0*y1 + x0^2*y1^2 - x0*x1*y2^2
sage: W = WehlerK3Surface([L, Q])
sage: T = W([-1, -1, 1, 1, 0, 1])
sage: W.nth_iterate_phi(T, -7)
(1 : 0 : 1 , -1 : 2 : 1)
```

```
sage: R.<x0,x1,x2,y0,y1,y2>=PolynomialRing(QQ, 6)
sage: L = (-y0 - y1)*x0 + (-y0*x1 - y2*x2)
sage: Q = (-y2*y0 - y1^2)*x0^2 + ((-y0^2 - y2*y0 + (-y2*y1 - y2^2))*x1 + (-y0^2 - y2*y1)*x2)*x0 \
+ ((-y0^2 - y2*y0 - y2^2)*x1^2 + (-y2*y0 - y1^2)*x2*x1 + (-y0^2 + (-y1 - y2))*y0)*x2^2)
sage: X = WehlerK3Surface([L, Q])
sage: P = X([1, 0, -1, 1, -1, 0])
sage: X.nth_iterate_phi(P, 8) == X.nth_iterate_psi(P, 8)
True
```

nth iterate psi(P, n, **kwds)

Computes the nth iterate for the psi function.

INPUT:

- \mathbb{P} a point in $\mathbb{P}^2 \times \mathbb{P}^2$
- n an integer

kwds:

- check (default: True) boolean, checks to see if point is on the surface
- normalize (default: False) boolean, normalizes the point

OUTPUT: The nth iterate of the point given the psi function (if n is positive), or the phi function (if n is negative)

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: L = x0*y0 + x1*y1 + x2*y2
sage: Q = x1^2*y0^2 + 2*x2^2*y0*y1 + x0^2*y1^2 - x0*x1*y2^2
sage: W = WehlerK3Surface([L, Q])
sage: T = W([-1, -1, 1, 1, 0, 1])
sage: W.nth_iterate_psi(T, -7)
(-1 : 0 : 1 , 1 : -2 : 1)
```

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: L = x0*y0 + x1*y1 + x2*y2
sage: Q = x1^2*y0^2 + 2*x2^2*y0*y1 + x0^2*y1^2 - x0*x1*y2^2
sage: W = WehlerK3Surface([L, Q])
```

```
sage: T = W([-1, -1, 1, 1, 0, 1])
sage: W.nth_iterate_psi(T, 7)
(1 : 0 : 1 , -1 : 2 : 1)
```

$orbit_phi(P, N, **kwds)$

Returns the orbit of the ϕ function defined by $\phi = \sigma_u \circ \sigma_x$ Function is defined in [CS1996].

INPUT:

- P Point on the K3 surface
- N a non-negative integer or list or tuple of two non-negative integers

kwds:

- check (default: True) boolean, checks to see if point is on the surface
- normalize (default: False) boolean, normalizes the point

OUTPUT: List of points in the orbit

EXAMPLES:

```
sage: PP. \langle x0, x1, x2, y0, y1, y2 \rangle = ProductProjectiveSpaces([2, 2], QQ)
sage: Z = x0^2 + y0^2 + 3 \times x0 \times x1 \times y0^2 + x1^2 \times y0^2 + 4 \times x0^2 \times y0 \times y1 + 
3*x0*x1*y0*y1 - 2*x2^2*y0*y1 - x0^2*y1^2 + 2*x1^2*y1^2 - x0*x2*y1^2 - 
4*x1*x2*y1^2 + 5*x0*x2*y0*y2 - 4*x1*x2*y0*y2 + 7*x0^2*y1*y2 + 4*x1^2*y1*y2 + 
x0*x1*y2^2 + 3*x2^2*y2^2
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: X = WehlerK3Surface([Z, Y])
sage: T = PP(0, 0, 1, 1, 0, 0)
sage: X.orbit_phi(T,2, normalize = True)
[(0:0:1,1:0:0),(-1:0:1,0:1:0),(-12816/6659:55413/6659]
\hookrightarrow: 1 , 1 : 1/9 : 1)]
sage: X.orbit_phi(T,[2,3], normalize = True)
[(-12816/6659 : 55413/6659 : 1 , 1 : 1/9 : 1),
(7481279673854775690938629732119966552954626693713001783595660989241/
\hookrightarrow 18550615454277582153932951051931712107449915856862264913424670784695
: 3992260691327218828582255586014718568398539828275296031491644987908/
→18550615454277582153932951051931712107449915856862264913424670784695 :
1 , -117756062505511/54767410965117 : -23134047983794359/37466994368025041 :..
\hookrightarrow 1)
```

$orbit_psi(P, N, **kwds)$

Returns the orbit of the ψ function defined by $\psi = \sigma_x \circ \sigma_y$.

Function is defined in [CS1996].

INPUT:

- P a point on the K3 surface
- N a non-negative integer or list or tuple of two non-negative integers

kwds:

- check (default: True) boolean, checks to see if point is on the surface
- normalize (default: False) boolean, normalizes the point

OUTPUT: a list of points in the orbit

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: Z = x0^2*y0^2 + 3*x0*x1*y0^2 + x1^2*y0^2 + 4*x0^2*y0*y1 + \
3*x0*x1*y0*y1 -2*x2^2*y0*y1 - x0^2*y1^2 + 2*x1^2*y1^2 - x0*x2*y1^2 - \
4*x1*x2*y1^2 + 5*x0*x2*y0*y2 -4*x1*x2*y0*y2 + 7*x0^2*y1*y2 + 4*x1^2*y1*y2 + \
x0*x1*y2^2 + 3*x2^2*y2^2
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: X = WehlerK3Surface([Z, Y])
sage: T = X(0, 0, 1, 1, 0, 0)
sage: X.orbit_psi(T, 2, normalize = True)
[(0 : 0 : 1 , 1 : 0 : 0), (0 : 0 : 1 , 0 : 1 : 0), (-1 : 0 : 1 , 1 : 1/9 : 1)]
sage: X.orbit_psi(T,[2,3], normalize = True)
[(-1 : 0 : 1 , 1 : 1/9 : 1),
(-12816/6659 : 55413/6659 : 1 , -117756062505511/54767410965117 : -
→23134047983794359/37466994368025041 : 1)]
```

phi (*a*, ***kwds*)

Evaluates the function $\phi = \sigma_u \circ \sigma_x$.

ALGORITHM:

Refer to Section 6: "An algorithm to compute σ_x , σ_y , ϕ , and ψ " in [CS1996].

For the degenerate case refer to [FH2015].

INPUT:

• a - Point in $\mathbb{P}^2 \times \mathbb{P}^2$

kwds:

- check (default: True) boolean checks to see if point is on the surface
- normalize (default: True) boolean normalizes the point

OUTPUT: A point on this surface

EXAMPLES:

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: Z = x0^2*y0^2 + 3*x0*x1*y0^2 + x1^2*y0^2 + 4*x0^2*y0*y1 + \
3*x0*x1*y0*y1 -2*x2^2*y0*y1 - x0^2*y1^2 + 2*x1^2*y1^2 - x0*x2*y1^2 \
- 4*x1*x2*y1^2 + 5*x0*x2*y0*y2 -4*x1*x2*y0*y2 + 7*x0^2*y1*y2 + 4*x1^2*y1*y2 \
+ x0*x1*y2^2 + 3*x2^2*y2^2
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: X = WehlerK3Surface([Z, Y])
sage: T = PP([0, 0, 1, 1, 0, 0])
sage: X.phi(T)
(-1 : 0 : 1 , 0 : 1 : 0)
```

psi (*a*, ***kwds*)

Evaluates the function $\psi = \sigma_x \circ \sigma_y$.

ALGORITHM:

Refer to Section 6: "An algorithm to compute σ_x , σ_y , ϕ , and ψ " in [CS1996].

For the degenerate case refer to [FH2015].

INPUT:

• a - Point in $\mathbb{P}^2 \times \mathbb{P}^2$

kwds:

- check (default: True) boolean checks to see if point is on the surface
- normalize (default: True) boolean normalizes the point

OUTPUT: A point on this surface

EXAMPLES:

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: Z = x0^2*y0^2 + 3*x0*x1*y0^2 + x1^2*y0^2 + 4*x0^2*y0*y1 + \
3*x0*x1*y0*y1 -2*x2^2*y0*y1 - x0^2*y1^2 + 2*x1^2*y1^2 - x0*x2*y1^2 \
- 4*x1*x2*y1^2 + 5*x0*x2*y0*y2 - 4*x1*x2*y0*y2 + 7*x0^2*y1*y2 + 4*x1^2*y1*y2 \
+ x0*x1*y2^2 + 3*x2^2*y2^2
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: X = WehlerK3Surface([Z, Y])
sage: T = PP([0, 0, 1, 1, 0, 0])
sage: X.psi(T)
(0 : 0 : 1 , 0 : 1 : 0)
```

sigmaX(P, **kwds)

Function returns the involution on the Wehler K3 surface induced by the double covers.

In particular, it fixes the projection to the first coordinate and swaps the two points in the fiber, i.e. $(x, y) \rightarrow (x, y')$. Note that in the degenerate case, while we can split fiber into pairs of points, it is not always possible distinguish them, using this algorithm.

ALGORITHM:

Refer to Section 6: "An algorithm to compute σ_x , σ_y , ϕ , and ψ " in [CS1996FH2015. For the degenerate case refer to [FH2015].

INPUT:

• P - a point in $\mathbb{P}^2 imes \mathbb{P}^2$

kwds:

- check (default: True) boolean checks to see if point is on the surface
- normalize (default: True) boolean normalizes the point

OUTPUT: A point on the K3 surface

EXAMPLES:

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: Z = x0^2*y0^2 + 3*x0*x1*y0^2 + x1^2*y0^2 + 4*x0^2*y0*y1 +\
3*x0*x1*y0*y1 -2*x2^2*y0*y1 - x0^2*y1^2 + 2*x1^2*y1^2 - x0*x2*y1^2 -\
4*x1*x2*y1^2 + 5*x0*x2*y0*y2 -4*x1*x2*y0*y2 + 7*x0^2*y1*y2 +\
4*x1^2*y1*y2 + x0*x1*y2^2 + 3*x2^2*y2^2
sage: Y = x0*y0 + x1*y1 + x2*y2
sage: X = WehlerK3Surface([Z, Y])
sage: T = PP(0, 0, 1, 1, 0, 0)
sage: X.sigmaX(T)
(0 : 0 : 1 , 0 : 1 : 0)
```

degenerate examples:

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: l = y0*x0 + y1*x1 + (y0 - y1)*x2
sage: q = (y1*y0)*x0^2 + ((y0^2)*x1 + (y0^2 + (y1^2 - y2^2))*x2)*x0\
+ (y2*y0 + y1^2)*x1^2 + (y0^2 + (-y1^2 + y2^2))*x2*x1
sage: X = WehlerK3Surface([1, q])
```

```
sage: X.sigmaX(X([1, 0, 0, 0, 1, -2]))
(1 : 0 : 0 , 0 : 1/2 : 1)
sage: X.sigmaX(X([1, 0, 0, 0, 0, 1]))
(1 : 0 : 0 , 0 : 0 : 1)
sage: X.sigmaX(X([-1, 1, 1, -1, -1, 1]))
(-1 : 1 : 1 , 2 : 2 : 1)
sage: X.sigmaX(X([0, 0, 1, 1, 1, 0]))
(0 : 0 : 1 , 1 : 1 : 0)
sage: X.sigmaX(X([0, 0, 1, 1, 1, 1]))
(0 : 0 : 1 , -1 : -1 : 1)
```

Case where we cannot distinguish the two points:

```
sage: PP.<y0,y1,y2,x0,x1,x2>=ProductProjectiveSpaces([2, 2], GF(3))
sage: 1 = x0*y0 + x1*y1 + x2*y2
sage: q=-3*x0^2*y0^2 + 4*x0*x1*y0^2 - 3*x0*x2*y0^2 - 5*x0^2*y0*y1 - \
190*x0*x1*y0*y1 - 5*x1^2*y0*y1 + 5*x0*x2*y0*y1 + 14*x1*x2*y0*y1 + \
5*x2^2*y0*y1 - x0^2*y1^2 - 6*x0*x1*y1^2 - 2*x1^2*y1^2 + 2*x0*x2*y1^2 - \
4*x2^2*y1^2 + 4*x0^2*y0*y2 - x1^2*y0*y2 + 3*x0*x2*y0*y2 + 6*x1*x2*y0*y2 - \
6*x0^2*y1*y2 - 4*x0*x1*y1*y2 - x1^2*y1*y2 + 51*x0*x2*y1*y2 - 7*x1*x2*y1*y2 - \
9*x2^2*y1*y2 - x0^2*y2^2 - 4*x0*x1*y2^2 + 4*x1^2*y2^2 - x0*x2*y2^2 + \
-13*x1*x2*y2^2 - x2^2*y2^2
sage: X = WehlerK3Surface([1, q])
sage: P = X([1, 0, 0, 0, 1, 1])
sage: X.sigmaX(X.sigmaX(P))
Traceback (most recent call last):
...
ValueError: cannot distinguish points in the degenerate fiber
```

sigmaY(P, **kwds)

Function returns the involution on the Wehler K3 surfaces induced by the double covers.

In particular, it fixes the projection to the second coordinate and swaps the two points in the fiber, i.e. $(x,y) \to (x',y)$. Note that in the degenerate case, while we can split the fiber into two points, it is not always possible distinguish them, using this algorithm.

ALGORITHM:

Refer to Section 6: "An algorithm to compute σ_x , σ_y , ϕ , and ψ " in [CS1996]. For the degenerate case refer to [FH2015].

INPUT:

• P - a point in $\mathbb{P}^2 \times \mathbb{P}^2$

kwds:

- check (default: True) boolean checks to see if point is on the surface
- normalize (default: True) boolean normalizes the point

OUTPUT: A point on the K3 surface

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: Z = x0^2 \times y0^2 + 3 \times x0 \times x1 \times y0^2 + x1^2 \times y0^2 + 4 \times x0^2 \times y0 \times y1 + \\ 3 \times x0 \times x1 \times y0 \times y1 - 2 \times x2^2 \times y0 \times y1 - x0^2 \times y1^2 + 2 \times x1^2 \times y1^2 - x0 \times x2 \times y1^2 \\ - 4 \times x1 \times x2 \times y1^2 + 5 \times x0 \times x2 \times y0 \times y2 - 4 \times x1 \times x2 \times y0 \times y2 + 7 \times x0^2 \times y1 \times y2 + 4 \times x1^2 \times y1 \times y2 \\ + x0 \times x1 \times y2^2 + 3 \times x2^2 \times y2^2 \\ sage: Y = x0 \times y0 + x1 \times y1 + x2 \times y2
```

```
sage: X = WehlerK3Surface([Z, Y])
sage: T = PP(0, 0, 1, 1, 0, 0)
sage: X.sigmaY(T)
(0 : 0 : 1 , 1 : 0 : 0)
```

degenerate examples:

```
sage: PP.<x0,x1,x2,y0,y1,y2> = ProductProjectiveSpaces([2, 2], QQ)
sage: 1 = y0*x0 + y1*x1 + (y0 - y1)*x2
sage: q = (y1*y0)*x0^2 + ((y0^2)*x1 + (y0^2 + (y1^2 - y2^2))*x2)*x0 +\
    (y2*y0 + y1^2)*x1^2 + (y0^2 + (-y1^2 + y2^2))*x2*x1
sage: X = WehlerK3Surface([1, q])
sage: X.sigmaY(X([1, -1, 0, -1, -1, 1]))
(1/10: -1/10: 1, -1: -1: 1)
sage: X.sigmaY(X([0, 0, 1, -1, -1, 1]))
(-4: 4: 1, -1: -1: 1)
sage: X.sigmaY(X([1, 2, 0, 0, 0, 1]))
(-3: -3: 1, 0: 0: 1)
sage: X.sigmaY(X([1, 1, 1, 0, 0, 1]))
(1: 0: 0, 0: 0: 1)
```

Case where we cannot distinguish the two points:

```
sage: PP. < x0, x1, x2, y0, y1, y2 > = ProductProjectiveSpaces([2, 2], GF(3))
sage: 1 = x0*y0 + x1*y1 + x2*y2
sage: q=-3*x0^2*y0^2 + 4*x0*x1*y0^2 - 3*x0*x2*y0^2 - 5*x0^2*y0*y1 - __
\hookrightarrow 190 \times x0 \times x1 \times y0 \times y1 \
-5*x1^2*y0*y1 + 5*x0*x2*y0*y1 + 14*x1*x2*y0*y1 + 5*x2^2*y0*y1 - x0^2*y1^2 - ...
\hookrightarrow 6*x0*x1*y1^2 \
-2*x1^2*y1^2 + 2*x0*x2*y1^2 - 4*x2^2*y1^2 + 4*x0^2*y0*y2 - x1^2*y0*y2 + ...
\rightarrow 3 * x 0 * x 2 * y 0 * y 2 \
\rightarrow 7*x1*x2*y1*y2 \
-9*x2^2*y1*y2 - x0^2*y2^2 - 4*x0*x1*y2^2 + 4*x1^2*y2^2 - x0*x2*y2^2 + ...
\hookrightarrow 13 \times x1 \times x2 \times y2^2 - x2^2 \times y2^2
sage: X = WehlerK3Surface([1 ,q])
sage: P = X([0, 1, 1, 1, 0, 0])
sage: X.sigmaY(X.sigmaY(P))
Traceback (most recent call last):
ValueError: cannot distinguish points in the degenerate fiber
```

sage.dynamics.arithmetic_dynamics.wehlerK3.random_WehlerK3Surface(PP)

Produces a random K3 surface in $\mathbb{P}^2 \times \mathbb{P}^2$ defined as the intersection of a bilinear and biquadratic form. [Weh1998]

INPUT: Projective space cartesian product

OUTPUT: WehlerK3Surface_ring

```
- x0^2*y0*y2 - x0*x1*y0*y2 + x0*x2*y0*y2 + x1*x2*y0*y2 + x0*x1*y1*y2
- x1^2*y1*y2 - x1*x2*y1*y2 - x0^2*y2^2 + x0*x1*y2^2 - x1^2*y2^2 - x0*x2*y2^2
```

See also:

- sage.schemes.affine.affine_morphism
- sage.schemes.projective.projective_morphism
- sage.schemes.product_projective.morphism

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