Sage Reference Manual: Statistics Release 7.4

The Sage Development Team

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CHAPTER

ONE

BASIC STATISTICS

This file contains basic descriptive functions. Included are the mean, median, mode, moving average, standard deviation, and the variance. When calling a function on data, there are checks for functions already defined for that data type.

The mean function returns the arithmetic mean (the sum of all the members of a list, divided by the number of members). Further revisions may include the geometric and harmonic mean. The median function returns the number separating the higher half of a sample from the lower half. The mode returns the most common occuring member of a sample, plus the number of times it occurs. If entries occur equally common, the smallest of a list of the most common entries is returned. The moving_average is a finite impulse response filter, creating a series of averages using a user-defined number of subsets of the full data set. The std and the variance return a measurement of how far data points tend to be from the arithmetic mean.

Functions are available in the namespace stats, i.e. you can use them by typing stats.mean, stats.median.etc.

REMARK: If all the data you are working with are floating point numbers, you may find finance. TimeSeries helpful, since it is extremely fast and offers many of the same descriptive statistics as in the module.

AUTHOR:

• Andrew Hou (11/06/2009)

```
sage.stats.basic_stats.mean (v) Return the mean of the elements of v.
```

We define the mean of the empty list to be the (symbolic) NaN, following the convention of MATLAB, Scipy, and R.

INPUT:

•v – a list of numbers

OUTPUT:

•a number

```
sage: mean([pi, e])
1/2*pi + 1/2*e
sage: mean([])
NaN
sage: mean([I, sqrt(2), 3/5])
1/3*sqrt(2) + 1/3*I + 1/5
sage: mean([RIF(1.0103,1.0103), RIF(2)])
1.50515000000000000?
sage: mean(range(4))
3/2
```

```
sage: v = finance.TimeSeries([1..100])
sage: mean(v)
50.5
```

```
sage.stats.basic_stats.median (v)
```

Return the median (middle value) of the elements of v

If v is empty, we define the median to be NaN, which is consistent with NumPy (note that R returns NULL). If v is comprised of strings, TypeError occurs. For elements other than numbers, the median is a result of sorted().

INPUT:

•v – a list

OUTPUT:

•median element of v

EXAMPLES:

```
sage.stats.basic_stats.mode(v)
```

Return the mode of v. The mode is the sorted list of the most frequently occurring elements in v. If n is the most times that any element occurs in v, then the mode is the sorted list of elements of v that occur n times.

NOTE: The elements of v must be hashable and comparable.

INPUT:

•v – a list

OUTPUT:

•a list

```
sage: v = [1,2,4,1,6,2,6,7,1]
sage: mode(v)
[1]
sage: v.count(1)
3
sage: mode([])
[]
sage: mode([1,2,3,4,5])
[1, 2, 3, 4, 5]
sage: mode([3,1,2,1,2,3])
[1, 2, 3]
```

```
sage: mode(['sage', 4, I, 3/5, 'sage', pi])
['sage']
sage: class MyClass:
...     def mode(self):
...         return [1]
sage: stats.mode(MyClass())
[1]
```

```
sage.stats.basic_stats.moving_average ( v, n)
```

Provides the moving average of a list v

The moving average of a list is often used to smooth out noisy data.

If v is empty, we define the entries of the moving average to be NaN.

INPUT:

- $\bullet v$ a list
- $\bullet n$ the number of values used in computing each average.

OUTPUT:

•a list of length len (v) -n+1, since we do not fabric any values

EXAMPLES:

```
sage: moving_average([1..10], 1)
[1, 2, 3, 4, 5, 6, 7, 8, 9, 10]
sage: moving_average([1..10], 4)
[5/2, 7/2, 9/2, 11/2, 13/2, 15/2, 17/2]
sage: moving_average([], 1)
[]
sage: moving_average([pi, e, I, sqrt(2), 3/5], 2)
[1/2*pi + 1/2*e, 1/2*e + 1/2*I, 1/2*sqrt(2) + 1/2*I, 1/2*sqrt(2) + 3/10]
```

We check if the input is a time series, and if so use the optimized simple_moving_average method, but with (slightly different) meaning as defined above (the point is that the simple_moving_average on time series returns n values:

```
sage: a = finance.TimeSeries([1..10])
sage: stats.moving_average(a, 3)
[2.0000, 3.0000, 4.0000, 5.0000, 6.0000, 7.0000, 8.0000, 9.0000]
sage: stats.moving_average(list(a), 3)
[2.0, 3.0, 4.0, 5.0, 6.0, 7.0, 8.0, 9.0]
```

```
sage.stats.basic_stats. std (v, bias=False)
```

Returns the standard deviation of the elements of v

We define the standard deviation of the empty list to be NaN, following the convention of MATLAB, Scipy, and R.

INPUT:

- •v a list of numbers
- •bias bool (default: False); if False, divide by len(v) 1 instead of len(v) to give a less biased estimator (sample) for the standard deviation.

OUTPUT:

•a number

EXAMPLES:

```
sage: std([1..6], bias=True)
1/2*sqrt(35/3)
sage: std([1..6], bias=False)
sqrt (7/2)
sage: std([e, pi])
sqrt(1/2)*sqrt((pi - e)^2)
sage: std([])
NaN
sage: std([I, sqrt(2), 3/5])
sqrt(1/450*(10*sqrt(2) - 5*I - 3)^2 + 1/450*(5*sqrt(2) - 10*I + 3)^2 + 1/450*(10*sqrt(2) - 10*I + 3)^3 + 1/450*(10*sqrt(2) - 10*I + 3)^4 + 1
  450*(5*sqrt(2) + 5*I - 6)^2
sage: std([RIF(1.0103, 1.0103), RIF(2)])
0.6998235813403261?
sage: import numpy
sage: x = numpy.array([1, 2, 3, 4, 5])
sage: std(x, bias=False)
1.5811388300841898
sage: x = finance.TimeSeries([1..100])
sage: std(x)
29.011491975882016
```

sage.stats.basic_stats.variance (v, bias=False)

Returns the variance of the elements of v

We define the variance of the empty list to be NaN, following the convention of MATLAB, Scipy, and R.

INPUT:

•v – a list of numbers

•bias - bool (default: False); if False, divide by len(v) - 1 instead of len(v) to give a less biased estimator (sample) for the standard deviation.

OUTPUT:

•a number

```
sage: variance([1..6])
 7/2
sage: variance([1..6], bias=True)
35/12
sage: variance([e, pi])
1/2*(pi - e)^2
sage: variance([])
sage: variance([I, sqrt(2), 3/5])
1/450*(10*sqrt(2) - 5*I - 3)^2 + 1/450*(5*sqrt(2) - 10*I + 3)^2 + 1/450*(5*sqrt(2) - 10*I + 3)^3 + 1/450*(10*sqrt(2) - 10*I + 3)^4 + 1/450*(
  450*(5*sqrt(2) + 5*I - 6)^2
sage: variance([RIF(1.0103, 1.0103), RIF(2)])
0.4897530450000000?
sage: import numpy
sage: x = numpy.array([1, 2, 3, 4, 5])
sage: variance(x, bias=False)
2.5
sage: x = finance.TimeSeries([1..100])
sage: variance(x)
841.666666666666
```

```
sage: variance(x, bias=True)
833.25
sage: class MyClass:
      def variance(self, bias = False):
           return 1
sage: stats.variance(MyClass())
sage: class SillyPythonList:
        def __init__(self):
. . .
           self.__list = [2L,4L]
. . .
      def __len__(self):
. . .
            return len(self.__list)
. . .
        def __iter__(self):
. . .
            return self.__list.__iter__()
. . .
        def mean(self):
. . .
            return 3L
. . .
sage: R = SillyPythonList()
sage: variance(R)
sage: variance(R, bias=True)
```

TESTS:

The performance issue from #10019 is solved:

```
sage: variance([1] * 2^18)
0
```

CHAPTER

TWO

CINTLISTS

This is a class for fast basic operations with lists of C ints. It is similar to the double precision TimeSeries class. It has all the standard C int semantics, of course, including overflow. It is also similar to the Python list class, except all elements are C ints, which makes some operations much, much faster. For example, concatenating two IntLists can be over 10 times faster than concatenating the corresponding Python lists of ints, and taking slices is also much faster.

AUTHOR:

• William Stein, 2010-03

```
class sage.stats.intlist. IntList
    Bases: object
    A list of C int's.
    list ()
```

Return Python list version of self with Python ints as entries.

EXAMPLES:

```
sage: a = stats.IntList([1..15]); a
[1, 2, 3, 4, 5 ... 11, 12, 13, 14, 15]
sage: a.list()
[1, 2, 3, 4, 5, 6, 7, 8, 9, 10, 11, 12, 13, 14, 15]
sage: list(a) == a.list()
True
sage: type(a.list()[0])
<type 'int'>
```

max (index=False)

Return the largest value in this time series. If this series has length 0 we raise a ValueError

INPUT:

•index – bool (default: False); if True, also return index of maximum entry.

OUTPUT:

•int – largest value

•int – index of largest value; only returned if index=True

```
sage: v = stats.IntList([1,-4,3,-2,-4,3])
sage: v.max()
3
sage: v.max(index=True)
(3, 2)
```

min (index=False)

Return the smallest value in this integer list. If this series has length 0 we raise a ValueError.

INPUT:

•index – bool (default: False); if True, also return index of minimal entry.

OUTPUT:

- •float smallest value
- •integer index of smallest value; only returned if index=True

EXAMPLES:

```
sage: v = stats.IntList([1,-4,3,-2,-4])
sage: v.min()
-4
sage: v.min(index=True)
(-4, 1)
```

plot (*args, **kwds)

Return a plot of this IntList. This just constructs the corresponding double-precision floating point Time-Series object, passing on all arguments.

EXAMPLES:

```
sage: stats.IntList([3,7,19,-2]).plot()
Graphics object consisting of 1 graphics primitive
sage: stats.IntList([3,7,19,-2]).plot(color='red',pointsize=50,points=True)
Graphics object consisting of 1 graphics primitive
```

plot_histogram (*args, **kwds)

Return a histogram plot of this IntList. This just constructs the corresponding double-precision floating point TimeSeries object, and plots it, passing on all arguments.

EXAMPLES:

```
sage: stats.IntList([1..15]).plot_histogram()
Graphics object consisting of 50 graphics primitives
```

prod()

Return the product of the entries of self.

EXAMPLES:

```
sage: a = stats.IntList([1..10]); a
[1, 2, 3, 4, 5, 6, 7, 8, 9, 10]
sage: a.prod()
3628800
sage: factorial(10)
3628800
```

Note that there can be overflow:

```
sage: a = stats.IntList([2^30, 2]); a
[1073741824, 2]
sage: a.prod()
-2147483648
```

sum ()

Return the sum of the entries of self.

EXAMPLES:

```
sage: stats.IntList([1..100]).sum()
5050
```

Note that there can be overflow, since the entries are C ints:

```
sage: a = stats.IntList([2^30,2^30]); a
[1073741824, 1073741824]
sage: a.sum()
-2147483648
```

time_series ()

Return TimeSeries version of self, which involves changing each entry to a double.

EXAMPLES:

```
sage: T = stats.IntList([-2,3,5]).time_series(); T
[-2.0000, 3.0000, 5.0000]
sage: type(T)
<type 'sage.finance.time_series.TimeSeries'>
```

sage.stats.intlist.unpickle_intlist_v1 (v, n)

Version 1 unpickle method.

INPUT: v - a raw char buffer

```
sage: v = stats.IntList([1,2,3])
sage: s = v.__reduce__()[1][0]
sage: type(s)
<type 'str'>
sage: sage.stats.intlist.unpickle_intlist_v1(s, 3)
[1, 2, 3]
sage: sage.stats.intlist.unpickle_intlist_v1(s+s,6)
[1, 2, 3, 1, 2, 3]
sage: sage.stats.intlist.unpickle_intlist_v1('',0)
[]
```

CHAPTER

THREE

HIDDEN MARKOV MODELS

This is a complete pure-Cython optimized implementation of Hidden Markov Models. It fully supports Discrete, Gaussian, and Mixed Gaussian emissions.

The best references for the basic HMM algorithms implemented here are:

- Tapas Kanungo's "Hidden Markov Models"
- Jackson's HMM tutorial: http://personal.ee.surrey.ac.uk/Personal/P.Jackson/tutorial/

LICENSE: Some of the code in this file is based on reading Kanungo's GPLv2+ implementation of discrete HMM's, hence the present code must be licensed with a GPLv2+ compatible license.

AUTHOR:

• William Stein, 2010-03

```
{\bf class} \; {\tt sage.stats.hmm.hmm.} \; {\tt DiscreteHiddenMarkovModel} \\
```

Bases: sage.stats.hmm.hmm.HiddenMarkovModel

A discrete Hidden Markov model implemented using double precision floating point arithmetic.

INPUT:

- \bullet A a list of lists or a square N x N matrix, whose (i,j) entry gives the probability of transitioning from state i to state j.
- •B a list of N lists or a matrix with N rows, such that B[i,k] gives the probability of emitting symbol k while in state i.
- •pi the probabilities of starting in each initial state, i.e., pi[i] is the probability of starting in state i.
- •emission_symbols None or list (default: None); if None, the emission_symbols are the ints [0..N-1], where N is the number of states. Otherwise, they are the entries of the list emissions_symbols, which must all be hashable.
- •normalize -bool (default: True); if given, input is normalized to define valid probability distributions, e.g., the entries of A are made nonnegative and the rows sum to 1, and the probabilities in pi are normalized.

```
Initial probabilities: [0.5000, 0.5000]
sage: m.log_likelihood([0,1,0,1,0,1])
-4.66693474691329...
sage: m.viterbi([0,1,0,1,0,1])
([1, 1, 1, 1, 1, 1], -5.378832842208748)
sage: m.baum_welch([0,1,0,1,0,1])
(0.0, 22)
sage: m # rel tol 1e-10
Discrete Hidden Markov Model with 2 States and 2 Emissions
Transition matrix:
[1.0134345614745788e-70
                                           1.01
                   1.0 3.9974352713558623e-191
Emission matrix:
[ 7.380221566254936e-54
                   1.0 3.9974352626002193e-19]
Initial probabilities: [0.0000, 1.0000]
sage: m.sample(10)
[0, 1, 0, 1, 0, 1, 0, 1, 0, 1]
sage: m.graph().plot()
Graphics object consisting of 6 graphics primitives
```

A 3-state model that happens to always outputs 'b':

baum_welch (obs, max_iter=100, log_likelihood_cutoff=0.0001, fix_emissions=False)

Given an observation sequence obs, improve this HMM using the Baum-Welch algorithm to increase the probability of observing obs.

INPUT:

- •obs list of emissions
- •max_iter integer (default: 100) maximum number of Baum-Welch steps to take
- •log_likelihood_cutoff positive float (default: 1e-4); the minimal improvement in likelihood with respect to the last iteration required to continue. Relative value to log likelihood.
- •fix_emissions bool (default: False); if True, do not change emissions when updating

OUTPUT:

•changes the model in places, and returns the log likelihood and number of iterations.

```
[ 0.0 1.0]
Initial probabilities: [0.0000, 1.0000]
```

The following illustrates how Baum-Welch is only a local optimizer, i.e., the above model is far more likely to produce the sequence [1,0]*20 than the one we get below:

```
sage: m = hmm.DiscreteHiddenMarkovModel([[0.5,0.5],[0.5,0.5]], [[.5,.5],[.5,.

→5]], [.5,.5])
sage: m.baum_welch([1,0]*20, log_likelihood_cutoff=0)
(-27.725887222397784, 1)
sage: m
Discrete Hidden Markov Model with 2 States and 2 Emissions
Transition matrix:
[0.5 0.5]
[0.5 0.5]
Emission matrix:
[0.5 0.5]
[0.5 0.5]
Initial probabilities: [0.5000, 0.5000]
```

We illustrate fixing emissions:

emission_matrix()

Return the matrix whose i-th row specifies the emission probability distribution for the i-th state. More precisely, the i,j entry of the matrix is the probability of the Markov model outputing the j-th symbol when it is in the i-th state.

OUTPUT:

•a Sage matrix with real double precision (RDF) entries.

EXAMPLES:

The returned matrix is mutable, but changing it does not change the transition matrix for the model:

```
sage: E[0,0] = 0; E[0,1] = 1
sage: m.emission_matrix()
```

```
[0.1 0.9]
[0.5 0.5]
```

generate_sequence (length, starting_state=None)

Return a sample of the given length from this HMM.

INPUT:

- •length positive integer
- •starting_state int (or None); if specified then generate a sequence using this model starting with the given state instead of the initial probabilities to determine the starting state.

OUTPUT:

- •an IntList or list of emission symbols
- •IntList of the actual states the model was in when emitting the corresponding symbols

EXAMPLES:

In this example, the emission symbols are not set:

Here the emission symbols are set:

Specify the starting state:

```
sage: set_random_seed(0); a.generate_sequence(5, starting_state=0)
(['up', 'up', 'down', 'down'], [0, 0, 1, 1, 1])
```

log_likelihood (obs, scale=True)

Return the logarithm of the probability that this model produced the given observation sequence. Thus the output is a non-positive number.

INPUT:

- •obs sequence of observations
- •scale boolean (default: True); if True, use rescaling to overoid loss of precision due to the very limited dynamic range of floats. You should leave this as True unless the obs sequence is very small.

```
sage: m = hmm.DiscreteHiddenMarkovModel([[0.4,0.6],[0.1,0.9]], [[0.1,0.9],[0.

→5,0.5]], [.2,.8])
sage: m.log_likelihood([0, 1, 0, 1, 1, 0, 1, 0, 0, 0])
-7.3301308009370825
sage: m.log_likelihood([0, 1, 0, 1, 1, 0, 1, 0, 0, 0], scale=False)
```

```
-7.330130800937082

sage: m.log_likelihood([])
0.0

sage: m = hmm.DiscreteHiddenMarkovModel([[0.4,0.6],[0.1,0.9]], [[0.1,0.9],[0.4,0.6]]), [0.1,0.9]], [[0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]], [0.1,0.9]]
```

Overflow from not using the scale option:

```
sage: m = hmm.DiscreteHiddenMarkovModel([[0.4,0.6],[0.1,0.9]], [[0.1,0.9],[0.

→5,0.5]], [.2,.8])
sage: m.log_likelihood([0,1]*1000, scale=True)
-1433.820666652728
sage: m.log_likelihood([0,1]*1000, scale=False)
-inf
```

viterbi (obs, log scale=True)

Determine "the" hidden sequence of states that is most likely to produce the given sequence seq of observations, along with the probability that this hidden sequence actually produced the observation.

INPUT:

- •seq sequence of emitted ints or symbols
- •log_scale bool (default: True) whether to scale the sequence in order to avoid numerical over-flow.

OUTPUT:

- •list "the" most probable sequence of hidden states, i.e., the Viterbi path.
- •float log of probability that the observed sequence was produced by the Viterbi sequence of states.

EXAMPLES:

We predict the state sequence when the emissions are 3/4 and 'abc'.:

Note that state 0 is common below, despite the model trying hard to switch to state 1:

```
sage: a.viterbi([3/4, 'abc', 'abc'] + [3/4]*10)
([0, 1, 1, 0, 0 ... 0, 0, 0, 0], -25.299405845367794)
```

class sage.stats.hmm.hmm. HiddenMarkovModel

Bases: object

Abstract base class for all Hidden Markov Models.

graph (eps=0.001)

Create a weighted directed graph from the transition matrix, not including any edge with a probability less than eps.

INPUT:

•eps – nonnegative real number

OUTPUT:

•a digraph

EXAMPLES:

initial_probabilities ()

Return the initial probabilities, which as a TimeSeries of length N, where N is the number of states of the Markov model.

EXAMPLES:

```
sage: m = hmm.DiscreteHiddenMarkovModel([[0.4,0.6],[0.1,0.9]], [[0.1,0.9],[0.

→5,0.5]], [.2,.8])
sage: pi = m.initial_probabilities(); pi
[0.2000, 0.8000]
sage: type(pi)
<type 'sage.finance.time_series.TimeSeries'>
```

The returned time series is a copy, so changing it does not change the model.

```
sage: pi[0] = .1; pi[1] = .9 sage: m.initial_probabilities() [0.2000, 0.8000]
```

Some other models:

sample (*length*, *number=None*, *starting_state=None*)

Return number samples from this HMM of given length.

INPUT:

- •length positive integer
- •number (default: None) if given, compute list of this many sample sequences
- •starting_state int (or None); if specified then generate a sequence using this model starting with the given state instead of the initial probabilities to determine the starting state.

OUTPUT:

- •if number is not given, return a single TimeSeries.
- •if number is given, return a list of TimeSeries.

EXAMPLES:

If the emission symbols are set:

Force a starting state:

```
sage: set_random_seed(0); a.sample(10, starting_state=0)
['up', 'up', 'down', 'down', 'down', 'up', 'up', 'up', 'up']
```

transition_matrix ()

Return the state transition matrix.

OUTPUT:

•a Sage matrix with real double precision (RDF) entries.

EXAMPLES:

The returned matrix is mutable, but changing it does not change the transition matrix for the model:

```
sage: T[0,0] = .1; T[0,1] = .9
sage: M.transition_matrix()
[0.7 0.3]
[0.9 0.1]
```

Transition matrices for other types of models:

sage.stats.hmm.hmm.unpickle_discrete_hmm_v0 (A, B, pi, emission_symbols, name)
TESTS:

sage.stats.hmm.hmm.unpickle_discrete_hmm_v1 (A, B, pi, n_out, emission_symbols, emission_symbols_dict)

TESTS:

CONTINUOUS EMISSION HIDDEN MARKOV MODELS

AUTHOR:

• William Stein, 2010-03

```
{\bf class} \; {\tt sage.stats.hmm.chmm.} \; {\tt GaussianHiddenMarkovModel} \\
```

Bases: sage.stats.hmm.hmm.HiddenMarkovModel

GaussianHiddenMarkovModel(A, B, pi)

Gaussian emissions Hidden Markov Model.

INPUT:

- •A matrix; the N x N transition matrix
- •B list of pairs (mu,sigma) that define the distributions
- •pi initial state probabilities
- •normalize -bool (default: True)

EXAMPLES:

We illustrate the primary functions with an example 2-state Gaussian HMM:

We query the defining transition matrix, emission parameters, and initial state probabilities:

```
sage: m.transition_matrix()
[0.1 0.9]
[0.5 0.5]
sage: m.emission_parameters()
[(1.0, 1.0), (-1.0, 1.0)]
sage: m.initial_probabilities()
[0.5000, 0.5000]
```

We obtain a sample sequence with 10 entries in it, and compute the logarithm of the probability of obtaining his sequence, given the model:

We compute the Viterbi path, and probability that the given path of states produced obs:

```
sage: m.viterbi(obs)
([1, 0, 1, 0, 1, 1, 0, 1], -16.67738270170788)
```

We use the Baum-Welch iterative algorithm to find another model for which our observation sequence is more likely:

```
sage: m.baum_welch(obs)
(-10.6103334957397..., 14)
sage: m.log_likelihood(obs)
-10.6103334957397...
```

Notice that running Baum-Welch changed our model:

```
sage: m # rel tol 3e-14
Gaussian Hidden Markov Model with 2 States
Transition matrix:
[    0.4154981366185841    0.584501863381416]
[    0.9999993174253741 6.825746258991804e-07]
Emission parameters:
[(0.4178882427119503, 0.5173109664360919), (-1.5025208631331122, 0.
→5085512836055119)]
Initial probabilities: [0.0000, 1.0000]
```

baum_welch (obs, max_iter=500, log_likelihood_cutoff=0.0001, min_sd=0.01, fix_emissions=False, v=False)

Given an observation sequence obs, improve this HMM using the Baum-Welch algorithm to increase the probability of observing obs.

INPUT:

- •obs a time series of emissions
- •max_iter integer (default: 500) maximum number of Baum-Welch steps to take
- •log_likelihood_cutoff positive float (default: 1e-4); the minimal improvement in likelihood with respect to the last iteration required to continue. Relative value to log likelihood.
- •min_sd positive float (default: 0.01); when reestimating, the standard deviation of emissions is not allowed to be less than min_sd.
- •fix_emissions bool (default: False); if True, do not change emissions when updating

OUTPUT:

•changes the model in places, and returns the log likelihood and number of iterations.

We illustrate bounding the standard deviation below. Note that above we had different emission parameters when the min sd was the default of 0.01:

We watch the log likelihoods of the model converge, step by step:

We illustrate fixing emissions:

emission_parameters ()

Return the parameters that define the normal distributions associated to all of the states.

OUTPUT

•a list B of pairs B[i] = (mu, std), such that the distribution associated to state i is normal with mean mu and standard deviation std.

generate_sequence (length, starting_state=None)

Return a sample of the given length from this HMM.

INPUT:

- •length positive integer
- •starting_state int (or None); if specified then generate a sequence using this model starting with the given state instead of the initial probabilities to determine the starting state.

OUTPUT:

- •an IntList or list of emission symbols
- •TimeSeries of emissions

EXAMPLES:

Example in which the starting state is 0 (see trac ticket #11452):

```
sage: set_random_seed(23); m.generate_sequence(2)
([0.6501, -2.0151], [0, 1])
```

Force a starting state of 1 even though as we saw above it would be 0:

```
sage: set_random_seed(23); m.generate_sequence(2, starting_state=1)
([-3.1491, -1.0244], [1, 1])
```

Verify numerically that the starting state is 0 with probability about 0.1:

```
sage: set_random_seed(0)
sage: v = [m.generate_sequence(1)[1][0] for i in range(10^5)]
sage: 1.0 * v.count(int(0)) / len(v)
0.0998200000000000
```

log_likelihood (obs)

Return the logarithm of a continuous analogue of the probability that this model produced the given observation sequence.

Note that the "continuous analogue of the probability" above can be bigger than 1, hence the logarithm can be positive.

INPUT:

•obs – sequence of observations

OUTPUT:

•float

EXAMPLES:

viterbi (obs)

Determine "the" hidden sequence of states that is most likely to produce the given sequence seq of observations, along with the probability that this hidden sequence actually produced the observation.

INPUT:

•seq – sequence of emitted ints or symbols

OUTPUT:

- •list "the" most probable sequence of hidden states, i.e., the Viterbi path.
- •float log of probability that the observed sequence was produced by the Viterbi sequence of states.

EXAMPLES:

We find the optimal state sequence for a given model:

Another example in which the most likely states change based on the last observation:

class sage.stats.hmm.chmm. GaussianMixtureHiddenMarkovModel

Bases: sage.stats.hmm.chmm.GaussianHiddenMarkovModel

GaussianMixtureHiddenMarkovModel(A, B, pi)

Gaussian mixture Hidden Markov Model.

INPUT:

- •A matrix: the N x N transition matrix
- •B list of mixture definitions for each state. Each state may have a varying number of gaussians with selection probabilities that sum to 1 and encoded as (p,(mu,sigma))
- •pi initial state probabilities
- •normalize -bool (default: True); if given, input is normalized to define valid probability distributions, e.g., the entries of A are made nonnegative and the rows sum to 1, and the probabilities in pi are normalized.

EXAMPLES:

```
sage: A = [[0.5,0.5],[0.5,0.5]]
sage: B = [[(0.9,(0.0,1.0)), (0.1,(1,10000))],[(1,(1,1)), (0,(0,0.1))]]
sage: hmm.GaussianMixtureHiddenMarkovModel(A, B, [1,0])
Gaussian Mixture Hidden Markov Model with 2 States
Transition matrix:
[0.5 0.5]
[0.5 0.5]
Emission parameters:
[0.9*N(0.0,1.0) + 0.1*N(1.0,10000.0), 1.0*N(1.0,1.0) + 0.0*N(0.0,0.1)]
Initial probabilities: [1.0000, 0.0000]
```

TESTS:

If a standard deviation is 0, it is normalized to be slightly bigger than 0.:

```
sage: hmm.GaussianMixtureHiddenMarkovModel([[1]], [[(1,(0,0))]], [1])
Gaussian Mixture Hidden Markov Model with 1 States
Transition matrix:
[1.0]
Emission parameters:
[1.0*N(0.0,1e-08)]
Initial probabilities: [1.0000]
```

We test that number of emission distributions must be the same as the number of states:

```
sage: hmm.GaussianMixtureHiddenMarkovModel([[1]], [], [1])
Traceback (most recent call last):
...
ValueError: number of GaussianMixtures must be the same as number of entries of pi
sage: hmm.GaussianMixtureHiddenMarkovModel([[1]], [[]], [1])
Traceback (most recent call last):
...
ValueError: must specify at least one component of the mixture model
```

We test that the number of initial probabilities must equal the number of states:

baum_welch (obs, max_iter=1000, log_likelihood_cutoff=1e-12, min_sd=0.01, fix_emissions=False) Given an observation sequence obs, improve this HMM using the Baum-Welch algorithm to increase the probability of observing obs.

INPUT:

- •obs a time series of emissions
- •max_iter integer (default: 1000) maximum number of Baum-Welch steps to take
- •log_likelihood_cutoff positive float (default: 1e-12); the minimal improvement in likelihood with respect to the last iteration required to continue. Relative value to log likelihood.
- •min_sd positive float (default: 0.01); when reestimating, the standard deviation of emissions is not allowed to be less than min_sd.

•fix_emissions – bool (default: False); if True, do not change emissions when updating

OUTPUT:

•changes the model in places, and returns the log likelihood and number of iterations.

EXAMPLES:

```
sage: m = hmm.GaussianMixtureHiddenMarkovModel([[.9,.1],[.4,.6]], [[(.
 \rightarrow 4, (0,1)), (.6, (1,0.1))], [(1, (0,1))]], [.7,.3])
sage: set_random_seed(0); v = m.sample(10); v
[0.3576, -0.9365, 0.9449, -0.6957, 1.0217, 0.9644, 0.9987, -0.5950, -1.0219, ]
 →0.6477]
sage: m.log_likelihood(v)
-8.31408655939536...
sage: m.baum_welch(v)
(2.18905068682..., 15)
sage: m.log_likelihood(v)
2.18905068682...
sage: m # rel tol 6e-14
Gaussian Mixture Hidden Markov Model with 2 States
Transition matrix:
              0.8746363339773399
                                                                                       0.125363666022660161
                                                                      1.0 1.451685202290174e-401
Emission parameters:
 [0.500161629343*N (-0.812298726239, 0.173329026744) + 0.499838370657*N (0.499838370657*N (0.4998387) (0.49983870657*N (0.49983870657*N (0.4998387) (0.49983870657*N (0.4998387) (0.49987067*N (0.
 \rightarrow 982433690378, 0.029719932009), \ 1.0*N(0.503260056832, 0.145881515324)]
Initial probabilities: [0.0000, 1.0000]
```

We illustrate bounding the standard deviation below. Note that above we had different emission parameters when the min_sd was the default of 0.01:

We illustrate fixing all emissions:

emission_parameters ()

Returns a list of all the emission distributions.

OUTPUT:

•list of Gaussian mixtures

```
sage: m = hmm.GaussianMixtureHiddenMarkovModel([[.9,.1],[.4,.6]], [[(.
\rightarrow4,(0,1)), (.6,(1,0.1))],[(1,(0,1))]], [.7,.3])
sage: m.emission_parameters()
[0.4*N(0.0,1.0) + 0.6*N(1.0,0.1), 1.0*N(0.0,1.0)]
```

sage.stats.hmm.chmm.unpickle_gaussian_hmm_v0 (A, B, pi, name)

EXAMPLES:

```
sage: m = hmm.GaussianHiddenMarkovModel([[1]], [(0,1)], [1])
sage: sage.stats.hmm.chmm.unpickle_gaussian_hmm_v0(m.transition_matrix(), m.

--emission_parameters(), m.initial_probabilities(), 'test')
Gaussian Hidden Markov Model with 1 States
Transition matrix:
[1.0]
Emission parameters:
[(0.0, 1.0)]
Initial probabilities: [1.0000]
```

sage.stats.hmm.chmm.unpickle_gaussian_hmm_v1 (A, B, pi, prob, n_out)

EXAMPLES:

```
sage: m = hmm.GaussianHiddenMarkovModel([[1]], [(0,1)], [1])
sage: loads(dumps(m)) == m # indirect test
True
```

sage.stats.hmm.chmm.unpickle_gaussian_mixture_hmm_v1 (A, B, pi, mixture)

```
sage: m = hmm.GaussianMixtureHiddenMarkovModel([[1]], [[(.4,(0,1)), (.6,(1,0.
\rightarrow1))]], [1])
sage: loads(dumps(m)) == m # indirect test
True
```

DISTRIBUTIONS USED IN IMPLEMENTING HIDDEN MARKOV MODELS

These distribution classes are designed specifically for HMM's and not for general use in statistics. For example, they have fixed or non-fixed status, which only make sense relative to being used in a hidden Markov model.

AUTHOR:

```
• William Stein, 2010-03
class sage.stats.hmm.distributions. DiscreteDistribution
     Bases: sage.stats.hmm.distributions.Distribution
class sage.stats.hmm.distributions. Distribution
     Bases: object
     A distribution.
     plot (*args, **kwds)
         Return a plot of the probability density function.
         INPUT:
            •args and kwds, passed to the Sage plot function
         OUTPUT:
            •a Graphics object
         EXAMPLES:
         sage: P = hmm.GaussianMixtureDistribution([(.2,-10,.5),(.6,1,1),(.2,20,.5)])
         sage: P.plot(-10,30)
         Graphics object consisting of 1 graphics primitive
     prob(x)
         The probability density function evaluated at x.
         INPUT:
            •x – object
         OUTPUT:
            •float
         EXAMPLES:
         This method must be defined in a derived class:
```

```
sage: import sage.stats.hmm.distributions
sage: sage.stats.hmm.distributions.Distribution().prob(0)
Traceback (most recent call last):
...
NotImplementedError
```

sample (n=None)

Return either a single sample (the default) or n samples from this probability distribution.

INPUT:

•n – None or a positive integer

OUTPUT:

•a single sample if n is 1; otherwise many samples

EXAMPLES:

This method must be defined in a derived class:

```
sage: import sage.stats.hmm.distributions
sage: sage.stats.hmm.distributions.Distribution().sample()
Traceback (most recent call last):
...
NotImplementedError
```

```
{\bf class} \; {\tt sage.stats.hmm.distributions.} \; {\tt GaussianDistribution}
```

Bases: sage.stats.hmm.distributions.Distribution

class sage.stats.hmm.distributions. GaussianMixtureDistribution

Bases: sage.stats.hmm.distributions.Distribution

A probability distribution defined by taking a weighted linear combination of Gaussian distributions.

EXAMPLES:

```
sage: P = hmm.GaussianMixtureDistribution([(.3,1,2),(.7,-1,1)]); P
0.3*N(1.0,2.0) + 0.7*N(-1.0,1.0)
sage: P[0]
(0.3, 1.0, 2.0)
sage: P.is_fixed()
False
sage: P.fix(1)
sage: P.is_fixed(0)
False
sage: P.is_fixed(1)
True
sage: P.unfix(1)
sage: P.is_fixed(1)
False
```

fix (*i=None*)

Set that this GaussianMixtureDistribution (or its ith component) is fixed when using Baum-Welch to update the corresponding HMM.

INPUT:

•i - None (default) or integer; if given, only fix the i-th component

```
sage: P = hmm.GaussianMixtureDistribution([(.2,-10,.5),(.6,1,1),(.2,20,.5)])
sage: P.fix(1); P.is_fixed()
False
sage: P.is_fixed(1)
True
sage: P.fix(); P.is_fixed()
True
```

is fixed (i=None)

Return whether or not this GaussianMixtureDistribution is fixed when using Baum-Welch to update the corresponding HMM.

INPUT:

•i - None (default) or integer; if given, only return whether the i-th component is fixed

EXAMPLES:

```
sage: P = hmm.GaussianMixtureDistribution([(.2,-10,.5),(.6,1,1),(.2,20,.5)])
sage: P.is_fixed()
False
sage: P.is_fixed(0)
False
sage: P.fix(0); P.is_fixed()
False
sage: P.is_fixed(0)
True
sage: P.fix(); P.is_fixed()
True
```

prob(x)

Return the probability of x.

Since this is a continuous distribution, this is defined to be the limit of the p's such that the probability of [x,x+h] is p*h.

INPUT:

 $\bullet x$ – float

OUTPUT:

•float

EXAMPLES:

```
sage: P = hmm.GaussianMixtureDistribution([(.2,-10,.5),(.6,1,1),(.2,20,.5)])
sage: P.prob(.5)
0.21123919605857971
sage: P.prob(-100)
0.0
sage: P.prob(20)
0.1595769121605731
```

$prob_m (x, m)$

Return the probability of x using just the m-th summand.

INPUT:

```
•x – float
```

•m – integer

OUTPUT:

•float

EXAMPLES:

```
sage: P = hmm.GaussianMixtureDistribution([(.2,-10,.5),(.6,1,1),(.2,20,.5)])
sage: P.prob_m(.5, 0)
2.7608117680508...e-97
sage: P.prob_m(.5, 1)
0.21123919605857971
sage: P.prob_m(.5, 2)
0.0
```

sample (n=None)

Return a single sample from this distribution (by default), or if n>1, return a TimeSeries of samples.

INPUT:

•n – integer or None (default: None)

OUTPUT:

•float if n is None (default); otherwise a TimeSeries

EXAMPLES:

```
sage: P = hmm.GaussianMixtureDistribution([(.2,-10,.5),(.6,1,1),(.2,20,.5)])
sage: P.sample()
19.65824361087513
sage: P.sample(1)
[-10.4683]
sage: P.sample(5)
[-0.1688, -10.3479, 1.6812, 20.1083, -9.9801]
sage: P.sample(0)
[]
sage: P.sample(-3)
Traceback (most recent call last):
...
ValueError: n must be nonnegative
```

unfix (*i=None*)

Set that this GaussianMixtureDistribution (or its ith component) is not fixed when using Baum-Welch to update the corresponding HMM.

INPUT:

•i - None (default) or integer; if given, only fix the i-th component

```
sage: P = hmm.GaussianMixtureDistribution([(.2,-10,.5),(.6,1,1),(.2,20,.5)])
sage: P.fix(1); P.is_fixed(1)
True
sage: P.unfix(1); P.is_fixed(1)
False
sage: P.fix(); P.is_fixed()
True
sage: P.unfix(); P.is_fixed()
False
```

```
sage.stats.hmm.distributions. unpickle_gaussian_mixture_distribution_v1 ( c0, c1, param, fixed)
```

Used in unpickling GaussianMixtureDistribution's.

```
sage: P = hmm.GaussianMixtureDistribution([(.2,-10,.5),(.6,1,1),(.2,20,.5)])
sage: loads(dumps(P)) == P  # indirect doctest
True
```



HIDDEN MARKOV MODELS – UTILITY FUNCTIONS

AUTHOR:

• William Stein, 2010-03

```
class sage.stats.hmm.util. HMM_Util
    Bases: object
```

A class used in order to share cdef's methods between different files.

```
initial_probs_to_TimeSeries ( pi, normalize)
```

This function is used internally by the __init__ methods of various Hidden Markov Models.

INPUT:

```
•pi – vector, list, or TimeSeries
```

•normalize – if True, replace negative entries by 0 and rescale to ensure that the sum of the entries in each row is equal to 1. If the sum of the entries in a row is 0, replace them all by 1/N.

OUTPUT:

• a TimeSeries of length N

EXAMPLES:

```
sage: import sage.stats.hmm.util
sage: u = sage.stats.hmm.util.HMM_Util()
sage: u.initial_probs_to_TimeSeries([0.1,0.2,0.9], True)
[0.0833, 0.1667, 0.7500]
sage: u.initial_probs_to_TimeSeries([0.1,0.2,0.9], False)
[0.1000, 0.2000, 0.9000]
```

$normalize_probability_TimeSeries (T, i, j)$

This function is used internally by the Hidden Markov Models code.

Replace entries of T[i:j] in place so that they are all nonnegative and sum to 1. Negative entries are replaced by 0 and T[i:j] is then rescaled to ensure that the sum of the entries in each row is equal to 1. If all entries are 0, replace them by 1/(j-i).

INPUT:

- •T a TimeSeries
- •i nonnegative integer
- •j nonnegative integer

OUTPUT:

•T is modified

EXAMPLES:

```
sage: import sage.stats.hmm.util
sage: T = stats.TimeSeries([.1, .3, .7, .5])
sage: u = sage.stats.hmm.util.HMM_Util()
sage: u.normalize_probability_TimeSeries(T,0,3)
sage: T
[0.0909, 0.2727, 0.6364, 0.5000]
sage: u.normalize_probability_TimeSeries(T,0,4)
sage: T
[0.0606, 0.1818, 0.4242, 0.3333]
sage: abs(T.sum()-1) < le-8  # might not exactly equal 1 due to rounding
True</pre>
```

state_matrix_to_TimeSeries (A, N, normalize)

This function is used internally by the __init__ methods of Hidden Markov Models to make a transition matrix from A.

INPUT:

- •A matrix, list, list of lists, or TimeSeries
- •N number of states
- •normalize if True, replace negative entries by 0 and rescale to ensure that the sum of the entries in each row is equal to 1. If the sum of the entries in a row is 0, replace them all by 1/N.

OUTPUT:

•a TimeSeries

```
sage: import sage.stats.hmm.util
sage: u = sage.stats.hmm.util.HMM_Util()
sage: u.state_matrix_to_TimeSeries([[.1,.7],[3/7,4/7]], 2, True)
[0.1250, 0.8750, 0.4286, 0.5714]
sage: u.state_matrix_to_TimeSeries([[.1,.7],[3/7,4/7]], 2, False)
[0.1000, 0.7000, 0.4286, 0.5714]
```

DISCRETE GAUSSIAN SAMPLERS OVER THE INTEGERS

This class realizes oracles which returns integers proportionally to $\exp(-(x-c)^2/(2\sigma^2))$. All oracles are implemented using rejection sampling. See <code>DiscreteGaussianDistributionIntegerSampler.__init__()</code> for which algorithms are available.

AUTHORS:

• Martin Albrecht (2014-06-28): initial version

EXAMPLE:

We construct a sampler for the distribution $D_{3,c}$ with width $\sigma = 3$ and center c = 0:

We ask for 100000 samples:

```
sage: from six.moves import range
sage: n=1000000; l = [D() for _ in range(n)]
```

These are sampled with a probability proportional to $\exp(-x^2/18)$. More precisely we have to normalise by dividing by the overall probability over all integers. We use the fact that hitting anything more than 6 standard deviations away is very unlikely and compute:

```
sage: norm_factor = sum([exp(-x^2/(2*sigma^2)) for x in range(-6*sigma, sigma*6+1)])
sage: norm_factor
7.519...
```

With this normalisation factor, we can now test if our samples follow the expected distribution:

```
sage: x=0; 1.count(x), ZZ(round(n*exp(-x^2/(2*sigma^2))/norm_factor))
(13355, 13298)
sage: x=4; 1.count(x), ZZ(round(n*exp(-x^2/(2*sigma^2))/norm_factor))
(5479, 5467)
sage: x=-10; 1.count(x), ZZ(round(n*exp(-x^2/(2*sigma^2))/norm_factor))
(53, 51)
```

We construct an instance with a larger width:

ask for 100000 samples:

```
sage: from six.moves import range
sage: n=100000; l = [D() for _ in range(n)] # long time
```

and check if the proportions fit:

We construct a sampler with c%1! = 0:

REFERENCES:

A Discrete Gaussian Sampler using rejection sampling.

```
__init__ ( sigma, c=0, tau=6, algorithm=None, precision='mp')

Construct a new sampler for a discrete Gaussian distribution.
```

INPUT:

- •sigma samples x are accepted with probability proportional to $\exp(-(x-c)/(2\sigma))$
- •c the mean of the distribution. The value of c does not have to be an integer. However, some algorithms only support integer-valued c (default: 0)
- •tau samples outside the range $(\lfloor c \rceil \lceil \sigma \tau \rceil, ..., \lfloor c \rceil + \lceil \sigma \tau \rceil)$ are considered to have probability zero. This bound applies to algorithms which sample from the uniform distribution (default: 6)

```
•algorithm - see list below (default: "uniform+table" for \sigma t bounded by DiscreteGaussianDistributionIntegerSampler.table_cutoff and "uniform+online" for bigger \sigma \tau)
```

•precision - either "mp" for multi-precision where the actual precision used is taken from sigma or "dp" for double precision. In the latter case results are not reproducible. (default: "mp")

ALGORITHMS:

- •"uniform+table" classical rejection sampling, sampling from the uniform distribution and accepted with probability proportional to $\exp(-(x-c)/(2\sigma))$ where $\exp(-(x-c)/(2\sigma))$ is precomputed and stored in a table. Any real-valued c is supported.
- •"uniform+logtable" samples are drawn from a uniform distribution and accepted with probability proportional to $\exp(-(x-c)/(2\sigma))$ where $\exp(-(x-c)/(2\sigma))$ is computed using logarith-

mically many calls to Bernoulli distributions. See [DDLL13] for details. Only integer-valued c are supported.

- •"uniform+online" samples are drawn from a uniform distribution and accepted with probability proportional to $\exp(-(x-c)/(2\sigma))$ where $\exp(-(x-c)/(2\sigma))$ is computed in each invocation. Typically this is very slow. See [DDLL13] for details. Any real-valued c is accepted.
- •"sigma2+logtable" samples are drawn from an easily samplable distribution with $\sigma=k\cdot\sigma_2$ with $\sigma_2=\sqrt{1/(2\log 2)}$ and accepted with probability proportional to $\exp(-(x-c)/(2\sigma))$ where $\exp(-(x-c)/(2\sigma))$ is computed using logarithmically many calls to Bernoulli distributions (but no calls to exp). See [DDLL13] for details. Note that this sampler adjusts σ to match $k\cdot\sigma_2$ for some integer k. Only integer-valued c are supported.

EXAMPLES:

Note that "sigma2+logtable" adjusts σ :

TESTS:

We are testing invalid inputs:

```
sage: from sage.stats.distributions.discrete gaussian integer import.
→ DiscreteGaussianDistributionIntegerSampler
sage: DiscreteGaussianDistributionIntegerSampler(-3.0)
Traceback (most recent call last):
ValueError: sigma must be > 0.0 but got -3.000000
sage: DiscreteGaussianDistributionIntegerSampler(3.0, tau=-1)
Traceback (most recent call last):
. . .
ValueError: tau must be >= 1 but got -1
sage: DiscreteGaussianDistributionIntegerSampler(3.0, tau=2, algorithm=
→"superfastalgorithmyouneverheardof")
Traceback (most recent call last):
ValueError: Algorithm 'superfastalgorithmyouneverheardof' not supported by ...
⇒class 'DiscreteGaussianDistributionIntegerSampler'
sage: DiscreteGaussianDistributionIntegerSampler(3.0, c=1.5, algorithm=
→"sigma2+logtable")
Traceback (most recent call last):
```

```
...

ValueError: algorithm 'uniform+logtable' requires c%1 == 0
```

We are testing correctness for multi-precision:

```
sage: from sage.stats.distributions.discrete gaussian integer import.
→DiscreteGaussianDistributionIntegerSampler
sage: D = DiscreteGaussianDistributionIntegerSampler(1.0, c=0, tau=2)
sage: from six.moves import range
sage: 1 = [D() \text{ for } \_ \text{ in } range(2^16)]
sage: min(1) == 0-2*1.0, max(1) == 0+2*1.0, abs(mean(1)) < 0.01
(True, True, True)
sage: from sage.stats.distributions.discrete_gaussian_integer import,
→DiscreteGaussianDistributionIntegerSampler
sage: D = DiscreteGaussianDistributionIntegerSampler(1.0, c=2.5, tau=2)
sage: from six.moves import range
sage: 1 = [D() \text{ for } \_ \text{ in } range(2^18)]
sage: min(1) == 2-2 \times 1.0, max(1) == 2+2 \times 1.0, mean(1).n()
(True, True, 2.45...)
sage: from sage.stats.distributions.discrete_gaussian_integer import_
→ DiscreteGaussianDistributionIntegerSampler
sage: D = DiscreteGaussianDistributionIntegerSampler(1.0, c=2.5, tau=6)
sage: from six.moves import range
sage: 1 = [D() \text{ for } \_ \text{ in } range(2^18)]
sage: min(1), max(1), abs(mean(1)-2.5) < 0.01
(-2, 7, True)
```

We are testing correctness for double precision:

```
sage: from sage.stats.distributions.discrete gaussian integer import...
{\bf \hookrightarrow} {\tt DiscreteGaussianDistributionIntegerSampler}
sage: D = DiscreteGaussianDistributionIntegerSampler(1.0, c=0, tau=2,...
→precision="dp")
sage: from six.moves import range
sage: 1 = [D() \text{ for } \_ \text{ in } range(2^16)]
sage: min(1) == 0-2*1.0, max(1) == 0+2*1.0, abs(mean(1)) < 0.05
(True, True, True)
sage: from sage.stats.distributions.discrete_gaussian_integer import.
→DiscreteGaussianDistributionIntegerSampler
sage: D = DiscreteGaussianDistributionIntegerSampler(1.0, c=2.5, tau=2,...
→precision="dp")
sage: from six.moves import range
sage: 1 = [D() \text{ for } \_ \text{ in } range(2^18)]
sage: min(1) == 2-2 \times 1.0, max(1) == 2+2 \times 1.0, mean(1).n()
(True, True, 2.4...)
sage: from sage.stats.distributions.discrete_gaussian_integer import__
→ DiscreteGaussianDistributionIntegerSampler
sage: D = DiscreteGaussianDistributionIntegerSampler(1.0, c=2.5, tau=6,...
→precision="dp")
sage: from six.moves import range
sage: 1 = [D() \text{ for } \_ \text{ in } range(2^18)]
sage: min(1) <=-1, max(1) >= 6, abs(mean(1) -2.5) < 0.1
(True, True, True)
sage: tuple(1.count(i) for i in range(-2,8)) # output random
```

```
(7, 242, 4519, 34120, 92714, 91700, 33925, 4666, 246, 5)
```

We plot a histogram:

These generators cache random bits for performance reasons. Hence, resetting the seed of the PRNG might not have the expected outcome. You can flush this cache with _flush_cache():

__call___()

Return a new sample.

EXAMPLES:

TESTS:

algorithm

c

sigma

tau

CHAPTER

EIGHT

DISCRETE GAUSSIAN SAMPLERS FOR $\mathbb{Z}[X]$

This class realizes oracles which returns polynomials in $\mathbf{Z}[x]$ where each coefficient is sampled independently with a probability proportional to $\exp(-(x-c)/(2\sigma))$.

AUTHORS:

• Martin Albrecht, Robert Fitzpatrick, Daniel Cabracas, Florian Göpfert, Michael Schneider: initial version

EXAMPLE:

class sage.stats.distributions.discrete_gaussian_polynomial. DiscreteGaussianDistributionPoly

Bases: sage.structure.sage_object.SageObject

Discrete Gaussian sampler for polynomials.

EXAMPLE:

```
__init__ ( P, n, sigma)
```

Construct a sampler for univariate polynomials of degree n-1 where coefficients are drawn independently with standard deviation sigma.

INPUT:

- •P a univariate polynomial ring over the Integers
- •n number of coefficients to be sampled

•sigma - coefficients x are accepted with probability proportional to $\exp(-x/(2\sigma))$. If an object of type $sage.stats.distributions.discrete_gaussian_integer.DiscreteGaussianDistribut$ is passed, then this sampler is used to sample coefficients.

EXAMPLE:

```
sage: from sage.stats.distributions.discrete_gaussian_polynomial import_

→DiscreteGaussianDistributionPolynomialSampler
sage: DiscreteGaussianDistributionPolynomialSampler(ZZ['x'], 8, 3.0)()
3*x^7 + 3*x^6 - 3*x^5 - x^4 - 5*x^2 + 3
sage: gs = DiscreteGaussianDistributionPolynomialSampler(ZZ['x'], 8, 3.0)
sage: [gs() for _ in range(3)]
[4*x^7 + 4*x^6 - 4*x^5 + 2*x^4 + x^3 - 4*x + 7, -5*x^6 + 4*x^5 - 3*x^3 + 4*x^6 + 2*x^7 + 2*x^6 + 2*x^5 - x^4 - 2*x^2 + 3*x + 1]
```

__call__ ()

Return a new sample.

CHAPTER

NINE

DISCRETE GAUSSIAN SAMPLERS OVER LATTICES

This file implements oracles which return samples from a lattice following a discrete Gaussian distribution. That is, if σ is big enough relative to the provided basis, then vectors are returned with a probability proportional to $\exp(-|x-c|_2^2/(2\sigma^2))$. More precisely lattice vectors in $x \in \Lambda$ are returned with probability:

$$\exp(-|x-c|_2^2/(2\sigma))/(\sum_{x\in\Lambda}\exp(-|x|_2^2/(2\sigma)))$$

AUTHORS:

• Martin Albrecht (2014-06-28): initial version

EXAMPLES:

```
sage: from sage.stats.distributions.discrete_gaussian_lattice import_
    →DiscreteGaussianDistributionLatticeSampler
sage: D = DiscreteGaussianDistributionLatticeSampler(ZZ^10, 3.0)
sage: D(), D(), D()
((3, 0, -5, 0, -1, -3, 3, 3, -7, 2), (4, 0, 1, -2, -4, -4, 4, 0, 1, -4), (-3, 0, 4, 4, 4, 5, 0, 1, 3, 2, 0, -1))
```

 ${\bf class} \; {\tt sage.stats.distributions.discrete_gaussian_lattice.} \; {\tt DiscreteGaussianDistributionLattice} \; {\tt class} \; {\tt sage.stats.distributions.discrete_gaussian_lattice}. \; {\tt DiscreteGaussianDistributionLattice} \; {\tt class} \; {\tt sage.stats.distributions.discrete_gaussian_lattice}. \; {\tt DiscreteGaussianDistributionLattice} \; {\tt class} \; {\tt c$

Bases: sage.structure.sage_object.SageObject

GPV sampler for Discrete Gaussians over Lattices.

```
[0 0 0 0 0 0 0 1 0]
[0 0 0 0 0 0 0 1]
```

We plot a histogram:

REFERENCES:

```
__init__ ( B, sigma=1, c=None, precision=None)

Construct a discrete Gaussian sampler over the lattice \Lambda(B) with parameter sigma and center c.
```

INPUT:

- $\bullet B a$ basis for the lattice, one of the following:
 - -an integer matrix,
 - -an object with a matrix () method, e.g. ZZ^n, or
 - -an object where matrix (B) succeeds, e.g. a list of vectors.
- •sigma Gaussian parameter $\sigma > 0$.
- •c center c, any vector in \mathbb{Z}^n is supported, but $c \in \Lambda(B)$ is faster.
- •precision bit precision > 53.

```
sage: from sage.stats.distributions.discrete_gaussian_lattice import.
→ Discrete Gaussian Distribution Lattice Sampler
sage: n = 2; sigma = 3.0; m = 5000
sage: D = DiscreteGaussianDistributionLatticeSampler(ZZ^n, sigma)
sage: f = D.f
sage: c = D._normalisation_factor_zz(); c
56.2162803067524
sage: l = [D() for _ in range(m)]
sage: v = vector(ZZ, n, (-3, -3))
sage: 1.count(v), ZZ(round(m*f(v)/c))
(39, 33)
sage: target = vector(ZZ, n, (0,0))
sage: 1.count(target), ZZ(round(m*f(target)/c))
(116, 89)
sage: from sage.stats.distributions.discrete_gaussian_lattice import,
→ Discrete Gaussian Distribution Lattice Sampler
sage: qf = QuadraticForm(matrix(3, [2, 1, 1, 1, 2, 1, 1, 1, 2]))
sage: D = DiscreteGaussianDistributionLatticeSampler(qf, 3.0); D
Discrete Gaussian sampler with \sigma = 3.000000, c=(0, 0, 0) over lattice with
-basis
```

```
[2 1 1]
[1 2 1]
[1 1 2]
sage: D()
(0, 1, -1)
```

```
__call___()
```

Return a new sample.

EXAMPLE:

```
sage: from sage.stats.distributions.discrete_gaussian_lattice import_

DiscreteGaussianDistributionLatticeSampler
sage: D = DiscreteGaussianDistributionLatticeSampler(ZZ^3, 3.0, c=(1,0,0))
sage: L = [D() for _ in range(2^12)]
sage: abs(mean(L).n() - D.c)
0.08303258...

sage: D = DiscreteGaussianDistributionLatticeSampler(ZZ^3, 3.0, c=(1/2,0,0))
sage: L = [D() for _ in range(2^12)] # long time
sage: mean(L).n() - D.c # long time
(0.0607910156250000, -0.128417968750000, 0.0239257812500000)
```

C

Center c.

Samples from this sampler will be centered at c.

EXAMPLE:

```
sage: from sage.stats.distributions.discrete_gaussian_lattice import_
    →DiscreteGaussianDistributionLatticeSampler
sage: D = DiscreteGaussianDistributionLatticeSampler(ZZ^3, 3.0, c=(1,0,0)); D
Discrete Gaussian sampler with σ = 3.000000, c=(1, 0, 0) over lattice with_
    →basis

[1 0 0]
[0 1 0]
[0 0 1]

sage: D.c
(1, 0, 0)
```

static compute_precision (precision, sigma)

Compute precision to use.

INPUT:

```
•precision - an integer > 53 nor None.
```

•sigma - if precision is None then the precision of sigma is used.

```
sage: DiscreteGaussianDistributionLatticeSampler.compute_precision(100, 3)

100

sage: DiscreteGaussianDistributionLatticeSampler.compute_precision(None, → RR(3))

53

sage: DiscreteGaussianDistributionLatticeSampler.compute_precision(None, → RealField(200)(3))

200

sage: DiscreteGaussianDistributionLatticeSampler.compute_precision(None, 3)

53
```

sigma

Gaussian parameter σ .

Samples from this sampler will have expected norm $\sqrt{n}\sigma$ where n is the dimension of the lattice.

CHAPTER

TEN

T-TEST USING R

TESTS:

CHAPTER

ELEVEN

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