Manual for the GSR/RST learner

last edited: 23 August, 2022

highlighted things are statements that may/will become untrue as the program fills out

flagged things may change later, and need to be double-checked

TO DO's are things that need to be implemented. Some are essential but most are nonessential, or extensions of functionality

The **Features** class:

to use the **Features** class, you will first need to create a features text document. It should be tab-delimited, and formatted as follows:

|  |  |  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- | --- | --- |
| segment names | Feature names | | | | | | |
|  | consonantal | sonorant | continuant | high | low | voice | nasal |
| t | 1 | 0 | 0 | x | x | 0 | 0 |
| n | 1 | 1 | 0 | x | x | 1 | 1 |
| a | 0 | 1 | 1 | 0 | 1 | 1 | 0 |
| u | 0 | 1 | 1 | 1 | 0 | 1 | 0 |
| ... | ... | | | | | | |

Feature values can be any character at all. Here, we are using 1's for [+ feature] and 0's for [- feature]. You can also use '+' and '-', or real number values. Features will be considered a match only if they **match exactly**. TO DO: 'x' values are special and will be interpreted as a segment is unspecified for that feature. (Treated as a wildcard in some cases, and skipped in others)

Provided sample files:

features.xlsx

features.txt

These provide features for all the phonemes of English. Column A of the spreadsheet is IPA, while column B is single-character Arpabet (<https://en.wikipedia.org/wiki/ARPABET>). The text file was made by copying columns B-AD of the spreadsheet.

**Initializing the Features() class:**

You must use a features file to initialize an object of the Features class.

import learner

f = learner.Features('features.txt')

This class now has attributes:

featureNames – a list containing the feature names from the first row of the text file

featureValues – a dictionary with segment names as keys, and lists of feature values as values

use:

f.featureNames

f.featureValues

to examine them.

You can use this object to convert back and forth between strings of segments and feature dictionaries.

Features.stringToF(string)

This function takes a simple string, and interprets each character as a separate segment. It returns two items: a dictionary containing the segments and their feature values, and a list containing the order of the segments.

added seglabels argument

so, running

f.stringToF('ta')

should return this:

﻿({'seg1': [('0', 'syllabic'), ('0', 'stress'),

('0', 'long'), ('1', 'consonantal'), ('0', 'sonorant'), ('0', 'continuous'), ('0', 'delayedRelease'), ('0', 'approximant'), ('0', 'tap'), ('0', 'trill'), ('0', 'nasal'), ('0', 'voice'), ('0', 'spreadGlottis'), ('0', 'constrictedGlottis'), ('0', 'labial'), ('0', 'round'), ('0', 'labioDental'), ('1', 'coronal'), ('1', 'anterior'), ('0', 'distributed'), ('0', 'strident'), ('0', 'lateral'), ('0', 'dorsal'), ('x', 'high'), ('x', 'low'), ('x', 'front'), ('x', 'back'), ('x', 'tense')],

'seg2': [('1', 'syllabic'), ('0', 'stress'), ('0', 'long'), ('0', 'consonantal'), ('1', 'sonorant'), ('1', 'continuous'), ('x', 'delayedRelease'), ('1', 'approximant'), ('0', 'tap'), ('0', 'trill'), ('0', 'nasal'), ('1', 'voice'), ('0', 'spreadGlottis'), ('0', 'constrictedGlottis'), ('0', 'labial'), ('0', 'round'), ('0', 'labioDental'), ('0', 'coronal'), ('x', 'anterior'), ('x', 'distributed'), ('x', 'strident'), ('0', 'lateral'), ('1', 'dorsal'), ('0', 'high'), ('1', 'low'), ('0', 'front'), ('1', 'back'), ('x', 'tense')]},

['seg1', 'seg2'])

Converting a feature set to a string:

Features.featureToS(segs,order)

This function does the reverse. Given a dictionary object, and a list defining the order, it will return a string of the segments in the Features object. To run it, you first need to set up a dictionary for the 'segs' object. Here's an example:

segs = ﻿{'s1':[('1','lateral')],

's2':[('1','high'),('1','back'),('1','tense'),('1','syllabic'),('0','long')]}

f.stringToF(segs,['s1','s2'])

Should return:

'lu'

If you give the function a 'segs' object where one of the segments is compatible with more than one segment in the Features object, it will ask you to manually pick the one you want.

If you give it a set of features that are incompatible with any entry in the Features object, then it will ask you to make up a character, and it will then add that character to it's featureNames object, with featureValues being set to the ones you gave. It will fill in 'x' for all the features not mentioned in your dictionary.

**Note:** You must avoid specifying features that are completely complementary: For example, if I specify every vowel as +back -front OR +front -back, then morphing back-to-front will not work well. Instead, if they are completely predictable from each other, choose one.

**To Do:**

- Check that both functions work for a variety of input types

- Create the 'wildcard' functionality so that 'x' is ignored in Features.featureToS()

- In Features.featureToS(), add a check to make sure that the character the user gives is not already in the object

- Create error message for when the user gives a feature that's not part of the object in Features.featureToS(segs,order)

﻿# Tableau

Call print() on a `Tableau` object to see a printout to the console

\*\*tag\*\* The human-readable name of the tableau, generally indicating what the input to the tableau was

\*\*prob\*\* The tableau probability, corresponding to how likely the learner is to use this tableau for learning. Should correspond to token frequency of the input. Can be set using the 'tab.prob' column in an input file.

\*\*candidates\*\* A list of candidate objects that make up the tableau

\*\*surfaceCands\*\* A list of unique surface forms (strings), which can be set using the 'surface' column in an input file. If the tableau has hidden structure (if more than one candidate corresponds to the same surfaceCand), this list will be shorter than the list of candidates.

\*\*obsProbsList\*\* A list of the observed probabilities for surface candidates in `surfaceCands`. Length corresponds to the length of `surfaceCands`.

\*\*predProbsList\*\* A list of each candidate's predicted probability based on the last time `getPredWinner()` was called.

\*\*HList\*\* A list of each candidate's Harmony score, based on the last time `getPredWinner()` was called.

\*\*winner\*\*

\*\*constraintNames\*\*

\*\*hiddenStructure\*\*

\*\*probDenom\*\*

`addCandidate(cand)` Adds `candidate` `cand` to the tableau

`rect()` Checks the tableau for candidate violation length and violation sign. Returns 1 if the tableau is well-formed, 0 otherwise

`getPredWinner(w,theory=MaxEnt)` Returns a single candidate object that wins under `theory`, with constraint weights `w`. If multiple candidates win, one is sampled. In MaxEnt, sampling is done according to the predicted probability distribution. In HG or Noisy HG, sampling is done from a uniform distribution. Options for specifying `theory` are 'MaxEnt', 'HG', and 'NoisyHG'.

`getObsCandidate(w,theory='MaxEnt')` Returns a single candidate object that is the 'observed' candidate for learning. If more than one candidate has observed probability, then an observed candidate is sampled from the observed distribution. If the tableau has hidden structure, then an observed candidate is calculated using Expected Interpretive Parsing (EIP, Jaroscz, 2013).

`compareObsPred(w,theory='MaxEnt')` Generates an observed candidate and a predicted candidate, and compares their surface forms. Returns a tuple containing the error value, the observed candidate, and the predicted candidate, in that order. The error value is 1 if there was a mismatch between the two surface forms, and 0 otherwise.

**Candidate:**

Simply holds all the info for one candidate. Must be initialized with:

c: a string representation of the candidate. Something that can interface correctly with a Features object you are using, or it can be human-readable any way you want if you're not using a Features object.

violations: a list of violations for the candidate. Must be numbers. The Tableau.rect() function will yell at you though if they are negative, or all zero, or something.

observedProb: the observed probability that this candidate surfaces in its tableau

surfaceForm: the surface form if using hidden structure. If you're not using hidden structure, this does not need to be initialized, it will be set to the same string as c.

After initializing, the candidate has these 4 attributes, plus:

harmony

predictedProb

^ these are both set to zero initially, and filled out during tableau evaluation.

**richCand()**

Inherits from class **candidate**, which means it has all of **candidate**'s attributes and methods/functions. It includes additional information about segment features, ordering, and activity levels. Can interface better with a **Features** object.

Initialize with:

c, violations, observedProb .... (not done)

The **lexeme()** class:

This class is meant to hold a single lexical entry. You can initialize it with a 'tag' and a list of segments, or with a simple string:

l = learner.lexeme('ami', ['t','z','n','a','m','i'])

Here, 'ami' is just a human-readable label, attached to the lexeme but ignored by the program.

l = learner.lexeme('ami')

Here, the class will immediately create a segment list using the string 'ami': ['a', 'm', 'i']

(one entry per segment)

You may also specify a 'kind', such as 'root', 'prefix', 'suffix', or something else. Currently this has no consequences however.

l = learner.lexeme('ami', ['t','z','n','a','m','i'], kind='N')

Once initialized, **lexeme()** has the following attributes:

tag – A human-readable label for the lexeme, not used by the program

segmentList – a list of segments, either given or generated from characters in the tag

segLabels – a list of unique segment names, generated from segmentList.

If segmentList is ['p','a','t','a','k','a','t','a'], segLabels will be ['p', 'a','t','a3','k','a5','t6','a7']. Numbers are added to second, third, etc. copies of segments based on their position in the original list

activitys – (note the unusual spelling) a list of float value activity levels for segmentList Set

to all 1 at initialization

linearSegOrder – a list of integers specifying the linear order of the segments in segmentList

* Starts at 1
* By default, specifies the same order that segments appear in the list: if segmentList is ['a','m','i'], linearSegOrder will be [1,2,3]
* Can be used to specify multiple 'options' for a position: ['t','z','n','a','m','i'] -> [1,1,1,2,3,4]
* Can in principle be used to express epenthesis, deletion, and metathesis, though unclear what use that would be

kind – A label for humans, specifying 'root', 'suffix', etc. Not used by the rest of the program

freq – the frequency of the lexeme, initialized to 0. Intended to increase during learning.

PFCs – A placeholder to contain a list of PFC objects. Set to None on initialization

Use print() to view a handy readout of the lexeme's contents

print(l)

**TO DO**: fix print function to print prettily when there are decimals in the activity values

A set of example lexemes is provided:

exlex\_joli()

exlex\_petit() - partial activation on the last segment

exlex\_ami() - competing segments in the first position, with partial activation

exlex\_hero()

**lexeme.toRichCand(featureSet)**

This function is used in tableau creation, when candidates are not given by the user. It will return a richCand() object representing the 'faithful' candidate for the lexeme. If there are multiple equally 'faithful' candidates possible, it will return a list of all of them.

The argument featureSet must be a Features() object in which all the segments in the lexeme's segmentList are represented.

The **constraint()** class:

To use this class, you must first define a function to assign violations to a candidate. You can define constraints a number of ways, from simple lambda functions to more complex functions.

All constraints must be formulated in such a way that they take as input some subset of these four classes:

string – a string of characters, like ‘pataka’ to which the function will directly refer. Example:

lambda string: re.search([aeiou][aeiou], string) <- one way to define \*Hiatus

def hiatus(string):

return bool(re.search([aeiou][aeiou], string)

richSR – a richCand() object which is the ‘target’ candidate being evaluated

stringFaith – a string of characters which can be compared to string to evaluate a correspondence constraint

richFaith – a richCand() object representing the ‘faithful’ candidate, to which richSR can be compared to evaluate a faithfulness constraint.

Cost of transposition??

can't just be '1'

must be less than deletion plus epenthesis

useListed\_input

**Dealing with Hidden Structure:**

You can use any theory with hidden structure. UR-constraints, GSR’s, and UseListed (in the ‘hidden structure’ version) require hidden structure to operate. This learner follows the principles of Jarosz (2013)’s Expectation-Maximization algorithm to learn hidden structure.

Tableau for ‘cat’ (Based on Pater et al, 2012) correct: [moda]

|  |  |  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- | --- | --- |
|  |  | \*VTV | Ident-Voice | cat --> /mot/ | cat --> /mod/ | H | p |
|  |  | 1 | 2 | 3 | 1 |  |  |
| UR1 | a. /mot/+/a/ → mota | 1 |  |  | 1 | -2 | 0.57 |
| b. /mot/+/a/ → moda |  | 1 |  | 1 | -3 | 0.21 |
| UR2 | c. /mod/+/a/ → mota | 1 | 1 | 1 |  | -6 | 0.01 |
| d. /mod/+/a/ → moda |  |  | 1 |  | -3 | 0.21 |

candidate sampled from p distribution:

say we sampled candidate a.

mota ≠ moda --> update

candidate a. is the Predicted candidate

To get the Observed candidate, we create a mini-tableau with only the candidates that match the observed form (moda). Re-calculate p, and sample as before:

|  |  |  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- | --- | --- |
|  |  | \*VTV | Ident-Voice | cat --> /mot/ | cat --> /mod/ | H | p |
|  |  | 1 | 2 | 3 | 1 |  |  |
| UR1 | b. /mot/+/a/ → moda |  | 1 |  | 1 | -3 | 0.5 |
| UR2 | d. /mod/+/a/ → moda |  |  | 1 |  | -3 | 0.5 |

so there's a 50-50 chance between the two. let's say we sampled b.

Now the update vector looks like this:

|  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- |
|  |  | \*VTV | Ident-Voice | cat --> /mot/ | cat --> /mod/ |
|  | *weights* | 1 | 2 | 3 | 1 |
| predicted | a. /mot/+/a/ → mota | 1 |  |  | 1 |
| observed | b. /mot/+/a/ → moda |  | 1 |  | 1 |
|  | *new weights* | 1.01 | 1.99 | - | - |

because the predicted and observed forms come from the same UR, neither UR constraint updates.

If candidate d. had been sampled as the observed form, we would update differently:

|  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- |
|  |  | \*VTV | Ident-Voice | cat --> /mot/ | cat --> /mod/ |
|  | *weights* | 1 | 2 | 3 | 1 |
| predicted | a. /mot/+/a/ → mota | 1 |  |  | 1 |
| observed | d. /mod/+/a/ → moda |  |  | 1 |  |
|  | *new weights* | 1.01 | - | 2.99 | 1.01 |

GSRs: Let's say that the activity on ta=0.7, and on tb=0.2

|  |  |  |  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- | --- | --- | --- |
|  |  | Max | Dep | Uniformity | NoCoda | \*Hiatus | H | p |
|  | /petit/a + /{t,z,n}ami/b | 2 | 2 | 3 | 5 | 5 |  |  |
| UR1 | a. petiami |  |  |  |  | 1 | -5 | 0.57 |
| b. petitaami | -0.7 | 0.8 |  |  |  | -1 | 0.21 |
| UR2 | c. petitbami | -0.2 | 0.3 |  |  |  | -1 | 0.01 |
| d. petita,bami |  | 0.1 | 1 |  |  | - | 0.21 |
|  | e. petitatbami |  |  |  | 1 |  | -5 |  |

**UR-Constraints**

* A lot like UseListed

|  |  |  |  |  |  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- |
|  |  | Assmilate | \*NC | \*nCoda |  |  |  |  | **H** | **P** |
|  |  | 6 | 3 | 3 | 5 | 10 | 10 | 10 |  |  |
|  | paŋ,poɁok |  |  |  |  |  |  |  |  |  |
| composed | /paŋ+poɁok/ pa**ŋp**oɁok | 1 | 1 | 1 | 0 | 0 | 0 | 0 | -7 |  |
| /paŋ+poɁok/ pa**mp**oɁok | 0 | 1 | 1 | 1 | 0 | 0 | 0 | -3 |  |
| /paŋ+poɁok/ pa**m**oɁok | 0 | 0 | 0 | 0 | 1 | 0 | 0 | -2 |  |
| listed | /pampoɁok/ pa**ŋp**oɁok | 1 | 1 | 1 | 0 | 0 | 1 | 0 | -11 |  |
| /pampoɁok/ pa**mp**oɁok | 0 | 1 | 1 | 0 | 0 | 0 | 0 | -2 |  |
| /pampoɁok/ pa**m**oɁok | 0 |  | 0 | 0 | 0 | 0 | 1 | -4 |  |

**Gradient Symbolic Representations**

* Learn activation values for URs

**Representational Strength theory**

* Learn PFC’s

**Lexically Indexed Constraints:**

Storage:

1. A new list, Grammar.lexCs contains lists of weights for each constraint, always starting at 0

|  |  |  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- | --- | --- |
|  | ℂ1 | ℂ2 | ℂ3 | ℂ4 | ℂ5 | ℂ6 |  |
| *0* | 0 | 0 | 0 | 0 | 0 | 0 | placeholder for general weights, stored in Grammar.w |
| *1* | 1 | 3.5 | 7 |  | 8.2 | 0.1 | weight of first lexically indexed version |
| *2* | 10 | 2 |  |  | 3 |  | weight of second lexically indexed version |
| *3* | 12 |  |  |  |  |  | etc. |
| *4* | 0.3 |  |  |  |  |  |  |

1. Each lexical item has lexeme.lexCindexes, which is a list of indices into those weight lists. It will be 0 if the lexeme has no indexed version of that constraint

paka.lexCindexes: [0, 0, 1, 0, 2, 0]  lexical item ‘paka’ uses weight 7 for ℂ3 and 3 for ℂ5

tapa.lexCindexes: [0, 2, 1, 0, 0, 0]  lexical item ‘tapa’ uses weight 2 for ℂ2 and 7 for ℂ3

 ‘paka’ and ‘tapa’ are both indexed to the same copy of ℂ3!

When its weight changes, it changes for both lexical items

akta.lexCindexes: [0, 2, 0, 0, 2, 0]  lexical item ‘akta’ uses weight 2 for ℂ2 and 3 for ℂ5

 ‘paka’ and ‘akta’ are both indexed to the same copy of ℂ5!

None of these use an indexed version of ℂ6 and an indexed version of ℂ4 does not exist.

In EVAL, the indexed weight REPLACES the general grammatical weight of the constraint

But note: the general weight gets updated on error anyway, since they should reflect the behaviour of the whole lexicon (more or less)

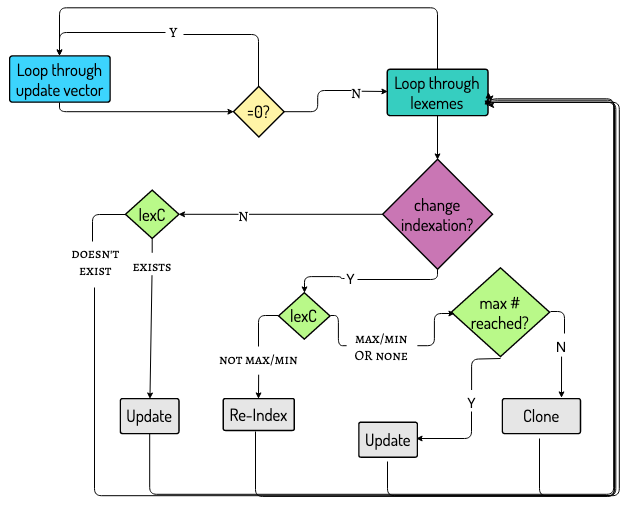
If input lexemes are indexed to different copies of the same constraint, the stricter one (with the higher weight) is chosen

*To edit, check out lines XXX*

Updating lexically-indexed constraints on error

Update vector is predictedCandidate.violations - observedCandidate.violations

[0, -1, 2, 0, 0, -1]  example where constraints 2 and 6 are predicted-preferring, and constraint 3 is observed-preferring. Constraints 1, 4, 5 don’t care and shouldn’t be cloned



random

Regular self.w becomes a vector of ‘baseline’ weights

**Grammar.makeTableau()**

This performs many functions

*Creating UseListed tableaux*

UseListed tableaux come in two flavors. We’ll call them ‘hidden structure’ and ‘probabilistic’

This is from the useListed\_input file, which is a simplified version of the table in Zuraw 2010, page 421. For this one, the assimilated, but unsubstituted form is the correct output: [pa**mp**oɁok].

**Hidden Structure**: Makes a tableau that chooses between the listed form and the surface form. The listed form must exist in the training lexicon, which will happen when the form has previously occurred in training, and an error was made on it.

Additionally, this parameter setting will cause a “UseListed” constraint to be added to the grammar.

|  |  |  |  |  |  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- |
|  |  | Assmilate | \*NC | \*nCoda | Ident-Placecomposed  Maxcomposed  Ident-Placelisted  Maxlisted |  |  |  | **H** | **P** |
|  |  | 6 | 3 | 3 | 5 | 10 | 10 | 10 |  |  |
|  | paŋ,poɁok |  |  |  |  |  |  |  |  |  |
| composed | /paŋ+poɁok/ pa**ŋp**oɁok | 1 | 1 | 1 | 0 | 0 | 0 | 0 | -7 |  |
| /paŋ+poɁok/ pa**mp**oɁok | 0 | 1 | 1 | 1 | 0 | 0 | 0 | -3 |  |
| /paŋ+poɁok/ pa**m**oɁok | 0 | 0 | 0 | 0 | 1 | 0 | 0 | -2 |  |
| listed | /pampoɁok/ pa**ŋp**oɁok | 1 | 1 | 1 | 0 | 0 | 1 | 0 | -11 |  |
| /pampoɁok/ pa**mp**oɁok | 0 | 1 | 1 | 0 | 0 | 0 | 0 | -2 |  |
| /pampoɁok/ pa**m**oɁok | 0 |  | 0 | 0 | 0 | 0 | 1 | -4 |  |

Ident-Placecomposed

Maxcomposed

Ident-Placelisted

Maxlisted

**Parsing liason consonants:**

- I am not writing a morphological parser, that is a big job, which is not the purpose of this here learner.

- Therefore, the user must specify in advance segments which they would like to designate as ambiguous (liason consonants)

When the model is initialized, these will be stored on both morphemes, with an activity value of 0

﻿input lexeme surface obs.prob tab.prob

1syll I I 1 1

2syll IO IO 1 1

3syll IOO IOO 1 1

3syll2 OIO OIO 1 1

Types of input file:

regular tableau without hidden structure – constraint violations prespecified

regular tableau *with* hidden structure – constraint violations prespecified

regular tableau with the ability to change activity values, and therefore max/dep violation values, but all other constraint violations prespecified

regular tableau with the ability to add PFCs, but markedness violations prespecified

simple training data, candidates to be generated on the fly

**Input files:**

Input files are read in as a trainingData object. This class contains:

* trainingData.lexicon - a dictionary where each key is a lexeme tag, and each entry is a lexeme object. The lexicon is mutable: entries may be added or (rarely) removed during learning.
* trainingData.learnData - a list derived from all rows of the input file. Each entry is itself a list: [lexemes, surface], where lexemes is *also* a list of lexeme objects, and surface is a string - the surface form that that combination of lexemes should take[[1]](#footnote-2). (the 'observed' form)
* trainingData.sampler - a list derived from all rows of the input file, indicating sampling probability for each entry in learnData. Derived from obs.prob, or obs.prob\*tab.prob in the input file.
* trainingData.constraintNames - a list of constraint names, taken from constraint columns of the input file

If your input contains candidates, the following will also be part of the trainingData object:

* trainingData.tableaux - a list of Tableau objects
* trainingData.tabProb - a sampler for the Tableau objects, used in very simple learning problems
* trainingData.tableauxTags - a list of tags that corresponds to the list of Tableau objects

possible input columns:

**input - required**

**obs.prob - required**

**candidate**

**surface - at least one of candidate, surface is required**

**tab.prob**

**lexeme**

**constraint names**

In all columns, morphemes are separated by underscore \_

**Input:**

The input column must contain one unique value for each phonological input you want the learner to experience. Inputs may be one or more morphemes, separated by \_. All rows with a particular input will be treated as the same learning datum.

|  |  |  |  |
| --- | --- | --- | --- |
|  | input | candidate | obs.prob |
| *input 1* | pat | pat | 1 |
| pat | pad | 0 |
| *input 2* | pat\_ka | pat\_ka | 0.7 |
| pat\_ka | pad\_ka | 0 |
| pat\_ka | pad\_ga | 0.3 |
| *input 3* | taka | taka | 0.3 |
| taka | taga | 0.7 |
| *input 4* | taka\_ka | taka\_ka | 0.3 |
| taka\_ka | taga\_ka | 0.7 |
| *input 2* | pat\_ka | pa\_ka | 0 |

In this example tableau, there are 4 unique inputs. Note that the final row is also an instance of input 2, although it isn't contiguous with the other input 2 rows. The learner will create a tableau with all four rows together.

In this example, there are just 3 unique morphemes, *pat*, *taka*, and *ka.* When this file is read in, the lexicon will contain these three entries.

t = l.trainingData("manual\_examples/firstInputExample.txt")

t.lexicon

>{'pat': pat, 'ka': ka, 'taka': taka}

t.learnData

>[[[pat], 'pat', 'pat'],

[[pat], 'pad', 'pat'],

[[pat, ka], 'pat\_ka', 'pat\_ka'],

[[pat, ka], 'pad\_ka', 'pat\_ka'],

[[pat, ka], 'pad\_ga', 'pat\_ka'],

[[taka], 'taka', 'taka'],

[[taka], 'taga', 'taka'],

[[taka, ka], 'taka\_ka', 'taka\_ka'],

[[taka, ka], 'taga\_ka', 'taka\_ka'],

[[pat, ka], 'pa\_ka', 'pat\_ka']]

t.sampler

> [0.25, 0.0, 0.175, 0.0, 0.075, 0.075, 0.175, 0.075, 0.175, 0.0]

note that outputs with 0 observed probability still appear in t.learnData, but they correspond to a value of 0 in the sampler.

**Delineating morpheme boundaries**

In all columns, morpheme boundaries are delineated using "\_". The input, lexeme, candidate, and surface columns must all have the same number of morphemes.

|  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- |
|  | input | lexeme[[2]](#footnote-3) | candidate | surface | obs.prob |
| *a* | PLAY\_INF | pat\_ka | pat\_ka | pat\_ka | 1 |
| *b* | PLAY\_INF | pat\_ka | patS\_ka | patS\_ka | 0 |
| *c* | PLAY\_1SG\_PL | pat\_i\_ka | pat\_i\_ka | pat\_i\_ka | 0 |
| *d* | PLAY\_1SG\_PL | pat\_i\_ka | patS\_i\_ka | patS\_i\_ka | 1 |
| *e* | PLAY\_1SG\_PL | pat\_i\_ka | pat\_\_ka | pat\_\_ka | 0 |
| *f* | PLAY | pat | pat | pat | 1 |
| *g* | PLAY | pat | patS | patS | 0 |

This input file will create a lexicon with four entries: PLAY, INF, 1SG, and PL. Note that in row e, the candidate and surface forms have two \_'s next to each other - indicating that in this case the 1SG morpheme surfaces as null.

The following illustrates some inputs that will not align properly. Problem cells are highlighted in red.

row a, input: the '+' character does not delineate morphemes. Instead the whole cell will be read in as a single morpheme

row a, lexeme: this will count as only one morpheme, which mismatches the candidate column and the surface column

row c, surface: missing morpheme divisions that are present in the other columns

row d, candidate: same problem

row e: candidate and surface are meant to represent the case where the 1SG morpheme is deleted, but they must be written with two \_'s in order for the 1SG morpheme to be aligned with null properly.

|  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- |
|  | input | lexeme[[3]](#footnote-4) | candidate | surface | obs.prob |
| *a* | PLAY+INF | patka | pat\_ka | pat\_ka | 1 |
| *b* | PLAY\_INF | pat\_ka | patS\_ka | patS\_ka | 0 |
| *c* | PLAY\_1SG\_PL | pat\_i\_ka | pat\_i\_ka | patika | 0 |
| *d* | PLAY\_1SG\_PL | pat\_i\_ka | patSi\_ka | patS\_i\_ka | 1 |
| *e* | PLAY\_1SG\_PL | pat\_i\_ka | pat\_ka | pat\_ka | 0 |
| *f* | PLAY | pat | pat | pat | 1 |
| *g* | PLAY | pat | patS | patS | 0 |

**when morpheme boundaries are uncertain:**

Sometimes, you would like the learner to consider multiple parses. This is generally necessary when learning with GSR's, since segments can be part of two morphemes at once.

when specialLex is true:

ambiguous material can be delineated with '-'

peti-t-ami will align with petit\_ami, but will generate UR's petit and tami, both with an activity of 0 on the t segment.

**Specifying hidden structure in the tableau**

You can specify hidden structure by including both candidate and surface columns

|  |  |  |  |
| --- | --- | --- | --- |
| input | candidate | surface | obs.prob |
| pataka | pa(ta)ka | patAka | 0.9 |
| pataka | pa(taka) | patAka | 0.9 |
| pataka | (pata)ka | pAtaka | 0.1 |

In this tableau, candidates contain feet, which in some cases produce an identical stress pattern as output. The first two candidates here produce the same stress pattern. Here, 'surface' specifies the surface form of the candidate, which should correspond to features that would be audible, or extractable from a corpus. Stress pattern is such a feature, but foot structure is not. Other hidden structures might include deletion+insertion vs. feature change, word-internal correspondence structure, floating features, and others.

When using hidden structure, you must give all copies of the same surface form the same obs.prob value. The above tableau indicates that patAka is observed 90% of the time, and pAtaka is observed 10% of the time.

Note that it is also possible to specify only surface forms and not candidates. This structure is useful when you want the learner to generate candidates for you.

|  |  |  |
| --- | --- | --- |
| input | surface | obs.prob |
| pataka | patAka | 1 |
| badupi | bAdupi | 1 |
| potolu | potolU | 1 |

You can also do variation this way, just making a row for each observed surface form:

|  |  |  |
| --- | --- | --- |
| input | surface | obs.prob |
| pataka | patAka | .9 |
| pataka | pAtaka | .1 |
| badupi | bAdupi | 1 |
| potolu | potolU | 1 |

**Labels vs. underlying forms**

For many problems, you may want the lexeme's label to be different from its underlying form. (This won't matter unless you are (a) assigning violations by formula, or (b) generating candidates during learning)

In this case, use the input column to define lexeme labels, and the lexeme column to define UR's. The material in lexeme should all be characters which are in your features file.

|  |  |  |  |
| --- | --- | --- | --- |
| input | lexeme | candidate | obs.prob |
| an\_apple | xn\_@pl | xn\_@pl | 1 |
| an\_apple | xn\_@pl | x\_@pl | 0 |
| an\_tree | xn\_tri | x\_tri | 1 |
| an\_tree | xn\_tri | xn\_tri | 0 |
| an | xn | xn | 1 |
| an | xn | x | 0 |
| apple | @pl | Q@apl | 1 |
| apple | @pl | @apl | 0 |
| apple\_PL | @pl\_z | Q@pl\_z | 1 |
| apple\_PL | @pl\_z | @pl\_z | 0 |

This input file will generate four lexemes:

'an', whose UR is schwa followed by n

'apple' whose UR is /aepl/

'tree' whose UR is /tri/

'PL' whose UR is /z/

These UR's can be used to generate candidates, and assess faithfulness violations. But, the actual labels of the lexemes can still be more human-readable.

**Observed probabilities:**

There are two columns used to specify how often a particular output is observed: obs.prob, and tab.prob.

obs.prob is obligatory, and determines the observed probability of each distinct surface form. This is how this column has been used so far in this section.

Here is another example:

|  |  |  |
| --- | --- | --- |
| input | surface | obs.prob |
| xn\_@pl | xn@pl | .9 |
| xn\_@pl | x@pl | .1 |
| xn\_tri | xtri | 1 |
| xn | xn | 1 |
| @pl | Q@apl | 1 |

This input file specifies that 'xn\_@pl' surfaces with the n 90% of the time, and without it 10% of the time. Meanwhile, xn\_tri surfaces as x\_tri always, 'xn' surfaces faithfully always, and '@pl' surfaces with glottal stop insertion (Q = ?) always.

With this input file, the learner will get each of these inputs equally often. Suppose I want the learner to train on 'xn\_@pl' more often than on 'xn\_tri'. I can do this two ways. The simplest is to use the tab.prob column.

|  |  |  |  |
| --- | --- | --- | --- |
| input | surface | obs.prob | tab.prob |
| xn\_@pl | xn@pl | .9 | 4 |
| xn\_@pl | x@pl | .1 | 4 |
| xn\_tri | xtri | 1 | 2 |
| xn | xn | 1 | 1 |
| @pl | Q@apl | 1 | 1 |

With this input file, the input 'xn\_@pl' will be trained 4 times as often as 'xn' or '@pl'. 'xn\_tri' will be trained twice as often as 'xn' or '@pl'.

In the code, these two columns are combined together to create the trainingData.sampler object. This means that you don't really have to use both of them; you can specify both within-input probabilities and input frequencies together:

|  |  |  |
| --- | --- | --- |
| input | surface | obs.prob |
| xn\_@pl | xn@pl | 3.6 |
| xn\_@pl | x@pl | .4 |
| xn\_tri | xtri | 2 |
| xn | xn | 1 |
| @pl | Q@apl | 1 |

**Constraints:**

Any column after the pre-defined columns will be interpreted as a constraint.

Constraint names can be whatever you want, except for the pre-defined names (input, lexeme, candidate, surface, obs.prob, and tab.prob)

Violations should be positive numbers, and cells with no violations should be filled with 0

|  |  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- | --- |
| input | lexeme | candidate | obs.prob | Onset | Dep | Max |
| an\_apple | xn\_@pl | xn\_@pl | 1 | 0 | 0 | 0 |
| an\_apple | xn\_@pl | x\_@pl | 0 | 1 | 0 | 1 |
| an\_tree | xn\_tri | x\_tri | 1 | 0 | 0 | 1 |
| an\_tree | xn\_tri | xn\_tri | 0 | 0 | 0 | 0 |
| an | xn | xn | 1 | 1 | 0 | 0 |
| an | xn | x | 0 | 1 | 0 | 1 |
| apple | @pl | Q@apl | 1 | 0 | 1 | 0 |
| apple | @pl | @apl | 0 | 1 | 0 | 0 |
| apple\_PL | @pl\_z | Q@pl\_z | 1 | 0 | 1 | 0 |
| apple\_PL | @pl\_z | @pl\_z | 0 | 1 | 0 | 0 |

Constraint names: if you append \_listed to the end of a constraint name (it should be a faithfulness constraint), then that constraint will apply only to listed whole-forms in the UseListed framework, and will not apply to composed forms, or single morphemes.

You need two constraints that are identical except for the \_listed designation. E.g. Ident-Voice\_listed, and Ident-Voice

Tableaux file:

The Tableaux file is similar to what you would use in other phonological learning software, such as OTSoft, OT-Help, or hgr.

Candidate Generation function:

**NOTE:** there are some odd dependencies with the morphFeature operation – as defined now, you won’t ever get a candidate that changes, say, backness AND rounding unless there’s a segment in the language that just does one of those. Need to fix!

In order to generate candidates, we traverse a tree, that begins with the faithful candidate. (Note that in GSR’s there are multiple faithful candidates)

Daughters of each node are computed via operations, which must be defined by the user of the program. Here is an example (part of) a tree, with the following operations;

Delete a single segment

Add a glottal stop [ʔ]

Change a single feature

pʌdəl

Add

Delete

Change-feature

pʌʔdəl

pʌdəɹ

pʌdəlʔ

...

... ʌdəl

pdəl

pʌəl

pʌdl

ʔpʌdəl

phʌdəl

pədəl

pʌɾəl

phdəl

pdʌl

pɾəl

...

...

pʔdəl

dəl

ʔənæpl̩

ʔənæp

ʔənæp

ʔnæpl̩

How tree search proceeds:

Beginning with the root node, the faithful candidate, operations are considered with probabili

Rules about tree search:

1. If a candidate is already in the candidate set, it is never added again
   1. This means its daughters won’t be generated again
2. Operations are considered as a wgo

Calculating the Gradient:

Need the softmax function. Did you know that the MaxEnt function is a softmax function???

- explanation and history here: <https://deepai.org/machine-learning-glossary-and-terms/softmax-layer> -- Boltzmann or Gibbs distribution!!!

<https://towardsdatascience.com/derivative-of-the-softmax-function-and-the-categorical-cross-entropy-loss-ffceefc081d1>

log likelihood:

Given any probability model, we can calculate the likelihood of our data given our model. Let's take a simple example with dice. Say, I have a fair 6-sided die. In this case, my probability model would be that each side of the die has an equal chance of being rolled. So, we expect to roll a 1 with probability of 1/6. Likewise, we expect to roll a 2 with a probability of 1/6, and a 3 a 4 and so on.

So, suppose our data are that we rolled the die 3 times and got: 1, 2, 1

The likelihood of rolling the first 1 is 1/6. So is the likelihood of rolling the 2, and the second 1. So, since P(A and B) = P(A)\*P(B), the likelihood of the data here is 1/6^3 = 0.00463

Just to illustrate what this would look like if we switched out the theory, which is what we'll be doing when we optimize weights: suppose our probability model was instead that the die only ever rolls 1's. So, the probability of a 1 is 1, and the probability of a 2 or any other number is 0. Now, our Likelihood is 1\*0\*1 = 0

One more: suppose our probability model was that the die was heavily biased towards 1's: p(1)= 0.9, prob of any other number is 0.016. Now, the likelihood of our data is .9\*.016\*.9 = 0.013]

So, according to our likelihoods, that third model is the best, since it yields the highest likelihood.

So, how do we scale up this likelihood calculation? Let's assume we've got a set of tableaux, with a bunch of observed counts, and predicted probabilities.

|  |  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- | --- |
|  | ℂ1 | ℂ2 | ℂ3 | 𝓗 | *predicted* | *observed* |
|  | 1.1 | 2.3 | 8.7 |  |  |  |
| /Input/1 |  |  |  |  |  |  |
| *cand1,1* | 1 | 1 | 0 | -3.4 | *0.25* | *415* |
| *cand1,2* | 1 | 0 | 1 | -9.8 | *0* | *12* |
| *cand1,3* | 0 | 1 | 0 | -2.3 | *0.75* | *891* |
| /Input/2 |  |  |  |  |  |  |
| *cand2,1* | 0 | 0 | 1 | -8.7 | *0* | *1* |
| *cand2,2* | 1 | 0 | 0 | -1.1 | *1* | *345* |

Now, for each candidate, we have a number of observations of that candidate, perhaps from a corpus or an experiment, and those observations together are our data. Each observation of each candidate has a likelihood of appearing equal to its predicted probability. So, *cand1,1* has a 0.24 chance of occurring. If it occurred twice, the chances of that happening would just be 0.24\*0.24. If it occurred *n* times, the chances of that are 0.24*n*. The same is true for each candidate:

|  |  |  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- | --- | --- |
|  | ℂ1 | ℂ2 | ℂ3 | 𝓗 | *predicted* | *observed* | **Likelihood** |
|  | 1.1 | 2.3 | 4.8 |  |  |  |  |
| /Input/1 |  |  |  |  |  |  |  |
| *cand1,1* | 1 | 1 | 0 | -3.4 | *0.24* | *415* | **0.24415** |
| *cand1,2* | 1 | 0 | 1 | -5.9 | *0.02* | *12* | **0.0212** |
| *cand1,3* | 0 | 1 | 0 | -2.3 | *0.74* | *891* | **0.74891** |
| /Input/2 |  |  |  |  |  |  |  |
| *cand2,1* | 0 | 0 | 1 | -4.8 | *0.02* | *1* | **0.021** |
| *cand2,2* | 1 | 0 | 0 | -1.1 | *0.98* | *345* | **0.98345** |

So how do we get the likelihood of the whole dataset? It's just each candidate's likelihood multiplied together:

or:

Before we move on, a brief note: You will often see models fit with observed probabilities, rather than observed counts. For the tableaux above, that would be 0.315 instead of 415, 0.009 instead of 12, 891 instead of 0.68, and so on. So then the likelihoods for each candidate would be 0.240.315, 0.020.009, and so on. Technically when you do this, you are not calculating the likelihood, but instead something called Cross Entropy. You can optimize Cross Entropy just like you would likelihood, and all the math below applies.

When to go with Likelihood and when to go with Cross Entropy:

You should use observed probabilities, and therefore cross entropy, when you want to fit to type frequency only, and you don't want any input's distribution to be more important than another's. You should use observed counts, and therefore likelihood, when you want to fit to token and type frequency. With observed counts, inputs with fewer total observations will contribute less to the likelihood calculation, and therefore may influence your learned constraint weights less.

Let's return to this likelihood function:

Now, 0.24415 is already a teeny tiny number, which most computers will round to zero. After we multiply all these probabilities together, the problem just gets worse. So, to fix this, we will optimize for the *log* of the likelihood instead of the likelihood itself.

How does this help? Well, log(x\*y) = log(x) + log(y), so:

And, log(xy) = ylog(x), so:

Now, we can plug in our actual calculation for the predicted probability, which is:

Recall also that the harmony of a candidate, is just the dot product of the constraint weights with that candidate's violations. If we also plug that into our equation, we get the full likelihood function, in terms of our weights:

Suppose we have *n* weights:

We need the gradient of this, which will be . In other words, we need a vector of the partial derivative of the likelihood function with respect to each weight. It will be the same length as .

First, let's simplify that objective function a little:

Now, let's rewrite, expanding those sums and dot products out:

This is the first term of the second input

Where r is the number of inputs

So each of these terms can be differentiated, and each entry in the gradient will be a sum of all of them. But, because they have the same form except for constants, let's just demo with a single term:

We have to do the chain rule here, that's what that last term is. Here

so: that last term in terms of *u* is

so:

and to get we actually have to apply the chain rule one more time, this time for each term of u(w1)

For a single term:

The derivative of this is just

So,

The rest of the gradient terms are just the same, but replace w1 with w2 and so on

1. Note that if there is variation, there may be multiple entries in learnData with the same set of lexemes but different surface strings. [↑](#footnote-ref-2)
2. see below for an explanation of this column [↑](#footnote-ref-3)
3. see below for an explanation of this column [↑](#footnote-ref-4)