

TPGMM

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Chapter 1

Hierarchical Index

1.1 Class Hierarchy

This inheritance list is sorted roughly, but not completely, alphabetically:

ABC

tpgmm.utils.learning_modules.ClassificationModule	5
tpgmm.TPGMM	17
tpgmm.utils.learning_modules.LearningModule	14
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tpgmm.utils.learning_modules.RegressionModel	16
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Chapter 2

Class Index

2.1 Class List

Here are the classes, structs, unions and interfaces with brief descriptions:

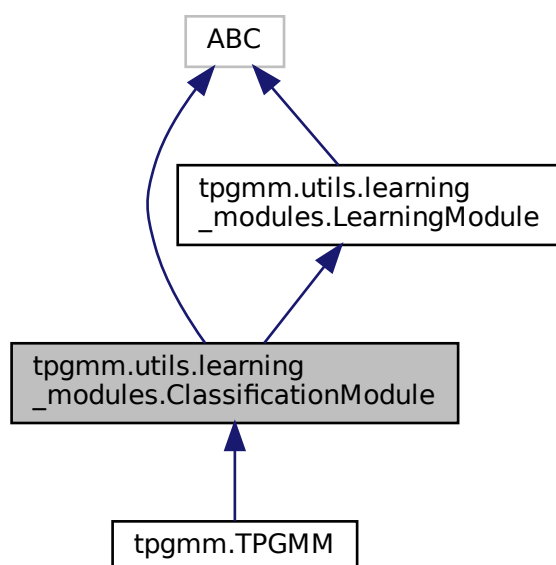
tpgmm.utils.learning_modules.ClassificationModule	5
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This class implements a gaussian mixture regression model	10
tpgmm.utils.learning_modules.LearningModule	
Basic abstract class for a generic learning module	14
tpgmm.utils.learning_modules.RegressionModel	
Basic Regression Model	16
tpgmm.TPGMM	
This class is an implementation of the task parameterized gaussian mixture model according to Calinon Paper @ https://calinon.ch/papers/Calinon-JIST2015.pdf	17

Chapter 3

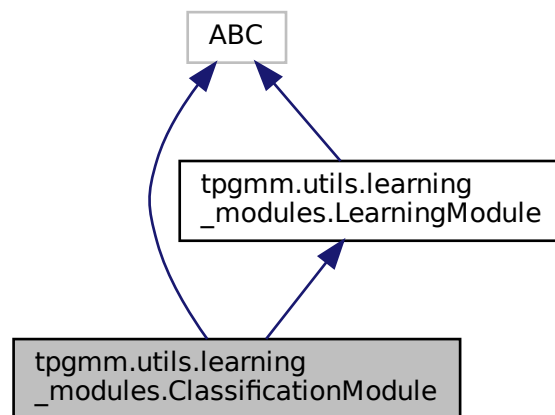
Class Documentation

3.1 tpgmm.utils.learning_modules.ClassificationModule Class Reference

Inheritance diagram for tpgmm.utils.learning_modules.ClassificationModule:



Collaboration diagram for `tpgmm.utils.learning_modules.ClassificationModule`:



Public Member Functions

- None `__init__` (self, int n_components)
- def `predict_proba` (self, *args, **kwargs)
Predict class probabilities for the input data.
- float `score` (self, ndarray X)
Calculate the score function from the descendant.
- float `silhouette_score` (self, ndarray X)
Calculate the silhouette score for the given data.
- float `inertia` (self, ndarray X)
Calculate the sum of squared distances of samples to their closest cluster center.
- float `davies_bouldin_score` (self, ndarray X)
calculates davies_bouldin_score on given data
- float `bic` (self, ndarray X)
calculates the bayesian information criterion as in
- def `aic` (self, ndarray X)
calculates the Akaike information criterion as in https://scikit-learn.org/stable/modules/linear_model.html#aic-bic
- Dict[str, Any] `config` (self)
Get the configuration parameters.

3.1.1 Member Function Documentation

3.1.1.1 `aic()`

```
def tpgmm.utils.learning_modules.ClassificationModule.aic (
    self,
    ndarray X )
```

calculates the Akaike information criterion as in https://scikit-learn.org/stable/modules/linear_model.html#aic-bic

Parameters

<i>X</i>	data tensor with expected shape same form self.score(X)
----------	---

Returns

float aic score

3.1.1.2 bic()

```
float tpgmm.utils.learning_modules.ClassificationModule.bic (
    self,
    ndarray X )
```

calculates the bayesian information criterion as in

https://scikit-learn.org/stable/modules/linear_model.html#aic-bic

Parameters

<i>X</i>	data tensor with expected shape (num_points, num_features)
----------	--

Returns

float bic score

Reimplemented in [tpgmm.TPGMM](#).

3.1.1.3 davies_bouldin_score()

```
float tpgmm.utils.learning_modules.ClassificationModule.davies_bouldin_score (
    self,
    ndarray X )
```

calculates davies_bouldin_score on given data

$$DB = \frac{1}{n} \sum_{i,j=1}^N \max_{j \neq i} \frac{\sigma_i + \sigma_j}{d(c_i, c_j)}$$

Where

N ... number of datapoints d() ... distance function c_i ... cluster center σ_i ... the average distance of all points in cluster i from the cluster centre c_i

Parameters

<i>X</i>	data in the shape for self.predict()
----------	--------------------------------------

Returns

float score value

3.1.1.4 inertia()

```
float tpgmm.utils.learning_modules.ClassificationModule.inertia (  
    self,  
    ndarray X )
```

Calculate the sum of squared distances of samples to their closest cluster center.

Parameters

<i>X</i>	data in local reference frames. Shape (num_frames, num_points, num_features)
----------	--

Returns

float inertia

3.1.1.5 score()

```
float tpgmm.utils.learning_modules.ClassificationModule.score (  
    self,  
    ndarray X )
```

Calculate the score function from the descendant.

Often the score is calculated based on the optimization objective.

Parameters

<i>X</i>	Data to calculate the score on.
----------	---------------------------------

Returns

float The calculated score.

Reimplemented in [tpgmm.TPGMM](#).

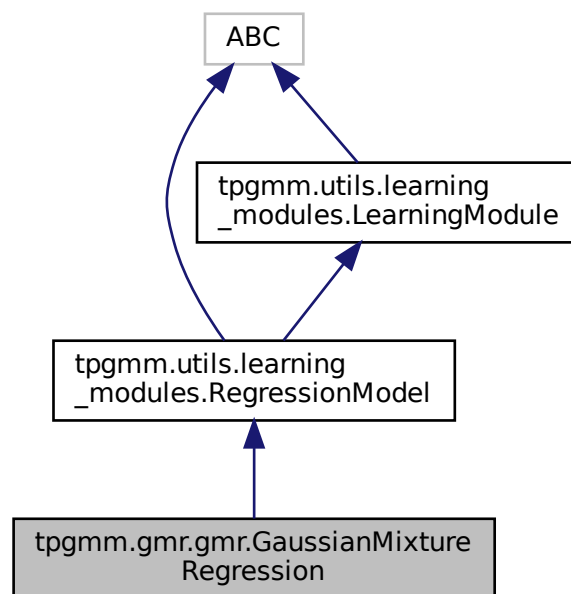
The documentation for this class was generated from the following file:

- `documentation/tmp/tpgmm/utils/learning_modules.py`

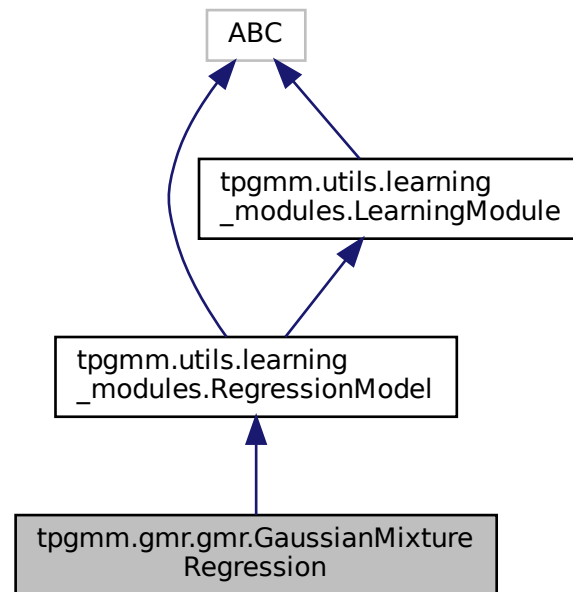
3.2 tpgmm.gmr.gmr.GaussianMixtureRegression Class Reference

This class implements a gaussian mixture regression model.

Inheritance diagram for `tpgmm.gmr.gmr.GaussianMixtureRegression`:



Collaboration diagram for tpgmm.gmr.gmr.GaussianMixtureRegression:



Public Member Functions

- None **__init__** (self, ndarray weights, ndarray means, ndarray covariances, Iterable[int] input_idx)
- "GaussianMixtureRegression" **from_tpgmm** ("GaussianMixtureRegression" cls, TPGMM tpgmm, Iterable[int] input_idx)
- None **fit** (self, ndarray translation, ndarray rotation_matrix)
Turns the task_parameterized gaussian mixture model into a single gaussian mixture model.
- Tuple[ndarray, ndarray] **predict** (self, ndarray input_data)
this function is inspired by formula 13 in Calinon paper it creates for each given data point its own parameterized gaussian distribution.
- int **num_input_features** (self)
get number of input features.
- int **num_output_features** (self)
get number of output features.
- List[int] **output_idx** (self)
- Dict[str, Any] **config** (self)
Get all model config parameters.

Public Attributes

- **tpgmm_means_**
- **tpgmm_covariances_**
- **gmm_weights**
- **input_idx**
- **xi_**
- **sigma_**

3.2.1 Detailed Description

This class implements a gaussian mixture regression model.

This model was described in 'A Tutorial on Task-Parameterized Movement Learning and Retrival' from S.Calvinon at: <https://calinon.ch/papers/Calinon-JIST2015.pdf>

The class fits a gaussian mixture regression on a given Gaussian Mixture model or a Task-Parameterized gaussian mixture model. Used equations are:

$$\mathcal{P}(\phi_t^{\mathcal{O}} | \phi_t^{\mathcal{I}}) \sim \sum_{i=1}^K h_i(\phi_t^{\mathcal{I}}) \mathcal{N}(\hat{\mu}_t^{\mathcal{O}}(\phi_t^{\mathcal{I}}), \hat{\Sigma}_t^{\mathcal{O}})$$

$$\hat{\mu}_i^{\mathcal{O}}(\phi_t^{\mathcal{I}}) = \mu_i^{\mathcal{O}} + \Sigma_i^{\mathcal{OI}} \Sigma_i^{\mathcal{I},-1} (\phi_t^{\mathcal{I}} - \mu_i^{\mathcal{I}})$$

$$\hat{\Sigma}_t^{\mathcal{O}} = \Sigma_i^{\mathcal{O}} - \Sigma_i^{\mathcal{OI}} \Sigma_i^{\mathcal{I},-1} \Sigma_i^{\mathcal{OI}}$$

$$h_i(\phi_t^{\mathcal{I}}) = \frac{\pi_i \mathcal{N}(\phi_t^{\mathcal{I}} | \mu_i^{\mathcal{I}}, \Sigma_i^{\mathcal{I}})}{\sum_k^K \pi_k \mathcal{N}(\phi_t^{\mathcal{I}} | \mu_k^{\mathcal{I}}, \Sigma_k^{\mathcal{I}})}$$

Examples

```
>>> # trajectory data in local reference frames with shape: (num_reference_frames, num_points, 4)
>>> # last dimension could be for example: x, y, z, time
>>> # in case you got multiple trajectories: concatenate all of them in axis 1
>>> trajectories = load_trajectories()
>>> tpgmm = TPGMM(n_components=5)
>>> tpgmm.fit(trajectories)
>>> # gmr for the first reference frame
>>> gmr = GaussianMixtureRegression(weights=tpgmm.weights_, means=tpgmm.means_[0],
>>>   covariances=tpgmm.covariances_[0], input_idx=[3])
>>> gmr.fit(trajectory[0])
```

3.2.2 Member Function Documentation

3.2.2.1 fit()

```
None tpgmm.gmr.gmr.GaussianMixtureRegression.fit (
    self,
    ndarray translation,
    ndarray rotation_matrix )
```

Turns the task_parameterized gaussian mixture model into a single gaussian mixture model.

function is performing equation (5) and (6) from calinon paper

TODO write formulas

Parameters

<i>translation</i>	translation matrix for translating into desired frames. Shape (num_frames, num_output_features)
<i>rotation_matrix</i>	rotation matrix for rotating into desired frames. Shape (num_frames, num_output_features, num_output_features)

3.2.2.2 num_input_features()

```
int tpgmm.gmr.gmr.GaussianMixtureRegression.num_input_features (
    self )
```

get number of input features.

Example

if you have 4 features: x, y, z, time. You can define with input_idx=[3] time as the input feature. This function will return then 1

Returns

int number of input features

3.2.2.3 num_output_features()

```
int tpgmm.gmr.gmr.GaussianMixtureRegression.num_output_features (
    self )
```

get number of output features.

Example

if you have 4 features: x, y, z, time. You can define with input_idx=[3] time as the input feature. This function will return then 3 for x,y and z

Returns

int number of output features

3.2.2.4 predict()

```

Tuple[ndarray, ndarray] tpgmm.gmr.gmr.GaussianMixtureRegression.predict (
    self,
    ndarray input_data )

```

this function is inspired by formula 13 in Calinon paper it creates for each given data point its own parameterized gaussian distribution.

The mechanics are described by:

$$\mathcal{P}(\phi_t^{\mathcal{O}} | \phi_t^{\mathcal{I}}) \sim \sum_{i=1}^K h_i(\phi_t^{\mathcal{I}}) \mathcal{N}(\hat{\mu}_t^{\mathcal{O}}(\phi_t^{\mathcal{I}}), \hat{\Sigma}_t^{\mathcal{O}})$$

in:
self.h()

$$h_i(\phi_t^{\mathcal{I}}) = \frac{\pi_i \mathcal{N}(\phi_t^{\mathcal{I}} | \mu_i^{\mathcal{I}}, \Sigma_i^{\mathcal{I}})}{\sum_k^K \pi_k \mathcal{N}(\phi_t^{\mathcal{I}} | \mu_k^{\mathcal{I}}, \Sigma_k^{\mathcal{I}})}$$

in:
self.mu_hat()

$$\hat{\mu}_i^{\mathcal{O}}(\phi_t^{\mathcal{I}}) = \mu_i^{\mathcal{O}} + \Sigma_i^{\mathcal{OI}} \Sigma_i^{\mathcal{I},-1} (\phi_t^{\mathcal{I}} - \mu_i^{\mathcal{I}})$$

in:
self.sigma_hat()

$$\hat{\Sigma}_t^{\mathcal{O}} = \Sigma_i^{\mathcal{O}} - \Sigma_i^{\mathcal{OI}} \Sigma_i^{\mathcal{I},-1} \Sigma_i^{\mathcal{OI}}$$

Parameters

<i>data</i>	Shape: (num_points, num_input_features)
-------------	---

Returns

Tuple[ndarray, ndarray] : mu: shape -> (num_points, num_output_features), cov: shape (num_points, num_output_features, num_output_features)

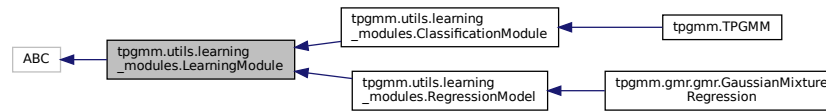
The documentation for this class was generated from the following file:

- documentation/tmp/tpgmm/gmr/gmr.py

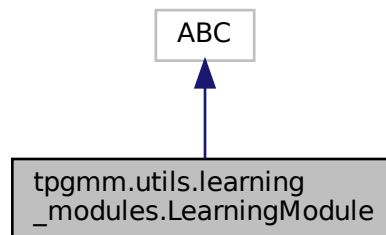
3.3 tpgmm.utils.learning_modules.LearningModule Class Reference

Basic abstract class for a generic learning module.

Inheritance diagram for tpgmm.utils.learning_modules.LearningModule:



Collaboration diagram for tpgmm.utils.learning_modules.LearningModule:



Public Member Functions

- None `__init__` (self)
- None `fit` (self, *args, **kwargs)
Basic abstract class for a generic learning module.
- def `predict` (self, *args, **kwargs)
Predict output using the fitted model.
- ndarray `fit_predict` (self, ndarray X)
Convenience method; equivalent to calling `fit(data)` followed by `predict(data)`.
- Dict[str, Any] `config` (self)
Get all model config parameters.

3.3.1 Detailed Description

Basic abstract class for a generic learning module.

3.3.2 Member Function Documentation

3.3.2.1 fit_predict()

```

ndarray tpgmm.utils.learning_modules.LearningModule.fit_predict (
    self,
    ndarray X )

```

Convenience method; equivalent to calling `fit(data)` followed by `predict(data)`.

Parameters

<i>X</i>	Data in local reference frames. Shape (num_frames, num_points, num_features).
----------	---

Returns

ndarray The label for each data-point. Shape (num_points).

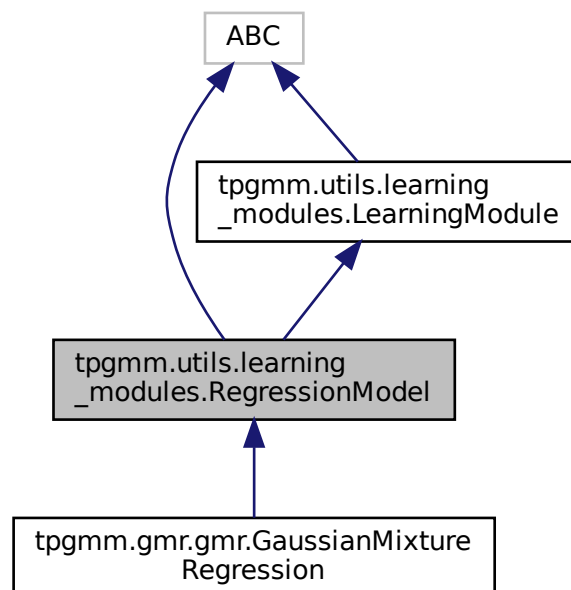
The documentation for this class was generated from the following file:

- documentation/tmp/tpgmm/utils/learning_modules.py

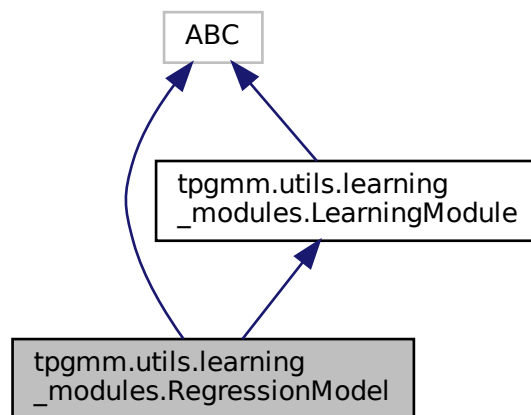
3.4 tpgmm.utils.learning_modules.RegressionModel Class Reference

Basic Regression Model.

Inheritance diagram for tpgmm.utils.learning_modules.RegressionModel:



Collaboration diagram for tpgmm.utils.learning_modules.RegressionModel:



Additional Inherited Members

3.4.1 Detailed Description

Basic Regression Model.

Implements all interfaces and common methods for regression models.

Basic Classification Model. Implements all interfaces and common methods for classification models.

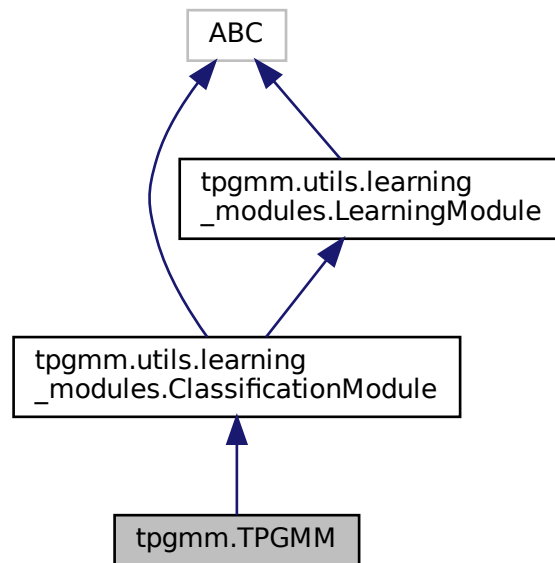
The documentation for this class was generated from the following file:

- documentation/tmp/tpgmm/utils/learning_modules.py

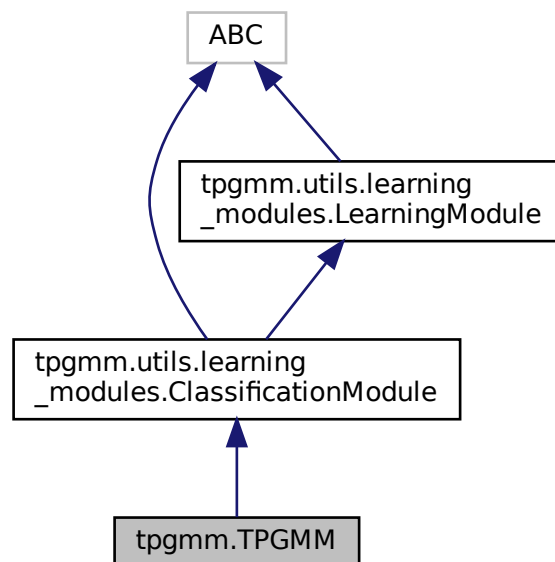
3.5 tpgmm.TPGMM Class Reference

This class is an implementation of the task parameterized gaussian mixture model according to Calinon Paper @<https://calinon.ch/papers/Calinon-JIST2015.pdf>.

Inheritance diagram for tpgmm.TPGMM:



Collaboration diagram for tpgmm.TPGMM:



Public Member Functions

- None `__init__` (self, int n_components, float threshold=1e-7, int max_iter=100, int min_iter=5, ndarray weights_init=None, ndarray means_init=None, float reg_factor=1e-5, bool verbose=False)
*Initialize the **TPGMM** class.*
- None `fit` (self, ndarray X)
*fits X on the task parameterized gaussian mixture model using K-Means clustering as default initialization method and executes the expectation maximization algorithm:
E-Step:*
- ndarray `predict` (self, ndarray X)
predict cluster labels for each data point in X
- ndarray `predict_proba` (self, ndarray X)
predict cluster labels for each data point
- float `silhouette_score` (self, ndarray X)
calculated the silhouette score of the model over the given metric and given data x TODO(Marco Todescato): please review this function if the merge for the silhouette score is correct
- float `score` (self, ndarray X)
calculate log likelihood score given data
- float `bic` (self, ndarray X)
calculates the bayesian information criterion as in
- ndarray `gauss_cdf` (self, ndarray X)
calculate the gaussian probability for a given data set.
- Dict[str, Any] `config` (self)
*Returns a dictionary containing the configuration parameters of the **TPGMM** model.*

Public Attributes

- `weights_`
ndarray of shape (n_components,) Weights between gaussian components.
- `means_`
ndarray of shape (num_frames, n_components, num_features) Mean matrix for each frame and component.
- `covariances_`
ndarray of shape (num_frames, n_components, num_features, num_features).

3.5.1 Detailed Description

This class is an implementation of the task parameterized gaussian mixture model according to Calinon Paper @<https://calinon.ch/papers/Calinon-JIST2015.pdf>.

It implements also an Expectation Maximization Algorithm with:

E-Step:

$$h_{t,i} = \frac{\pi_i \prod_{j=1}^P \mathcal{N}(X_t^{(j)} | \mu_i^{(j)}, \Sigma_i^{(j)})}{\sum_{k=1}^K \pi_k \prod_{j=1}^P \mathcal{N}(X_t^{(j)} | \mu_k^{(j)}, \Sigma_k^{(j)})}$$

M-Step:

$$\pi_i \leftarrow \frac{\sum_{t=1}^N h_{t,i}}{N}$$

$$\mu_i^{(j)} \leftarrow \frac{\sum_{t=1}^N h_{t,i} X_t^{(j)}}{\sum_{t=1}^N h_{t,i}}$$

$$\Sigma_i^{(j)} \leftarrow \frac{\sum_{t=1}^N h_{t,i} \left(X_t^{(j)} - \mu_i^{(j)} \right) \left(X_t^{(j)} - \mu_i^{(j)} \right)^T}{\sum_{t=1}^N h_{t,i}}$$

The optimization criterion is the log-likelihood implemented with:

$$LL = \frac{\sum_{t=1}^N \log \left(\sum_{k=1}^K \pi_k \prod_{j=1}^J \mathcal{N} \left(X_t^{(j)} \mid \mu_k^{(j)}, \Sigma_k^{(j)} \right) \right)}{N}$$

Variable explanation:

N ... number of components

π ... weights between components

i ... component index

j ... frame index (like pick or place frame)

μ ... mean

Σ ... variance / covariance matrix

LL ... log likelihood

Examples

```
>> trajectories = load_trajectories() # trajectory data in local reference frames with shape:
    (num_reference_frames, num_points, 3)
>> tpgmm = TPGMM(n_components=5)
>> tpgmm.fit(trajectories)
```

Parameters

<i>n_components</i>	number of components
<i>tol</i>	threshold to break from EM algorithm. Defaults to 1e-3.
<i>max_iter</i>	maximum number of iterations to perform the expectation maximization algorithm. Defaults to 100.
<i>min_iter</i>	minimum number of iterations to perform the expectation maximization algorithm. Defaults to 5.
<i>weights_init</i>	initial weights between each component. If set replaces initialization from K-Means. Defaults to None.
<i>means_init</i>	initial means between each component. If set replaces initialization from K-Means. Defaults to None.
<i>reg_factor</i>	regularization factor for empirical covariance matrix. Defaults to 1e-5.
<i>verbose</i>	Triggers print of learning stats. Defaults to False.

3.5.2 Constructor & Destructor Documentation

3.5.2.1 `__init__()`

```
None tpgmm.TPGMM.__init__ (
    self,
    int n_components,
    float threshold = 1e-7,
    int max_iter = 100,
    int min_iter = 5,
    ndarray weights_init = None,
```



```

ndarray means_init = None,
float reg_factor = 1e-5,
bool verbose = False )

```

Initialize the [TPGMM](#) class.

Parameters

<i>n_components</i>	Number of Gaussian multidimensional distributions to mix.
<i>threshold</i>	Threshold to break from EM algorithm. Defaults to 1e-3.
<i>max_iter</i>	Maximum number of iterations to perform the expectation maximization algorithm. Defaults to 100.
<i>min_iter</i>	Minimum number of iterations to perform the expectation maximization algorithm. Defaults to 5.
<i>weights_init</i>	Initial weights between each component. If set, it replaces initialization from K-Means. Defaults to None.
<i>means_init</i>	Initial means between each component. If set, it replaces initialization from K-Means. Defaults to None.
<i>reg_factor</i>	Regularization factor for the empirical covariance matrix. Defaults to 1e-5.
<i>verbose</i>	Triggers print of learning stats. Defaults to False.

3.5.3 Member Function Documentation

3.5.3.1 bic()

```

float tpgmm.TPGMM.bic (
    self,
    ndarray X )

```

calculates the bayesian information criterion as in

https://scikit-learn.org/stable/modules/linear_model.html#aic-bic

Parameters

<i>X</i>	data tensor with expected shape (num_frames, num_points, num_features)
----------	--

Returns

float bic score

Reimplemented from [tpgmm.utils.learning_modules.ClassificationModule](#).

3.5.3.2 config()

```
Dict[str, Any] tpgmm.TPGMM.config (
    self )
```

Returns a dictionary containing the configuration parameters of the [TPGMM](#) model.

Returns

Dict[str, Any]: Dictionary containing the configuration parameters.

Reimplemented from [tpgmm.utils.learning_modules.ClassificationModule](#).

3.5.3.3 fit()

```
None tpgmm.TPGMM.fit (
    self,
    ndarray X )
```

fits X on the task parameterized gaussian mixture model using K-Means clustering as default initialization method and executes the expectation maximization algorithm:

E-Step:

```
self._update_h()
```

M-Step:

```
self._update_weights_()
self._update_means_()
self._update_covariances()
```

The optimization criterion is the log-likelihood implemented in:

```
self._log_likelihood()
```

The algorithm stops if $LL_{t-1} - LL_t < self.tol$ with LL_t as the log-likelihood at time t.

Parameters

X	data tensor to fit the the task parameterized gaussian mixture model on. Expected shape: (num_frames, num_points, num_features)
-----	---

3.5.3.4 gauss_cdf()

```
ndarray tpgmm.TPGMM.gauss_cdf (
    self,
    ndarray X )
```

calculate the gaussian probability for a given data set.

$$\mathcal{N}\left(X_t^{(j)} \mid \mu_i^{(j)}, \Sigma_i^{(j)}\right) = \frac{1}{\sqrt{(2\pi)^D |\Sigma_i^{(j)}|}} \exp\left(\frac{1}{2}(X_t^{(j)} - \mu_i^{(j)})\Sigma_i^{(j),-1}(X_t^{(j)} - \mu_i^{(j)})^T\right)$$

Variable explanation: D ... number of features

Parameters

<i>X</i>	data with shape: (num_frames, num_points, num_features)
----------	---

Returns

ndarray probability shape (num_frames, n_components, num_points)

3.5.3.5 predict()

```
ndarray tpgmm.TPGMM.predict (
    self,
    ndarray X )
```

predict cluster labels for each data point in X

Parameters

<i>X</i>	data in local reference frames. Shape (num_frames, num_points, num_features)
----------	--

Returns

ndarray the label for each data-point. Shape (num_points)

3.5.3.6 predict_proba()

```
ndarray tpgmm.TPGMM.predict_proba (
    self,
    ndarray X )
```

predict cluster labels for each data point

Parameters

<i>X</i>	data in local reference frames. Shape (num_frames, num_points, num_features)
----------	--

Returns

ndarray cluster probabilities for each data_point. Shape: (num_points, num_components)

3.5.3.7 score()

```
float tpgmm.TPGMM.score (
    self,
    ndarray X )
```

calculate log likelihood score given data

Parameters

<i>X</i>	data tensor with expected shape (num_frames, num_points, num_features)
----------	--

Returns

float log likelihood of given data

Reimplemented from [tpgmm.utils.learning_modules.ClassificationModule](#).

3.5.3.8 silhouette_score()

```
float tpgmm.TPGMM.silhouette_score (
    self,
    ndarray X )
```

calculated the silhouette score of the model over the given metric and given data x TODO(Marco Todescato): please review this function if the merge for the silhouette score is correct

Parameters

<i>X</i>	data in expected shape: (num_frames, num_points, num_features)
<i>metric</i>	<i>description</i> . Defaults to "euclidean".

Reimplemented from [tpgmm.utils.learning_modules.ClassificationModule](#).

3.5.4 Member Data Documentation

3.5.4.1 covariances_

```
tpgmm.TPGMM.covariances_
```

ndarray of shape (num_frames, n_components, num_features, num_features).

Covariance matrix for each frame and component

The documentation for this class was generated from the following file:

- documentation/tmp/tpgmm/tpgmm/tpgmm.py

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