
LUMIN

LUMIN Contributors

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Lumin Unifies Many Improvements for Networks

LUMIN.DATA_PROCESSING PACKAGE

1.1 Submodules

1.2 lumin.data_processing.file_proc module

`lumin.data_processing.file_proc.save_to_grp` (*arr, grp, name*)

Save Numpy array as a dataset in an h5py Group

Parameters

- **arr** (ndarray) – array to be saved
- **grp** (Group) – group in which to save arr
- **name** (str) – name of dataset to create

Return type None

`lumin.data_processing.file_proc.fold2foldfile` (*df, out_file, fold_idx, cont_feats, cat_feats, targ_feats, targ_type, misc_feats=None, wgt_feat=None*)

Save fold of data into an h5py Group

Parameters

- **df** (DataFrame) – Dataframe from which to save data
- **out_file** (File) – h5py file to save data in
- **fold_idx** (int) – ID for the fold; used name h5py group according to ‘fold_{fold_idx}’
- **cont_feats** (List[str]) – list of columns in df to save as continuous variables
- **cat_feats** (List[str]) – list of columns in df to save as discreet variables
- **targ_feats** (*list of column(s) in df to save as target feature(s)*) –
- **targ_type** (Any) – type of target feature, e.g. int, float32
- **misc_feats** (*optional*) – any extra columns to save
- **wgt_feat** (*optional*) – column to save as data weights

Return type None

`lumin.data_processing.file_proc.df2foldfile` (*df, n_folds, cont_feats, cat_feats, targ_feats, savename, targ_type, strat_key=None, misc_feats=None, wgt_feat=None*)

Convert dataframe into h5py file by splitting data into sub-folds to be accessed by a *FoldYielder*

Parameters

- **df** (DataFrame) – Dataframe from which to save data
- **n_folds** (int) – number of folds to split df into
- **cont_feats** (List[str]) – list of columns in df to save as continuous variables
- **cat_feats** (List[str]) – list of columns in df to save as discrete variables
- **targ_feats** (list of) column(s) in df to save as target feature(s) –
- **savename** (Union[Path, str]) – name of h5py file to create (.h5py extension not required)
- **targ_type** (str) – type of target feature, e.g. int, float32
- **strat_key** (optional) – column to use for stratified splitting
- **misc_feats** (optional) – any extra columns to save
- **wgt_feat** (optional) – column to save as data weights

1.3 lumin.data_processing.hep_proc module

`lumin.data_processing.hep_proc.to_cartesian(df, vec, drop=False)`

Vectorised conversion of 3-momenta to Cartesian coordinates inplace, optionally dropping old pT,eta,phi features

Parameters

- **df** (DataFrame) – DataFrame to alter
- **vec** (str) – column prefix of vector components to alter, e.g. ‘muon’ for columns [‘muon_pt’, ‘muon_phi’, ‘muon_eta’]
- **drop** (bool) – Whether to remove original columns and just keep the new ones

Return type None

`lumin.data_processing.hep_proc.to_pt_eta_phi(df, vec, eta=None, drop=False)`

Vectorised conversion of 3-momenta to pT,eta,phi coordinates inplace, optionally dropping old px,py,pz features

Attention: eta is now deprecated as it is now inferred from df. Will be removed in V0.4

Parameters

- **df** (DataFrame) – DataFrame to alter
- **vec** (str) – column prefix of vector components to alter, e.g. ‘muon’ for columns [‘muon_px’, ‘muon_py’, ‘muon_pz’]
- **eta** (Optional[bool]) – deprecated as now inferred
- **drop** (bool) – Whether to remove original columns and just keep the new ones

Return type None

`lumin.data_processing.hep_proc.delta_phi(arr_a, arr_b)`

Vectorised computation of modulo 2π angular separation of array of angles `b` from array of angles `a`, in range $[-\pi, \pi]$

Parameters

- **arr_a** (Union[float, ndarray]) – reference angles
- **arr_b** (Union[float, ndarray]) – final angles

Return type Union[float, ndarray]

Returns angular separation as float or np.array

`lumin.data_processing.hep_proc.twist(dphi, deta)`

Vectorised computation of twist between vectors (<https://arxiv.org/abs/1010.3698>)

Parameters

- **dphi** (Union[float, ndarray]) – delta phi separations
- **deta** (Union[float, ndarray]) – delta eta separations

Return type Union[float, ndarray]

Returns angular separation as float or np.array

`lumin.data_processing.hep_proc.add_abs_mom(df, vec, z=True)`

Vectorised computation 3-momenta magnitude, adding new column in place. Currently only works for Cartesian vectors

Parameters

- **df** (DataFrame) – DataFrame to alter
- **vec** (str) – column prefix of vector components, e.g. ‘muon’ for columns [‘muon_px’, ‘muon_py’, ‘muon_pz’]
- **z** (bool) – whether to consider the z-component of the momenta

Return type None

`lumin.data_processing.hep_proc.add_mass(df, vec)`

Vectorised computation of mass of 4-vector, adding new column in place.

Parameters

- **df** (DataFrame) – DataFrame to alter
- **vec** (str) – column prefix of vector components, e.g. ‘muon’ for columns [‘muon_px’, ‘muon_py’, ‘muon_pz’]

Return type None

`lumin.data_processing.hep_proc.add_energy(df, vec)`

Vectorised computation of energy of 4-vector, adding new column in place.

Parameters

- **df** (DataFrame) – DataFrame to alter
- **vec** (str) – column prefix of vector components, e.g. ‘muon’ for columns [‘muon_px’, ‘muon_py’, ‘muon_pz’]

Return type None

`lumin.data_processing.hep_proc.add_mt(df, vec, mpt_name='mpt')`

Vectorised computation of transverse mass of 4-vector with respect to missing transverse momenta, adding new column in place. Currently only works for pT, eta, phi vectors

Parameters

- **df** (`DataFrame`) – DataFrame to alter
- **vec** (`str`) – column prefix of vector components, e.g. ‘muon’ for columns [‘muon_px’, ‘muon_py’, ‘muon_pz’]
- **mpt_name** (`str`) – column prefix of vector of missing transverse momenta components, e.g. ‘mpt’ for columns [‘mpt_pT’, ‘mpt_phi’]

`lumin.data_processing.hep_proc.get_vecs(feats, strict=True)`

Filter list of features to get list of 3-momenta defined in the list. Works for both pT, eta, phi and Cartesian coordinates. If strict, return only vectors with all coordinates present in feature list.

Parameters

- **feats** (`List[str]`) – list of features to filter
- **strict** (`bool`) – whether to require all 3-momenta components to be present in the list

Return type `Set[str]`

Returns set of unique 3-momneta prefixes

`lumin.data_processing.hep_proc.fix_event_phi(df, ref_vec)`

Rotate event in phi such that ref_vec is at phi == 0. Performed inplace. Currently only works on vectors defined in pT, eta, phi

Parameters

- **df** (`DataFrame`) – DataFrame to alter
- **ref_vec** (`str`) – column prefix of vector components to use as reference, e.g. ‘muon’ for columns [‘muon_pT’, ‘muon_eta’, ‘muon_phi’]

Return type `None`

`lumin.data_processing.hep_proc.fix_event_z(df, ref_vec)`

Flip event in z-axis such that ref_vec is in positive z-direction. Performed inplace. Works for both pT, eta, phi and Cartesian coordinates.

Parameters

- **df** (`DataFrame`) – DataFrame to alter
- **ref_vec** (`str`) – column prefix of vector components to use as reference, e.g. ‘muon’ for columns [‘muon_pT’, ‘muon_eta’, ‘muon_phi’]

Return type `None`

`lumin.data_processing.hep_proc.fix_event_y(df, ref_vec_0, ref_vec_1)`

Flip event in y-axis such that ref_vec_1 has a higher py than ref_vec_0. Performed in place. Works for both pT, eta, phi and Cartesian coordinates.

Parameters

- **df** (`DataFrame`) – DataFrame to alter
- **ref_vec_0** (`str`) – column prefix of vector components to use as reference 0, e.g. ‘muon’ for columns [‘muon_pT’, ‘muon_eta’, ‘muon_phi’]

- **ref_vec_1** (str) – column prefix of vector components to use as reference 1, e.g. ‘muon’ for columns [‘muon_pT’, ‘muon_eta’, ‘muon_phi’]

Return type None

`lumin.data_processing.hep_proc.event_to_cartesian` (*df*, *drop=False*, *ignore=None*)

Convert entire event to Cartesian coordinates, except vectors listed in *ignore*. Optionally, drop old pT,eta,phi features. Performed inplace.

Parameters

- **df** (DataFrame) – DataFrame to alter
- **drop** (bool) – whether to drop old coordinates
- **ignore** (Optional[List[str]]) – vectors to ignore when converting

Return type None

`lumin.data_processing.hep_proc.proc_event` (*df*, *fix_phi=False*, *fix_y=False*, *fix_z=False*, *use_cartesian=False*, *ref_vec_0=None*, *ref_vec_1=None*, *keep_feats=None*, *default_vals=None*)

Process event: Pass data through inplace various conversions and drop unneeded columns. Data expected to consist of vectors defined in pT, eta, phi.

Parameters

- **df** (DataFrame) – DataFrame to alter
- **fix_phi** (bool) – whether to rotate events using `fix_event_phi()`
- **fix_y** – whether to flip events using `fix_event_y()`
- **fix_z** – whether to flip events using `fix_event_z()`
- **use_cartesian** – whether to convert vectors to Cartesian coordinates
- **ref_vec_0** (Optional[str]) – column prefix of vector components to use as reference (0) for `meth:~lumin.data_processing.hep_proc.fix_event_phi()`, `fix_event_y()`, and `fix_event_z()` e.g. ‘muon’ for columns [‘muon_pT’, ‘muon_eta’, ‘muon_phi’]
- **ref_vec_1** (Optional[str]) – column prefix of vector components to use as reference 1 for `fix_event_z()`, e.g. ‘muon’ for columns [‘muon_pT’, ‘muon_eta’, ‘muon_phi’]
- **keep_feats** (Optional[List[str]]) – columns to keep which would otherwise be dropped
- **default_vals** (Optional[List[str]]) – list of default values which might be used to represent missing vector components. These will be replaced with `np.nan`.

Return type None

`lumin.data_processing.hep_proc.calc_pair_mass` (*df*, *masses*, *feat_map*)

Vectorised computation of invariant mass of pair of particles with given masses, using 3-momenta. Only works for vectors defined in Cartesian coordinates.

Parameters

- **df** (DataFrame) – DataFrame vector components
- **masses** (Union[Tuple[float, float], Tuple[ndarray, ndarray]]) – tuple of masses of particles (either constant or different pair of masses per pair of particles)
- **feat_map** (Dict[str, str]) – dictionary mapping of requested momentum components to the features in *df*

Return type ndarray

Returns np.array of invariant masses

1.4 lumin.data_processing.pre_proc module

```
lumin.data_processing.pre_proc.get_pre_proc_pipes (norm_in=True, norm_out=False,  
                                                  pca=False, whitening=False,  
                                                  with_mean=True, with_std=True,  
                                                  n_components=None)
```

Configure SKLearn Pipelines for processing inputs and targets with the requested transformations.

Parameters

- **norm_in** (bool) – whether to apply StandardScaler to inputs
- **norm_out** (bool) – whether to apply StandardScaler to outputs
- **pca** (bool) – whether to apply PCA to inputs. Performed prior to StandardScaler. No dimensionality reduction is applied, purely rotation.
- **whiten** (bool) – whether PCA should whiten inputs.
- **with_mean** (bool) – whether StandardScalers should shift means to 0
- **with_std** (bool) – whether StandardScalers should scale standard deviations to 1
- **n_components** (Optional[int]) – if set, causes PCA to reduce the dimensionality of the input data

Return type Tuple[Pipeline, Pipeline]

Returns Pipeline for input data Pipeline for target data

```
lumin.data_processing.pre_proc.fit_input_pipe (df, cont_feats, savename=None, in-  
                                              put_pipe=None, norm_in=True,  
                                              pca=False, whitening=False,  
                                              with_mean=True, with_std=True,  
                                              n_components=None)
```

Fit input pipeline to continuous features and optionally save.

Parameters

- **df** (DataFrame) – DataFrame with data to fit pipeline
- **cont_feats** (Union[str, List[str]]) – (list of) column(s) to use as input data for fitting
- **savename** (Optional[str]) – if set will save the fitted Pipeline to with that name as Pickle (.pkl extension added automatically)
- **input_pipe** (Optional[Pipeline]) – if set will fit, otherwise will instantiate a new Pipeline
- **norm_in** (bool) – whether to apply StandardScaler to inputs. Only used if input_pipe is not set.
- **pca** (bool) – whether to apply PCA to inputs. Performed prior to StandardScaler. No dimensionality reduction is applied, purely rotation. Only used if input_pipe is not set.
- **whiten** (bool) – whether PCA should whiten inputs. Only used if input_pipe is not set.

- **with_mean** (bool) – whether StandardScalers should shift means to 0. Only used if input_pipe is not set.
- **with_std** (bool) – whether StandardScalers should scale standard deviations to 1. Only used if input_pipe is not set.
- **n_components** (Optional[int]) – if set, causes PCA to reduce the dimensionality of the input data. Only used if input_pipe is not set.

Return type Pipeline

Returns Fitted Pipeline

`lumin.data_processing.pre_proc.fit_output_pipe(df, targ_feats, savename=None, output_pipe=None, norm_out=True)`

Fit output pipeline to target features and optionally save. Have you thought about using a y_range for regression instead?

Parameters

- **df** (DataFrame) – DataFrame with data to fit pipeline
- **targ_feats** (Union[str, List[str]]) – (list of) column(s) to use as input data for fitting
- **savename** (Optional[str]) – if set will save the fitted Pipeline to with that name as Pickle (.pkl extension added automatically)
- **output_pipe** (Optional[Pipeline]) – if set will fit, otherwise will instantiate a new Pipeline
- **norm_out** (bool) – whether to apply StandardScaler to outputs . Only used if output_pipe is not set.

Return type Pipeline

Returns Fitted Pipeline

`lumin.data_processing.pre_proc.proc_cats(train_df, cat_feats, val_df=None, test_df=None)`

Process categorical features in train_df to be valued 0->cardinality-1. Applied inplace. Applies same transformation to validation and testing data is passed. Will complain if validation or testing sets contain categories which are not present in the training data.

Parameters

- **train_df** (DataFrame) – DataFrame with the training data, which will also be used to specify all the categories to consider
- **cat_feats** (List[str]) – list of columns to use as categorical features
- **val_df** (Optional[DataFrame]) – if set will apply the same category to code mapping to the validation data as was performed on the training data
- **test_df** (Optional[DataFrame]) – if set will apply the same category to code mapping to the testing data as was performed on the training data

Return type Tuple[OrderedDict, OrderedDict]

Returns ordered dictionary mapping categorical features to dictionaries mapping categories to codes
ordered dictionary mapping categorical features to their cardinalities

1.5 Module contents

LUMIN.EVALUATION PACKAGE

2.1 Submodules

2.2 `lumin.evaluation.ams` module

`lumin.evaluation.ams.calc_ams` (*s*, *b*, *br*=0, *unc_b*=0)

Compute Approximate Median Significance (<https://arxiv.org/abs/1007.1727>)

Parameters

- **s** (float) – signal weight
- **b** (float) – background weight
- **br** (float) – background offset bias
- **unc_b** (float) – fractional systematic uncertainty on background

Return type float

Returns Approximate Median Significance if $b > 0$ else -1

`lumin.evaluation.ams.calc_ams_torch` (*s*, *b*, *br*=0, *unc_b*=0)

Compute Approximate Median Significance (<https://arxiv.org/abs/1007.1727>) using Tensor inputs

Parameters

- **s** (Tensor) – signal weight
- **b** (Tensor) – background weight
- **br** (float) – background offset bias
- **unc_b** (float) – fractional systematic uncertainty on background

Return type Tensor

Returns Approximate Median Significance if $b > 0$ else $1e-18 * s$

`lumin.evaluation.ams.ams_scan_quick` (*df*, *wgt_factor*=1, *br*=0, *syst_unc_b*=0,
pred_name='pred', *targ_name*='gen_target',
wgt_name='gen_weight')

Scan across a range of possible prediction thresholds in order to maximise the Approximate Median Significance (<https://arxiv.org/abs/1007.1727>). Note that whilst this method is quicker than `ams_scan_slow()`, it suffers from float precision. Not recommended for final evaluation.

Parameters

- **df** (DataFrame) – DataFrame containing prediction data

- **wgt_factor** (float) – factor to reweight signal and background weights
- **br** (float) – background offset bias
- **syst_unc_b** (float) – fractional systematic uncertainty on background
- **pred_name** (str) – column to use as predictions
- **targ_name** (str) – column to use as truth labels for signal and background
- **wgt_name** (str) – column to use as weights for signal and background events

Return type Tuple[float, float]

Returns maximum AMS prediction threshold corresponding to maximum AMS

```
lumin.evaluation.ams.ams_scan_slow(df, wgt_factor=1, br=0, syst_unc_b=0,
                                   use_stat_unc=False, start_cut=0.9, min_events=10,
                                   pred_name='pred', targ_name='gen_target',
                                   wgt_name='gen_weight', show_prog=True)
```

Scan across a range of possible prediction thresholds in order to maximise the Approximate Median Significance (<https://arxiv.org/abs/1007.1727>). Note that whilst this method is slower than `ams_scan_quick()`, it does not suffer as much from float precision. Additionally it allows one to account for statistical uncertainty in AMS calculation.

Parameters

- **df** (DataFrame) – DataFrame containing prediction data
- **wgt_factor** (float) – factor to reweight signal and background weights
- **br** (float) – background offset bias
- **syst_unc_b** (float) – fractional systematic uncertainty on background
- **use_stat_unc** (bool) – whether to account for the statistical uncertainty on the background
- **start_cut** (float) – minimum prediction to consider; useful for speeding up scan
- **min_events** (int) – minimum number of background unscaled events required to pass threshold
- **pred_name** (str) – column to use as predictions
- **targ_name** (str) – column to use as truth labels for signal and background
- **wgt_name** (str) – column to use as weights for signal and background events
- **show_prog** (bool) – whether to display progress and ETA of scan

Return type Tuple[float, float]

Returns maximum AMS prediction threshold corresponding to maximum AMS

2.3 Module contents

LUMIN.INFERENCE PACKAGE

3.1 Submodules

3.2 `lumin.inference.summary_stat` module

```
lumin.inference.summary_stat.bin_binary_class_pred(df, max_unc, con-  
sider_samples=None, step_sz=0.001, pred_name='pred', sam-  
ple_name='gen_sample', compact_samples=False, class_name='gen_target',  
add_pure_signal_bin=False, max_unc_pure_signal=0.1)
```

Define bin-edges for binning particle process samples as a function of event class prediction (signal | background) such that the statistical uncertainties on per bin yields are below `max_unc` for each considered sample.

Parameters

- **df** (`DataFrame`) – DataFrame containing the data
- **max_unc** (`float`) – maximum fractional statistical uncertainty to allow when defining bins
- **consider_samples** (`Optional[List[str]]`) – if set, only listed samples are considered when defining bins
- **step_sz** (`float`) – resolution of scan along event prediction
- **pred_name** (`str`) – column to use as event class prediction
- **sample_name** (`str`) – column to use as particle process for each event
- **compact_samples** (`bool`) – if true, will not consider samples when computing bin edges, only the class
- **class_name** (`str`) – name of column to use as class indicator
- **add_pure_signal_bin** (`bool`) – if true will attempt to add a bin which only contains signal (class 1) if the fractional bin-fill uncertainty would be less than `max_unc_pure_signal`
- **max_unc_pure_signal** (`float`) – maximum fractional statistical uncertainty to allow when defining pure-signal bins

Return type `List[float]`

Returns list of bin edges

3.3 Module contents

LUMIN.NN PACKAGE

4.1 Subpackages

4.1.1 lumin.nn.callbacks package

Submodules

`lumin.nn.callbacks.callback` module

class `lumin.nn.callbacks.callback.Callback` (*model=None, plot_settings=<lumin.plotting.plot_settings.PlotSettings object>*)

Bases: `lumin.nn.callbacks.abs_callback.AbsCallback`

Base callback class from which other callbacks should inherit.

Parameters

- **model** (`Optional[AbsModel]`) – model to refer to during training
- **plot_settings** (`PlotSettings`) – `PlotSettings` class

set_model (*model*)

Sets the callback's model in order to allow the callback to access and adjust model parameters

Parameters **model** (`AbsModel`) – model to refer to during training

Return type `None`

set_plot_settings (*plot_settings*)

Sets the plot settings for any plots produced by the callback

Parameters **plot_settings** (`PlotSettings`) – `PlotSettings` class

Return type `None`

lumin.nn.callbacks.cyclic_callbacks module

```
class lumin.nn.callbacks.cyclic_callbacks.AbsCyclicCallback (interp,
                                                         param_range,
                                                         cycle_mult=1, decrease_param=False,
                                                         scale=1,
                                                         model=None,
                                                         nb=None,
                                                         plot_settings=<lumin.plotting.plot_settings.P
                                                         object>)
```

Bases: *lumin.nn.callbacks.callback.Callback*

Abstract class for callbacks affecting lr or mom

Parameters

- **interp** (*str*) – string representation of interpolation function. Either ‘linear’ or ‘cosine’.
- **param_range** (*Tuple[float, float]*) – minimum and maximum values for parameter
- **cycle_mult** (*int*) – multiplicative factor for adjusting the cycle length after each cycle. E.g *cycle_mult=1* keeps the same cycle length, *cycle_mult=2* doubles the cycle length after each cycle.
- **decrease_param** (*bool*) – whether to begin by decreasing the parameter, otherwise begin by increasing it
- **scale** (*int*) – multiplicative factor for setting the initial number of epochs per cycle. E.g *scale=1* means 1 epoch per cycle, *scale=5* means 5 epochs per cycle.
- **model** (*Optional[AbsModel]*) – model to refer to during training
- **nb** (*Optional[int]*) – number of minibatches (iterations) to expect per epoch
- **plot_settings** (*PlotSettings*) – *PlotSettings* class

on_batch_begin (***kargs*)

Computes the new value for the optimiser parameter and returns it

Return type *float*

Returns new value for optimiser parameter

on_batch_end (***kargs*)

Increments the callback’s progress through the cycle

Return type *None*

on_epoch_begin (***kargs*)

Ensures the *cycle_end* flag is false when the epoch starts

Return type *None*

plot ()

Plots the history of the parameter evolution as a function of iterations

Return type *None*

set_nb (*nb*)

Sets the callback’s internal number of iterations per cycle equal to *nb*scale*

Parameters **nb** (*int*) – number of minibatches per epoch

Return type *None*

```
class lumin.nn.callbacks.cyclic_callbacks.CycleLR (lr_range,          interp='cosine',
                                                  cycle_mult=1,          de-
                                                  crease_param='auto',    scale=1,
                                                  model=None,             nb=None,
                                                  plot_settings=<lumin.plotting.plot_settings.PlotSettings
                                                  object>)
```

Bases: *lumin.nn.callbacks.cyclic_callbacks.AbsCyclicCallback*

Callback to cycle learning rate during training according to either: cosine interpolation for SGDR <https://arxiv.org/abs/1608.03983> or linear interpolation for Smith cycling <https://arxiv.org/abs/1506.01186>

Parameters

- **lr_range** (Tuple[float, float]) – tuple of initial and final LRs
- **interp** (str) – ‘cosine’ or ‘linear’ interpolation
- **cycle_mult** (int) – Multiplicative constant for altering the cycle length after each complete cycle
- **decrease_param** (Union[str, bool]) – whether to increase or decrease the LR (effectively reverses lr_range order), ‘auto’ selects according to interp
- **scale** (int) – Multiplicative constant for altering the length of a cycle. 1 corresponds to one cycle = one (sub-)epoch
- **model** (Optional[AbsModel]) – *Model* to alter, alternatively call `set_model()`.
- **nb** (Optional[int]) – Number of batches in a (sub-)epoch
- **plot_settings** (*PlotSettings*) – *PlotSettings* class to control figure appearance

Examples::

```
>>> cosine_lr = CycleLR(lr_range=(0, 2e-3), cycle_mult=2, scale=1,
...                    interp='cosine', nb=100)
>>>
>>> cyclical_lr = CycleLR(lr_range=(2e-4, 2e-3), cycle_mult=1, scale=5,
...                      interp='linear', nb=100)
```

on_batch_begin (**kargs)

Computes the new lr and assigns it to the optimiser

Return type None

```
class lumin.nn.callbacks.cyclic_callbacks.CycleMom (mom_range,      interp='cosine',
                                                  cycle_mult=1,          de-
                                                  crease_param='auto',    scale=1,
                                                  model=None,             nb=None,
                                                  plot_settings=<lumin.plotting.plot_settings.PlotSettings
                                                  object>)
```

Bases: *lumin.nn.callbacks.cyclic_callbacks.AbsCyclicCallback*

Callback to cycle momentum (beta 1) during training according to either: cosine interpolation for SGDR <https://arxiv.org/abs/1608.03983> or linear interpolation for Smith cycling <https://arxiv.org/abs/1506.01186> By default is set to evolve in opposite direction to learning rate, a la <https://arxiv.org/abs/1803.09820>

Parameters

- **mom_range** (Tuple[float, float]) – tuple of initial and final momenta

- **interp** (*str*) – ‘cosine’ or ‘linear’ interpolation
- **cycle_mult** (*int*) – Multiplicative constant for altering the cycle length after each complete cycle
- **decrease_param** (*Union[str, bool]*) – whether to increase or decrease the momentum (effectively reverses *mom_range* order), ‘auto’ selects according to *interp*
- **scale** (*int*) – Multiplicative constant for altering the length of a cycle. 1 corresponds to one cycle = one (sub-)epoch
- **model** (*Optional[AbsModel]*) – *Model* to alter, alternatively call *set_model()*
- **nb** (*Optional[int]*) – Number of batches in a (sub-)epoch
- **plot_settings** (*PlotSettings*) – *PlotSettings* class to control figure appearance

Examples::

```
>>> cyclical_mom = CycleMom(mom_range=(0.85 0.95), cycle_mult=1,
...                          scale=5, interp='linear', nb=100)
```

on_batch_begin (***kargs*)

Computes the new momentum and assigns it to the optimiser

Return type *None*

```
class lumin.nn.callbacks.cyclic_callbacks.OneCycle (lengths, lr_range,
...                                               mom_range=(0.85,
...                                               0.95), interp='cosine',
...                                               model=None, nb=None,
...                                               plot_settings=<lumin.plotting.plot_settings.PlotSettings
...                                               object>)
```

Bases: *lumin.nn.callbacks.cyclic_callbacks.AbsCyclicCallback*

Callback implementing Smith 1-cycle evolution for lr and momentum (*beta_1*) <https://arxiv.org/abs/1803.09820>
Default interpolation uses fastai-style cosine function. Automatically triggers early stopping on cycle completion.

Parameters

- **lengths** (*Tuple[int, int]*) – tuple of number of (sub-)epochs in first and second stages of cycle
- **lr_range** (*List[float]*) – tuple of initial and final LRs
- **mom_range** (*Tuple[float, float]*) – tuple of initial and final momenta
- **interp** (*str*) – ‘cosine’ or ‘linear’ interpolation
- **model** (*Optional[AbsModel]*) – *Model* to alter, alternatively call *set_model()*
- **nb** (*Optional[int]*) – Number of batches in a (sub-)epoch
- **plot_settings** (*PlotSettings*) – *PlotSettings* class to control figure appearance

Examples::

```
>>> onecycle = OneCycle(lengths=(15, 30), lr_range=[1e-4, 1e-2],
...                  mom_range=(0.85, 0.95), interp='cosine', nb=100)
```

on_batch_begin (**kargs)

Computes the new lr and momentum and assigns them to the optimiser

Return type None

plot ()

Plots the history of the lr and momentum evolution as a function of iterations

lumin.nn.callbacks.data_callbacks module

class lumin.nn.callbacks.data_callbacks.**BinaryLabelSmooth** (coefs=0, model=None)

Bases: *lumin.nn.callbacks.callback.Callback*

Callback for applying label smoothing to binary classes, based on <https://arxiv.org/abs/1512.00567> Applies smoothing during both training and inference.

Parameters

- **coefs** (Union[float, Tuple[float, float]]) – Smoothing coefficients: 0->coef[0] 1->1-coef[1]. if passed float, coef[0]=coef[1]
- **model** (Optional[AbsModel]) – not used, only for compatability

Examples::

```
>>> lbl_smooth = BinaryLabelSmooth(0.1)
>>>
>>> lbl_smooth = BinaryLabelSmooth((0.1, 0.02))
```

on_epoch_begin (by, **kargs)

Apply smoothing at train-time

Return type None

on_eval_begin (targets, **kargs)

Apply smoothing at test-time

Return type None

class lumin.nn.callbacks.data_callbacks.**SequentialReweight** (reweight_func,

scale=0.1,

model=None)

Bases: *lumin.nn.callbacks.callback.Callback*

Caution: Experiemntal proceedure

During ensemble training, sequentially reweight training data in last validation fold based on prediction performance of last trained model. Reweighting highlights data which are easier or more difficult to predict to the next model being trained.

Parameters

- **reweight_func** (Callable[[Tensor, Tensor], Tensor]) – callable function returning a tensor of same shape as targets, ideally quantifying model-prediction performance
- **scale** (float) – multiplicative factor for rescaling returned tensor of reweight_func
- **model** (Optional[AbsModel]) – *Model* to provide predictions, alternatively call `set_model()`

Examples::

```
>>> seq_reweight = SequentialReweight (
...     reweight_func=nn.BCELoss(reduction='none'), scale=0.1)
```

on_train_end (*fy, val_id, **kargs*)

Reweights the validation fold once training is finished

Parameters

- **fy** (*FoldYielder*) – FoldYielder providing the training and validation data
- **fold_id** – Fold index which was used for validation

Return type None

```
class lumin.nn.callbacks.data_callbacks.SequentialReweightClasses (reweight_func,
...                                     scale=0.1,
...                                     model=None)
```

Bases: *lumin.nn.callbacks.data_callbacks.SequentialReweight*

Caution: Experimental procedure

Version of *SequentialReweight* designed for classification, which renormalises class weights to original weight-sum after reweighting. During ensemble training, sequentially reweight training data in last validation fold based on prediction performance of last trained model. Reweighting highlights data which are easier or more difficult to predict to the next model being trained.

Parameters

- **reweight_func** (Callable[[Tensor, Tensor], Tensor]) – callable function returning a tensor of same shape as targets, ideally quantifying model-prediction performance
- **scale** (float) – multiplicative factor for rescaling returned tensor of *reweight_func*
- **model** (Optional[AbsModel]) – *Model* to provide predictions, alternatively call *set_model()*

Examples::

```
>>> seq_reweight = SequentialReweight (
...     reweight_func=nn.BCELoss(reduction='none'), scale=0.1)
```

```
class lumin.nn.callbacks.data_callbacks.BootstrapResample (n_folds,
...                                     bag_each_time=False,
...                                     reweight=True,
...                                     model=None)
```

Bases: *lumin.nn.callbacks.callback.Callback*

Callback for bootstrap sampling new training datasets from original training data during (ensemble) training.

Parameters

- **n_folds** (int) – the number of folds present in training *FoldYielder*
- **bag_each_time** (bool) – whether to sample a new set for each sub-epoch or to use the same sample each time
- **reweight** (bool) – whether to reweight the sampled data to match the weight sum (per class) of the original data

- **model** (Optional[AbsModel]) – not used, only for compatability

Examples::

```
>>> bs_resample BootstrapResample(n_folds=len(train_fy))
```

on_epoch_begin (*by*, ***kargs*)

Resamples training data for new epoch

Parameters by (*BatchYielder*) – BatchYielder providing data for the upcoming epoch

Return type None

on_train_begin (***kargs*)

Resets internal parameters to prepare for a new training

Return type None

class lumin.nn.callbacks.data_callbacks.**FeatureSubsample** (*cont_feats*,
model=None)

Bases: *lumin.nn.callbacks.callback.Callback*

Callback for training a model on a random sub-sample of the range of possible input features. Only sub-samples continuous features. Number of continuous inputs inferred from model. Associated *Model* will automatically mask its inputs during inference; simply provide inputs with the same number of columns as training data.

Parameters

- **cont_feats** (List[str]) – list of all continuous features in input data. Order must match.
- **model** (Optional[AbsModel]) – *Model* being trained, alternatively call `set_model()`

Examples::

```
>>> feat_subsample = FeatureSubsample(cont_feats=['pT', 'eta', 'phi'])
```

on_epoch_begin (*by*, ***kargs*)

Masks input data to remove non-selected features

Parameters by (*BatchYielder*) – BatchYielder providing data for the upcoming epoch

Return type None

on_train_begin (***kargs*)

Subsamples features for use in training and sets model's input mask for inference

Return type None

lumin.nn.callbacks.loss_callbacks module

class lumin.nn.callbacks.loss_callbacks.**GradClip** (*clip*, *clip_norm=True*, *model=None*)

Bases: *lumin.nn.callbacks.callback.Callback*

Callback for clipping gradients by norm or value.

Parameters

- **clip** (float) – value to clip at

- **clip_norm** (bool) – whether to clip according to norm (*torch.nn.utils.clip_grad_norm_*) or value (*torch.nn.utils.clip_grad_value_*)
- **model** (Optional[AbsModel]) – *Model* with parameters to clip gradients, alternatively call `set_model()`

Examples::

```
>>> grad_clip = GradClip(1e-5)
```

on_backwards_end (**kargs)

Clips gradients prior to parameter updates

Return type None

lumin.nn.callbacks.model_callbacks module

```
class lumin.nn.callbacks.model_callbacks.SWA (start_epoch, renewal_period=-1, model=None, val_fold=None, cyclic_callback=None, update_on_cycle_end=None, verbose=False, plot_settings=<lumin.plotting.plot_settings.PlotSettings object>)
```

Bases: *lumin.nn.callbacks.model_callbacks.AbsModelCallback*

Callback providing Stochastic Weight Averaging based on (<https://arxiv.org/abs/1803.05407>) This adapted version allows the tracking of a pair of average models in order to avoid having to hardcode a specific start point for averaging:

- Model average x0 will begin to be tracked start_epoch (sub-)epochs/cycles after training begins.
- *cycle_since_replacement* is set to 1
- Renewal_period (sub-)epochs/cycles later, a second average x1 will be tracked.
- At the next renewal period, the performance of x0 and x1 will be compared on data contained in val_fold.
 - **If x0 is better than x1:**
 - * x1 is replaced by a copy of the current model
 - * *cycle_since_replacement* is increased by 1
 - * *renewal_period* is multiplied by *cycle_since_replacement*
 - **Else:**
 - * x0 is replaced by x1
 - * x1 is replaced by a copy of the current model
 - * *cycle_since_replacement* is set to 1
 - * *renewal_period* is set back to its original value

Additionally, will optionally (default True) lock-in to any cyclical callbacks to only update at the end of a cycle.

Parameters

- **start_epoch** (int) – (sub-)epoch/cycle to begin averaging
- **renewal_period** (int) – How often to check performance of averages, and renew tracking of least performant

- **model** (Optional[AbsModel]) – *Model* to provide parameters, alternatively call `set_model()`
- **val_fold** (Optional[Dict[str, ndarray]]) – Dictionary containing inputs, targets, and weights (or None) as Numpy arrays
- **cyclic_callback** (Optional[AbsCyclicCallback]) – Optional for any cyclical callback which is running
- **update_on_cycle_end** (Optional[bool]) – Whether to lock in to the cyclic callback and only update at the end of a cycle. Default yes, if cyclic callback present.
- **verbose** (bool) – Whether to print out update information for testing and operation confirmation
- **plot_settings** (*PlotSettings*) – *PlotSettings* class to control figure appearance

Examples::

```
>>> swa = SWA(start_epoch=5, renewal_period=5)
```

get_loss()

Evaluates SWA model and returns loss

Return type float

Returns Loss on validation fold for oldest SWA average

on_epoch_begin (kargs)**

Resets loss to prepare for new epoch

Return type None

on_epoch_end (kargs)**

Checks whether averages should be updated (or reset) and increments counters

Return type None

on_train_begin (kargs)**

Initialises model variables to begin tracking new model averages

Return type None

```
class lumin.nn.callbacks.model_callbacks.AbsModelCallback (model=None,
                                                         val_fold=None,
                                                         cyclic_callback=None,
                                                         up-
                                                         date_on_cycle_end=None,
                                                         plot_settings=<lumin.plotting.plot_settings.Plot
                                                         object>)
```

Bases: *lumin.nn.callbacks.callback.Callback*

Abstract class for callbacks which provide alternative models during training

Parameters

- **model** (Optional[AbsModel]) – *Model* to provide parameters, alternatively call `set_model()`
- **val_fold** (Optional[Dict[str, ndarray]]) – Dictionary containing inputs, targets, and weights (or None) as Numpy arrays

- **cyclic_callback** (Optional[AbsCyclicCallback]) – Optional for any cyclical callback which is running
- **update_on_cycle_end** (Optional[bool]) – Whether to lock in to the cyclic callback and only update at the end of a cycle. Default yes, if cyclic callback present.
- **plot_settings** (PlotSettings) – PlotSettings class to control figure appearance

abstract get_loss ()

Return type float

set_cyclic_callback (cyclic_callback)

Sets the cyclical callback to lock into for updating new models

Return type None

set_val_fold (val_fold)

Sets the validation fold used for evaluating new models

Return type None

lumin.nn.callbacks.opt_callbacks module

```
class lumin.nn.callbacks.opt_callbacks.LRFinder (nb, lr_bounds=[1e-07, 10], model=None, plot_settings=<lumin.plotting.plot_settings.PlotSettings object>)
```

Bases: *lumin.nn.callbacks.callback.Callback*

Callback class for Smith learning-rate range test (<https://arxiv.org/abs/1803.09820>)

Parameters

- **nb** (int) – number of batches in a (sub-)epoch
- **lr_bounds** (Tuple[float, float]) – tuple of initial and final LR
- **model** (Optional[AbsModel]) – Model to alter, alternatively call `set_model()`
- **plot_settings** (PlotSettings) – PlotSettings class to control figure appearance

on_batch_end (loss, **kargs)

Records loss and increments LR

Parameters **loss** (float) – training loss for most recent batch

on_train_begin (**kargs)

Prepares variables and optimiser for new training

plot (n_skip=0, n_max=None, lim_y=None)

Plot the loss as a function of the LR.

Parameters

- **n_skip** (int) – Number of initial iterations to skip in plotting
- **n_max** (Optional[int]) – Maximum iteration number to plot
- **lim_y** (Optional[Tuple[float, float]]) – y-range for plotting

plot_lr ()

Plot the LR as a function of iterations.

Module contents

4.1.2 lumin.nn.data package

Submodules

`lumin.nn.data.batch_yielder` module

```
class lumin.nn.data.batch_yielder.BatchYielder (inputs, targets, bs, objective,  
weights=None, shuffle=True,  
use_weights=True, bulk_move=True)
```

Bases: object

Yields minibatches to model during training. Iteration provides one minibatch as tuple of tensors of inputs, targets, and weights.

Parameters

- **inputs** (ndarray) – input array for (sub-)epoch
- **targets** (ndarray) – target array for (sub-)epoch
- **bs** (int) – batchsize, number of data to include per minibatch
- **objective** (str) – ‘classification’, ‘multiclass classification’, or ‘regression’. Used for casting target dtype.
- **weights** (Optional[ndarray]) – Optional weight array for (sub-)epoch
- **shuffle** – whether to shuffle the data at the beginning of an iteration
- **use_weights** (bool) – if passed weights, whether to actually pass them to the model
- **bulk_move** – whether to move all data to device at once. Default is true (saves time), but if device has low memory you can set to False.

`lumin.nn.data.fold_yielder` module

```
class lumin.nn.data.fold_yielder.FoldYielder (foldfile, cont_feats, cat_feats, ig-  
nore_feats=None, input_pipe=None,  
output_pipe=None)
```

Bases: object

Interface class for accessing data from foldfiles created by `df2foldfile()`

Parameters

- **foldfile** (File) – filename of hdf5 file
- **cont_feats** (List[str]) – list of names of continuous features present in input data
- **cat_feats** (List[str]) – list of names of categorical features present in input data
- **ignore_feats** (Optional[List[str]]) – optional list of input features which should be ignored
- **input_pipe** (Union[str, Pipeline, None]) – optional Pipeline, or filename for pickled Pipeline, which was used for processing the inputs
- **output_pipe** (Union[str, Pipeline, None]) – optional Pipeline, or filename for pickled Pipeline, which was used for processing the targets

Examples::

```
>>> fy = FoldYielder('train.h5', cont_feats=['pT','eta','phi','mass'],
...                 cat_feats=['channel'], ignore_feats=['phi'],
...                 input_pipe='input_pipe.pkl')
```

add_ignore (*feats*)

Add features to ignored features.

Parameters **feats** (List[str]) – list of feature names to ignore

Return type None

add_input_pipe (*input_pipe*)

Adds an input pipe to the FoldYielder for use when deprocessing data

Parameters **input_pipe** (Pipeline) – Pipeline which was used for preprocessing the input data

Return type None

add_input_pipe_from_file (*name*)

Adds an input pipe from a pkl file to the FoldYielder for use when deprocessing data

Parameters **name** (str) – name of pkl file containing Pipeline which was used for preprocessing the input data

Return type None

add_output_pipe (*output_pipe*)

Adds an output pipe to the FoldYielder for use when deprocessing data

Parameters **output_pipe** (Pipeline) – Pipeline which was used for preprocessing the target data

Return type None

add_output_pipe_from_file (*name*)

Adds an output pipe from a pkl file to the FoldYielder for use when deprocessing data

Parameters **name** (str) – name of pkl file containing Pipeline which was used for preprocessing the target data

Return type None

get_column (*column, n_folds=None, fold_idx=None, add_newaxis=False*)

Load column (h5py group) from foldfile. Used for getting arbitrary data which isn't automatically grabbed by other methods.

Parameters

- **column** (str) – name of h5py group to get
- **n_folds** (Optional[int]) – number of folds to get data from. Default all folds. Not compatible with `fold_idx`
- **fold_idx** (Optional[int]) – Only load group from a single, specified fold. Not compatible with `n_folds`
- **add_newaxis** (bool) – whether expand shape of returned data if data shape is ()

Return type Optional[ndarray]

Returns Numpy array of column data

get_data (*n_folds=None, fold_idx=None*)

Get data for single, specified fold or several of folds. Data consists of dictionary of inputs, targets, and weights. Does not accounts for ignored features. Inputs are passed through `np.nan_to_num` to deal with nans and infs.

Parameters

- **n_folds** (Optional[int]) – number of folds to get data from. Default all folds. Not compatible with `fold_idx`
- **fold_idx** (Optional[int]) – Only load group from a single, specified fold. Not compatible with `n_folds`

Return type Dict[str, ndarray]

Returns tuple of inputs, targets, and weights as Numpy arrays

get_df (*pred_name='pred', targ_name='targets', wgt_name='weights', n_folds=None, fold_idx=None, inc_inputs=False, inc_ignore=False, deprocess=False, verbose=True, suppress_warn=False*)

Get a Pandas DataFrame of the data in the foldfile. Will add columns for inputs (if requested), targets, weights, and predictions (if present)

Parameters

- **pred_name** (str) – name of prediction group
- **targ_name** (str) – name of target group
- **wgt_name** (str) – name of weight group
- **n_folds** (Optional[int]) – number of folds to get data from. Default all folds. Not compatible with `fold_idx`
- **fold_idx** (Optional[int]) – Only load group from a single, specified fold. Not compatible with `n_folds`
- **inc_inputs** (bool) – whether to include input data
- **inc_ignore** (bool) – whether to include ignored features
- **deprocess** (bool) – whether to deprocess inputs and targets if pipelines have been
- **verbose** (bool) – whether to print the number of datapoints loaded
- **suppress_warn** (bool) – whether to suppress the warning about missing columns

Return type DataFrame

Returns Pandas DataFrame with requested data

get_fold (*idx*)

Get data for single fold. Data consists of dictionary of inputs, targets, and weights. Accounts for ignored features. Inputs are passed through `np.nan_to_num` to deal with nans and infs.

Parameters **idx** (int) – fold index to load

Return type Dict[str, ndarray]

Returns tuple of inputs, targets, and weights as Numpy arrays

get_ignore ()

Returns list of ignored features

Return type List[str]

Returns Features removed from training data

save_fold_pred (*pred*, *fold_idx*, *pred_name*='pred')

Save predictions for given fold as a new column in the foldfile

Parameters

- **pred** (ndarray) – array of predictions in the same order as data appears in the file
- **fold_idx** (int) – index for fold
- **pred_name** (str) – name of column to save predictions under

Return type None

set_foldfile (*foldfile*)

Sets the file from which to access data

Parameters **foldfile** (File) – opened h5py file

Return type None

```
class lumin.nn.data.fold_yielder.HEPAugFoldYielder (foldfile, cont_feats, cat_feats,
                                                    ignore_feats=None,
                                                    targ_feats=None, rot_mult=2,
                                                    random_rot=False, re-
                                                    flect_x=False, re-
                                                    flect_y=True, reflect_z=True,
                                                    train_time_aug=True,
                                                    test_time_aug=True, in-
                                                    put_pipe=None, out-
                                                    put_pipe=None)
```

Bases: *lumin.nn.data.fold_yielder.FoldYielder*

Specialised version of *FoldYielder* providing HEP specific data augmentation at train and test time.

Parameters

- **foldfile** (File) – filename of hdf5 file
- **cont_feats** (List[str]) – list of names of continuous features present in input data
- **cat_feats** (List[str]) – list of names of categorical features present in input data
- **ignore_feats** (Optional[List[str]]) – optional list of input features which should be ignored
- **targ_feats** (Optional[List[str]]) – optional list of target features to also be transformed
- **rot_mult** (int) – number of rotations of event in phi to make at test-time (currently must be even). Greater than zero will also apply random rotations during train-time
- **random_rot** (bool) – whether test-time rotation angles should be random or in steps of $2\pi/\text{rot_mult}$
- **reflect_x** (bool) – whether to reflect events in x axis at train and test time
- **reflect_y** (bool) – whether to reflect events in y axis at train and test time
- **reflect_z** (bool) – whether to reflect events in z axis at train and test time
- **train_time_aug** (bool) – whether to apply augmentations at train time
- **test_time_aug** (bool) – whether to apply augmentations at test time
- **input_pipe** (Optional[Pipeline]) – optional Pipeline, or filename for pickled Pipeline, which was used for processing the inputs

- **output_pipe** (Optional[Pipeline]) – optional Pipeline, or filename for pickled Pipeline, which was used for processing the targets

Examples::

```
>>> fy = HEPAugFoldYielder('train.h5',
...                          cont_feats=['pT', 'eta', 'phi', 'mass'],
...                          rot_mult=2, reflect_y=True, reflect_z=True,
...                          input_pipe='input_pipe.pkl')
```

get_fold (*idx*)

Get data for single fold applying random train-time data augmentation. Data consists of dictionary of inputs, targets, and weights. Accounts for ignored features. Inputs are passed through `np.nan_to_num` to deal with nans and infs.

Parameters *idx* (int) – fold index to load

Return type Dict[str, ndarray]

Returns tuple of inputs, targets, and weights as Numpy arrays

get_test_fold (*idx*, *aug_idx*)

Get test data for single fold applying test-time data augmentation. Data consists of dictionary of inputs, targets, and weights. Accounts for ignored features. Inputs are passed through `np.nan_to_num` to deal with nans and infs.

Parameters

- *idx* (int) – fold index to load
- *aug_idx* (int) – index for the test-time augmentation (ignored if random test-time augmentation requested)

Return type Dict[str, ndarray]

Returns tuple of inputs, targets, and weights as Numpy arrays

Module contents**4.1.3 lumin.nn.ensemble package****Submodules****lumin.nn.ensemble.ensemble module**

class `lumin.nn.ensemble.ensemble.Ensemble` (*input_pipe=None*, *output_pipe=None*,
model_builder=None)
Bases: `lumin.nn.ensemble.abs_ensemble.AbsEnsemble`

Standard class for building an ensemble of collection of trained networks produced by `fold_train_ensemble()`. Input and output pipelines can be added to provide easy saving and loading of exported ensembles. Currently, the input pipeline is not used, so input data is expected to be preprocessed. However the output pipeline will be used to deprocess model predictions.

Once instantiated, `lumin.nn.ensemble.ensemble.Ensemble.build_ensemble()` or `:meth:load` should be called. Alternatively, class methods `lumin.nn.ensemble.ensemble.Ensemble.from_save()` or `lumin.nn.ensemble.ensemble.Ensemble.from_results()` may be used.

Parameters

- **input_pipe** (Optional[Pipeline]) – Optional input pipeline, alternatively call `lumin.nn.ensemble.ensemble.Ensemble.add_input_pipe()`
- **output_pipe** (Optional[Pipeline]) – Optional output pipeline, alternatively call `lumin.nn.ensemble.ensemble.Ensemble.add_output_pipe()`
- **model_builder** (Optional[*ModelBuilder*]) – Optional *ModelBuilder* for constructing models from saved weights.

Examples::

```
>>> ensemble = Ensemble()
>>>
>>> ensemble = Ensemble(input_pipe, output_pipe, model_builder)
```

add_input_pipe (*pipe*)

Add input pipeline for saving

Parameters **pipe** (Pipeline) – pipeline used for preprocessing input data**Return type** None**add_output_pipe** (*pipe*)

Add output pipeline for saving

Parameters **pipe** (Pipeline) – pipeline used for preprocessing target data**Return type** None

build_ensemble (*results*, *size*, *model_builder*, *metric='loss'*, *weighting='reciprocal'*, *higher_metric_better=False*, *snapshot_args=None*, *location=PosixPath('train_weights')*, *verbose=True*)

Load up an instantiated *Ensemble* with outputs of `fold_train_ensemble()`**Parameters**

- **results** (List[Dict[str, float]]) – results saved/returned by `fold_train_ensemble()`
- **size** (int) – number of models to load as ranked by metric
- **model_builder** (*ModelBuilder*) – *ModelBuilder* used for building *Model* from saved models
- **metric** (str) – metric name listed in results to use for ranking and weighting trained models
- **weighting** (str) – ‘reciprocal’ or ‘uniform’ how to weight model predictions during prediction. ‘reciprocal’ = models weighted by 1/metric ‘uniform’ = models treated with equal weighting
- **higher_metric_better** (bool) – whether metric should be maximised or minimised
- **snapshot_args** (Optional[Dict[str, Any]]) – Dictionary potentially containing: ‘cycle_losses’: returned/save by `fold_train_ensemble()` when using an *AbsCyclicCallback* ‘patience’: patience value that was passed to `fold_train_ensemble()` ‘n_cycles’: number of cycles to load per model ‘load_cycles_only’: whether to only load cycles, or also the best performing model ‘weighting_pwr’: weight cycles according to $(n+1)^{*}weighting_pwr$, where n is the number of cycles loaded so far.

Models are loaded youngest to oldest

- **location** (`Path`) – Path to save location passed to `fold_train_ensemble()`
- **verbose** (`bool`) – whether to print out information of models loaded

Examples::

```

>>> ensemble.build_ensemble(results, 10, model_builder,
...                          location=Path('train_weights'))
>>>
>>> ensemble.build_ensemble(
...     results, 1, model_builder,
...     location=Path('train_weights'),
...     snapshot_args={'cycle_losses':cycle_losses,
...                    'patience':patience,
...                    'n_cycles':8,
...                    'load_cycles_only':True,
...                    'weighting_pwr':0})

```

Return type `None`

export2onnx (`base_name`, `bs=1`)

Export all *Model* contained in *Ensemble* to ONNX format. Note that ONNX expects a fixed batch size (bs) which is the number of datapoints your wish to pass through the model concurrently.

Parameters

- **base_name** (`str`) – Exported models will be called `{base_name}_{model_num}.onnx`
- **bs** (`int`) – batch size for exported models

Return type `None`

export2tfpb (`base_name`, `bs=1`)

Export all *Model* contained in *Ensemble* to Tensorflow ProtocolBuffer format, via ONNX. Note that ONNX expects a fixed batch size (bs) which is the number of datapoints your wish to pass through the model concurrently.

Parameters

- **base_name** (`str`) – Exported models will be called `{base_name}_{model_num}.pb`
- **bs** (`int`) – batch size for exported models

Return type `None`

classmethod from_results (`results`, `size`, `model_builder`, `metric='loss'`, `weighting='reciprocal'`, `higher_metric_better=False`, `snapshot_args=None`, `location=PosixPath('train_weights')`, `verbose=True`)

Instantiate *Ensemble* from a outputs of `fold_train_ensemble()`. If cycle models are loaded, then only uniform weighting between models is supported.

Parameters

- **results** (`List[Dict[str, float]]`) – results saved/returned by `fold_train_ensemble()`
- **size** (`int`) – number of models to load as ranked by metric
- **model_builder** (`ModelBuilder`) – *ModelBuilder* used for building *Model* from saved models
- **metric** (`str`) – metric name listed in results to use for ranking and weighting trained models

- **weighting** (str) – ‘reciprocal’ or ‘uniform’ how to weight model predictions during prediction. ‘reciprocal’ = models weighted by 1/metric ‘uniform’ = models treated with equal weighting
- **higher_metric_better** (bool) – whether metric should be maximised or minimised
- **snapshot_args** (Optional[Dict[str, Any]]) – Dictionary potentially containing: ‘cycle_losses’: returned/save by `fold_train_ensemble()` when using an `AbsCyclicCallback` ‘patience’: patience value that was passed to `fold_train_ensemble()` ‘n_cycles’: number of cycles to load per model ‘load_cycles_only’: whether to only load cycles, or also the best performing model ‘weighting_pwr’: weight cycles according to $(n+1)**weighting_pwr$, where n is the number of cycles loaded so far.

Models are loaded youngest to oldest

- **location** (Path) – Path to save location passed to `fold_train_ensemble()`
- **verbose** (bool) – whether to print out information of models loaded

Return type AbsEnsemble

Returns Built `Ensemble`

Examples::

```
>>> ensemble = Ensemble.from_results(results, 10, model_builder,
...                                 location=Path('train_weights'))
>>>
>>> ensemble = Ensemble.from_results(
...     results, 1, model_builder,
...     location=Path('train_weights'),
...     snapshot_args={'cycle_losses':cycle_losses,
...                    'patience':patience,
...                    'n_cycles':8,
...                    'load_cycles_only':True,
...                    'weighting_pwr':0})
```

classmethod `from_save` (name)

Instantiate `Ensemble` from a saved `Ensemble`

Parameters `name` (str) – base filename of ensemble

Return type AbsEnsemble

Returns Loaded `Ensemble`

Examples::

```
>>> ensemble = Ensemble.from_save('weights/ensemble')
```

get_feat_importance (fy, eval_metric=None)

Call `get_ensemble_feat_importance()`, passing this `Ensemble` and provided arguments

Parameters

- **fy** (`FoldYielder`) – `FoldYielder` interfacing to data on which to evaluate importance

- **eval_metric** (Optional[*EvalMetric*]) – Optional *EvalMetric* to use for quantifying performance

Return type *DataFrame*

load (*name*)

Load an instantiated *Ensemble* with weights and *Model* from save.

Arguments; *name*: base name for saved objects

Examples::

```
>>> ensemble.load('weights/ensemble')
```

Return type *None*

static load_trained_model (*model_idx*, *model_builder*, *name*='train_weights/train_')

Load trained model from save file of the form {*name*}{*model_idx*}.h5

Arguments *model_idx*: index of model to load *model_builder*: *ModelBuilder* used to build the model
name: base name of file from which to load model

Return type *Model*

Returns Model loaded from save

predict (*inputs*, *n_models*=None, *pred_name*='pred')

Compatibility method for predicting data contained in either a Numpy array or a *FoldYielder*. Will either pass inputs to `lumin.nn.ensemble.ensemble.Ensemble.predict_array()` or `lumin.nn.ensemble.ensemble.Ensemble.predict_folds()`.

Parameters

- **inputs** (Union[*ndarray*, *FoldYielder*, List[*ndarray*]]) – either a *FoldYielder* interfacing with the input data, or the input data as an array
- **n_models** (Optional[int]) – number of models to use in predictions as ranked by the metric which was used when constructing the *Ensemble*. By default, entire ensemble is used.
- **pred_name** (str) – name for new group of predictions if passed a *FoldYielder*

Return type Union[None, *ndarray*]

Returns If passed a Numpy array will return predictions.

Examples::

```
>>> preds = ensemble.predict(input_array)
>>>
>>> ensemble.predict(test_fy)
```

predict_array (*arr*, *n_models*=None, *parent_bar*=None, *display*=True)

Apply ensemble to Numpy array and get predictions. If an output pipe has been added to the ensemble, then the predictions will be deprocessed. Inputs are expected to be preprocessed; i.e. any input pipe added to the ensemble is not used.

Parameters

- **arr** (*ndarray*) – input data

- **n_models** (Optional[int]) – number of models to use in predictions as ranked by the metric which was used when constructing the *Ensemble*. By default, entire ensemble is used.
- **parent_bar** (Optional[ConsoleMasterBar]) – not used when calling the method directly
- **display** (bool) – whether to display a progress bar for model evaluations

Return type ndarray

Returns Numpy array of predictions

Examples::

```
>>> preds = ensemble.predict_array(inputs)
```

predict_folds (*fy*, *n_models=None*, *pred_name='pred'*)

Apply ensemble to data accessed by a *FoldYielder* and save predictions as a new group per fold in the foldfile. If an output pipe has been added to the ensemble, then the predictions will be deprocessed. Inputs are expected to be preprocessed; i.e. any input pipe added to the ensemble is not used. If *foldyielder* has test-time augmentation, then predictions will be averaged over all augmented forms of the data.

Parameters

- **fy** (*FoldYielder*) – *FoldYielder* interfacing with the input data
- **n_models** (Optional[int]) – number of models to use in predictions as ranked by the metric which was used when constructing the *Ensemble*. By default, entire ensemble is used.
- **pred_name** (str) – name for new group of predictions

Examples::

```
>>> ensemble.predict_array(test_fy, pred_name='pred_tta')
```

Return type None

save (*name*, *feats=None*, *overwrite=False*)

Save ensemble and associated objects

Parameters

- **name** (str) – base name for saved objects
- **feats** (Optional[Any]) – optional list of input features
- **overwrite** (bool) – if existing objects are found, whether to overwrite them

Examples::

```
>>> ensemble.save('weights/ensemble')
>>>
>>> ensemble.save('weights/ensemble', ['pt', 'eta', 'phi'])
```

Return type None

Module contents

4.1.4 lumin.nn.interpretation package

Submodules

lumin.nn.interpretation.features module

```
lumin.nn.interpretation.features.get_nn_feat_importance(model, fy,
                                                         eval_metric=None,
                                                         pb_parent=None,
                                                         plot=True, save-
                                                         name=None, set-
                                                         tings=<lumin.plotting.plot_settings.PlotSettings
                                                         object>)
```

Compute permutation importance of features used by a *Model* on provided data using either loss or an *EvalMetric* to quantify performance. Returns bootstrapped mean importance from sample constructed by computing importance for each fold in *fy*.

Parameters

- **model** (*AbsModel*) – *Model* to use to evaluate feature importance
- **fy** (*FoldYielder*) – *FoldYielder* interfacing to data used to train model
- **eval_metric** (*Optional[EvalMetric]*) – Optional *EvalMetric* to use to quantify performance in place of loss
- **pb_parent** (*Optional[ConsoleMasterBar]*) – Not used if calling method directly
- **plot** (*bool*) – whether to plot resulting feature importances
- **savename** (*Optional[str]*) – Optional name of file to which to save the plot of feature importances
- **settings** (*PlotSettings*) – *PlotSettings* class to control figure appearance

Return type *DataFrame*

Returns *Pandas DataFrame* containing mean importance and associated uncertainty for each feature

Examples::

```
>>> fi = get_nn_feat_importance(model, train_fy)
>>>
>>> fi = get_nn_feat_importance(model, train_fy, savename='feat_import')
>>>
>>> fi = get_nn_feat_importance(model, train_fy,
...                             eval_metric=AMS(n_total=100000))
```

```
lumin.nn.interpretation.features.get_ensemble_feat_importance(ensemble, fy,
                                                             eval_metric=None,
                                                             save-
                                                             name=None, set-
                                                             tings=<lumin.plotting.plot_settings.PlotSe
                                                             object>)
```

Compute permutation importance of features used by an *Ensemble* on provided data using either loss or an *EvalMetric* to quantify performance. Returns bootstrapped mean importance from sample constructed by computing importance for each *Model* in ensemble.

Parameters

- **ensemble** (`AbsEnsemble`) – *Ensemble* to use to evaluate feature importance
- **fy** (`FoldYielder`) – *FoldYielder* interfacing to data used to train models in ensemble
- **eval_metric** (`Optional[EvalMetric]`) – Optional `EvalMetric` to use to quantify performance in place of loss
- **savename** (`Optional[str]`) – Optional name of file to which to save the plot of feature importances
- **settings** (`PlotSettings`) – *PlotSettings* class to control figure appearance

Return type `DataFrame`

Returns Pandas `DataFrame` containing mean importance and associated uncertainty for each feature

Examples::

```
>>> fi = get_ensemble_feat_importance(ensemble, train_fy)
>>>
>>> fi = get_ensemble_feat_importance(ensemble, train_fy,
...                                 savename='feat_import')
>>>
>>> fi = get_ensemble_feat_importance(ensemble, train_fy,
...                                 eval_metric=AMS(n_total=100000))
...

```

Module contents**4.1.5 lumin.nn.losses package****Submodules****lumin.nn.losses.basic_weighted module**

class `lumin.nn.losses.basic_weighted.WeightedMSE` (*weight=None*)

Bases: `torch.nn.modules.loss.MSELoss`

Class for computing Mean Squared-Error loss with optional weights per prediction. For compatibility with using basic PyTorch losses, weights are passed during initialisation rather than when computing the loss.

Parameters **weight** (`Optional[Tensor]`) – sample weights as PyTorch `Tensor`, to be used with data to be passed when computing the loss

Examples::

```
>>> loss = WeightedMSE()
>>>
>>> loss = WeightedMSE(weights)

```

forward (*input, target*)

Evaluate loss for given predictions

Parameters

- **input** (`Tensor`) – prediction tensor

- **target** (Tensor) – target tensor

Return type Tensor

Returns (weighted) loss

class lumin.nn.losses.basic_weighted.**WeightedMAE** (*weight=None*)

Bases: torch.nn.modules.loss.L1Loss

Class for computing Mean Absolute-Error loss with optional weights per prediction. For compatibility with using basic PyTorch losses, weights are passed during initialisation rather than when computing the loss.

Parameters **weight** (Optional[Tensor]) – sample weights as PyTorch Tensor, to be used with data to be passed when computing the loss

Examples::

```
>>> loss = WeightedMAE()
>>>
>>> loss = WeightedMAE(weights)
```

forward (*input, target*)

Evaluate loss for given predictions

Parameters

- **input** (Tensor) – prediction tensor
- **target** (Tensor) – target tensor

Return type Tensor

Returns (weighted) loss

class lumin.nn.losses.basic_weighted.**WeightedCCE** (*weight=None*)

Bases: torch.nn.modules.loss.NLLLoss

Class for computing Categorical Cross-Entropy loss with optional weights per prediction. For compatibility with using basic PyTorch losses, weights are passed during initialisation rather than when computing the loss.

Parameters **weight** (Optional[Tensor]) – sample weights as PyTorch Tensor, to be used with data to be passed when computing the loss

Examples::

```
>>> loss = WeightedCCE()
>>>
>>> loss = WeightedCCE(weights)
```

forward (*input, target*)

Evaluate loss for given predictions

Parameters

- **input** (Tensor) – prediction tensor
- **target** (Tensor) – target tensor

Return type Tensor

Returns (weighted) loss

lumin.nn.losses.hep_losses module

```
class lumin.nn.losses.hep_losses.SignificanceLoss (weight, sig_wgt=<class 'float'>,
                                                bkg_wgt=<class 'float'>,
                                                func=typing.Callable[[torch.Tensor,
                                                                    torch.Tensor], torch.Tensor])
```

Bases: torch.nn.modules.module.Module

General class for implementing significance-based loss functions, e.g. Asimov Loss (<https://arxiv.org/abs/1806.00322>). For compatibility with using basic PyTorch losses, event weights are passed during initialisation rather than when computing the loss.

Parameters

- **weight** (Tensor) – sample weights as PyTorch Tensor, to be used with data to be passed when computing the loss
- **sig_wgt** – total weight of signal events
- **bkg_wgt** – total weight of background events
- **func** – callable which returns a float based on signal and background weights

Examples::

```
>>> loss = SignificanceLoss(weight, sig_weight=sig_weight,
...                         bkg_weight=bkg_weight, func=calc_ams_torch)
>>>
>>> loss = SignificanceLoss(weight, sig_weight=sig_weight,
...                         bkg_weight=bkg_weight,
...                         func=partial(calc_ams_torch, br=10))
```

forward (*input*, *target*)

Evaluate loss for given predictions

Parameters

- **input** (Tensor) – prediction tensor
- **target** (Tensor) – target tensor

Return type Tensor

Returns (weighted) loss

Module contents

4.1.6 lumin.nn.metrics package

Submodules

lumin.nn.metrics.class_eval module

```
class lumin.nn.metrics.class_eval.AMS (n_total, wgt_name, targ_name='targets', br=0,
                                         syst_unc_b=0, use_quick_scan=True)
```

Bases: *lumin.nn.metrics.eval_metric.EvalMetric*

Class to compute maximum Approximate Median Significance (<https://arxiv.org/abs/1007.1727>) using classifier which directly predicts the class of data in a binary classification problem. AMS is computed on a single fold of

data provided by a *FoldYielder* and automatically reweights data by event multiplicity to account missing weights.

Parameters

- **n_total** (int) – total number of events in entire data set
- **wgt_name** (str) – name of weight group in fold file to use. N.B. if you have reweighted to balance classes, be sure to use the un-reweighted weights.
- **targ_name** (str) – name of target group in fold file
- **br** (float) – constant bias offset for background yield
- **syst_unc_b** (float) – fractional systematic uncertainty on background yield
- **use_quick_scan** (bool) – whether to optimise AMS by the *ams_scan_quick()* method (fast but suffers floating point precision) if False use *ams_scan_slow()* (slower but more accurate)

Examples::

```
>>> ams_metric = AMS(n_total=250000, br=10, wgt_name='gen_orig_weight')
>>>
>>> ams_metric = AMS(n_total=250000, syst_unc_b=0.1,
...                 wgt_name='gen_orig_weight', use_quick_scan=False)
```

evaluate (*fy, idx, y_pred*)

Compute maximum AMS on fold using provided predictions.

Parameters

- **fy** (*FoldYielder*) – *FoldYielder* interfacing to data
- **idx** (int) – fold index corresponding to fold for which *y_pred* was computed
- **y_pred** (ndarray) – predictions for fold

Return type float

Returns Maximum AMS computed on reweighted data from fold

Examples::

```
>>> ams = ams_metric.evaluate(train_fy, val_id, val_preds)
```

```
class lumin.nn.metrics.class_eval.MultiAMS (n_total, wgt_name, targ_name, zero_preds,
                                             one_preds, br=0, syst_unc_b=0,
                                             use_quick_scan=True)
```

Bases: *lumin.nn.metrics.class_eval.AMS*

Class to compute maximum Approximate Median Significance (<https://arxiv.org/abs/1007.1727>) using classifier which predicts the class of data in a multiclass classification problem which can be reduced to a binary classification problem AMS is computed on a single fold of data provided by a *FoldYielder* and automatically reweights data by event multiplicity to account missing weights.

Parameters

- **n_total** (int) – total number of events in entire data set
- **wgt_name** (str) – name of weight group in fold file to use. N.B. if you have reweighted to balance classes, be sure to use the un-reweighted weights.

- **targ_name** (`str`) – name of target group in fold file which indicates whether the event is signal or background
- **zero_preds** (`List[str]`) – list of predicted classes which correspond to class 0 in the form `pred_[i]`, where `i` is a NN output index
- **one_preds** (`List[str]`) – list of predicted classes which correspond to class 1 in the form `pred_[i]`, where `i` is a NN output index
- **br** (`float`) – constant bias offset for background yield
- **syst_unc_b** (`float`) – fractional systematic uncertainty on background yield
- **use_quick_scan** (`bool`) – whether to optimise AMS by the `ams_scan_quick()` method (fast but suffers floating point precision) if `False` use `ams_scan_slow()` (slower but more accurate)

Examples::

```

>>> ams_metric = MultiAMS(n_total=250000, br=10, targ_name='gen_target',
...                       wgt_name='gen_orig_weight',
...                       zero_preds=['pred_0', 'pred_1', 'pred_2'],
...                       one_preds=['pred_3'])
>>>
>>> ams_metric = MultiAMS(n_total=250000, syst_unc_b=0.1,
...                       targ_name='gen_target',
...                       wgt_name='gen_orig_weight',
...                       use_quick_scan=False,
...                       zero_preds=['pred_0', 'pred_1', 'pred_2'],
...                       one_preds=['pred_3'])

```

evaluate (`fy, idx, y_pred`)

Compute maximum AMS on fold using provided predictions.

Parameters

- **fy** (`FoldYielder`) – `FoldYielder` interfacing to data
- **idx** (`int`) – fold index corresponding to fold for which `y_pred` was computed
- **y_pred** (`ndarray`) – predictions for fold

Return type `float`

Returns Maximum AMS computed on reweighted data from fold

Examples::

```

>>> ams = ams_metric.evaluate(train_fy, val_id, val_preds)

```

lumin.nn.metrics.eval_metric module

class `lumin.nn.metrics.eval_metric.EvalMetric` (`targ_name, wgt_name=None`)

Bases: `abc.ABC`

Abstract class for evaluating performance of a model using some metric

Parameters

- **targ_name** (`str`) – name of group in fold file containing regression targets

- **wgt_name** (Optional[str]) – name of group in fold file containing datapoint weights

abstract evaluate (*fy, idx, y_pred*)

Evaluate the required metric for a given fold and set of predictions

Parameters

- **fy** (*FoldYielder*) – *FoldYielder* interfacing to data
- **idx** (int) – fold index corresponding to fold for which *y_pred* was computed
- **y_pred** (ndarray) – predictions for fold

Return type float

Returns metric value

get_df (*fy, idx, y_pred*)

Returns a DataFrame for the given fold containing targets, weights, and predictions

Parameters

- **fy** (*FoldYielder*) – *FoldYielder* interfacing to data
- **idx** (int) – fold index corresponding to fold for which *y_pred* was computed
- **y_pred** (ndarray) – predictions for fold

Return type DataFrame

Returns DataFrame for the given fold containing targets, weights, and predictions

lumin.nn.metrics.reg_eval module

```
class lumin.nn.metrics.reg_eval.RegPull (return_mean, use_bootstrap=False,
                                         use_weights=True, use_pull=True,
                                         targ_name='targets', wgt_name=None)
```

Bases: *lumin.nn.metrics.eval_metric.EvalMetric*

Compute mean or standard deviation of delta or pull of some feature which is being directly regressed to. Optionally, use bootstrap resampling on validation data.

Parameters

- **return_mean** (bool) – whether to return the mean or the standard deviation
- **use_bootstrap** (bool) – whether to bootstrap resamples validation fold when computing statistic
- **use_weights** (bool) – whether to actually use weights if *wgt_name* is set
- **use_pull** (bool) – whether to return the pull (differences / targets) or delta (differences)
- **targ_name** (str) – name of group in fold file containing regression targets
- **wgt_name** (Optional[str]) – name of group in fold file containing datapoint weights

Examples::

```
>>> mean_pull = RegPull(return_mean=True, use_bootstrap=True,
...                     use_pull=True)
>>>
>>> std_delta = RegPull(return_mean=False, use_bootstrap=True,
...                     use_pull=False)
```

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```
>>>
>>> mean_pull = RegPull(return_mean=True, use_bootstrap=False,
...                     use_pull=True, wgt_name='weights')
```

evaluate (*fy, idx, y_pred*)

Compute statistic on fold using provided predictions.

Parameters

- **fy** (*FoldYielder*) – *FoldYielder* interfacing to data
- **idx** (int) – fold index corresponding to fold for which *y_pred* was computed
- **y_pred** (ndarray) – predictions for fold

Return type float

Returns Statistic set in initialisation computed on the chosen fold

Examples::

```
>>> mean = mean_pull.evaluate(train_fy, val_id, val_preds)
```

```
class lumin.nn.metrics.reg_eval.RegAsProxyPull (proxy_func,          return_mean,
                                                use_bootstrap=False,
                                                use_weights=True,    use_pull=True,
                                                targ_name='targets',
                                                wgt_name=None)
```

Bases: *lumin.nn.metrics.reg_eval.RegPull*

Compute mean or standard deviation of delta or pull of some feature which is being indirectly regressed to via a proxy function. Optionally, use bootstrap resampling on validation data.

Parameters

- **proxy_func** (Callable[[Dataframe], None]) – function which acts on regression predictions and adds *pred* and *gen_target* columns to the Pandas DataFrame it is passed which contains prediction columns *pred_{i}*
- **return_mean** (bool) – whether to return the mean or the standard deviation
- **use_bootstrap** (bool) – whether to bootstrap resamples validation fold when computing statistic
- **use_weights** (bool) – whether to actually use weights if *wgt_name* is set
- **use_pull** (bool) – whether to return the pull (differences / targets) or delta (differences)
- **targ_name** (str) – name of group in fold file containing regression targets
- **wgt_name** (Optional[str]) – name of group in fold file containing datapoint weights

Examples::

```
>>> def reg_proxy_func(df):
>>>     df['pred'] = calc_pair_mass(df, (1.77682, 1.77682),
...                               {targ[targ.find('_t')+3:]:
...                               f'pred_{i}' for i, targ
...                               in enumerate(targ_feats)})
>>>     df['gen_target'] = 125
```

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```
>>>
>>> std_delta = RegAsProxyPull(proxy_func=reg_proxy_func,
...                             return_mean=False, use_pull=False)
```

evaluate (*fy, idx, y_pred*)

Compute statistic on fold using provided predictions.

Parameters

- **fy** (*FoldYielder*) – *FoldYielder* interfacing to data
- **idx** (int) – fold index corresponding to fold for which *y_pred* was computed
- **y_pred** (ndarray) – predictions for fold

Return type float

Returns Statistic set in initialisation computed on the chosen fold

Examples::

```
>>> mean = mean_pull.evaluate(train_fy, val_id, val_preds)
```

Module contents

4.1.7 lumin.nn.models package

Subpackages

[lumin.nn.models.blocks package](#)

Submodules

[lumin.nn.models.blocks.body module](#)

```
class lumin.nn.models.blocks.body.FullyConnected (n_in, feat_map, depth, width, do=0,  
                                                    bn=False, act='relu', res=False,  
                                                    dense=False, growth_rate=0,  
                                                    lookup_init=<function  
                                                    lookup_normal_init>,  
                                                    lookup_act=<function lookup_act>,  
                                                    freeze=False)
```

Bases: `lumin.nn.models.blocks.body.AbsBody`

Fully connected set of hidden layers. Designed to be passed as a ‘body’ to *ModelBuilder*. Supports batch normalisation and dropout. Order is dense->activation->BN->DO, except when *res* is true in which case the BN is applied after the addition. Can optionally have skip connections between each layer (*res*=true). Alternatively can concatenate layers (*dense*=true) *growth_rate* parameter can be used to adjust the width of layers according to $width+(width*(depth-1)*growth_rate)$

Parameters

- **n_in** (int) – number of inputs to the block

- **feat_map** (Dict[str, List[int]]) – dictionary mapping input features to the model to outputs of head block
- **depth** (int) – number of hidden layers. If `res==True` and depth is even, depth will be increased by one.
- **width** (int) – base width of each hidden layer
- **do** (float) – if not None will add dropout layers with dropout rates do
- **bn** (bool) – whether to use batch normalisation
- **act** (str) – string representation of argument to pass to `lookup_act`
- **res** (bool) – whether to add an additive skip connection every two dense layers. Mutually exclusive with `dense`.
- **dense** (bool) – whether to perform layer-wise concatenations after every layer. Mutually exclusion with `res`.
- **growth_rate** (int) – rate at which width of dense layers should increase with depth beyond the initial layer. Ignored if `res=True`. Can be negative.
- **lookup_init** (Callable[[str, Optional[int], Optional[int]], Callable[[Tensor], None]]) – function taking choice of activation function, number of inputs, and number of outputs and returning a function to initialise layer weights.
- **lookup_act** (Callable[[str], Any]) – function taking choice of activation function and returning an activation function layer
- **freeze** (bool) – whether to start with module parameters set to untrainable

Examples::

```

>>> body = FullyConnected(n_in=32, feat_map=head.feat_map, depth=4,
...                        width=100, act='relu')
>>>
>>> body = FullyConnected(n_in=32, feat_map=head.feat_map, depth=4,
...                        width=200, act='relu', growth_rate=-0.3)
>>>
>>> body = FullyConnected(n_in=32, feat_map=head.feat_map, depth=4,
...                        width=100, act='swish', do=0.1, res=True)
>>>
>>> body = FullyConnected(n_in=32, feat_map=head.feat_map, depth=6,
...                        width=32, act='selu', dense=True,
...                        growth_rate=0.5)
>>>
>>> body = FullyConnected(n_in=32, feat_map=head.feat_map, depth=6,
...                        width=50, act='prelu', bn=True,
...                        lookup_init=lookup_uniform_init)

```

forward (*x*)

Pass tensor through body

Parameters *x* (Tensor) – incoming tensor**Returns** Resulting tensor**Return type** Tensor

`get_out_size()`
Get size width of output layer

Return type `int`

Returns Width of output layer

```
class lumin.nn.models.blocks.body.MultiBlock(n_in, feat_map, blocks,
                                              feats_per_block, bottleneck_sz=0, bot-
                                              tleneck_act=None, lookup_init=<function
                                              lookup_normal_init>,
                                              lookup_act=<function lookup_act>,
                                              freeze=False)
```

Bases: `lumin.nn.models.blocks.body.AbsBody`

Body block allowing outputs of head block to be split amongst a series of body blocks. Output is the concatenation of all sub-body blocks. Optionally, single-neuron ‘bottleneck’ layers can be used to pass an input to each sub-block based on a learned function of the input features that block would otherwise not receive, i.e. a highly compressed representation of the rest of the feature space.

Parameters

- **n_in** (`int`) – number of inputs to the block
- **feat_map** (`Dict[str, List[int]]`) – dictionary mapping input features to the model to outputs of head block
- **blocks** (`List[partial]`) – list of uninstantiated `AbsBody` blocks to which to pass a subsection of the total inputs. Note that partials should be used to set any relevant parameters at initialisation time
- **feats_per_block** (`List[List[str]]`) – list of lists of names of features to pass to each `AbsBody`, not that the `feat_map` provided by `AbsHead` will map features to their relevant head outputs
- **bottleneck** – if true, each block will receive the output of a single neuron which takes as input all the features which each given block does not directly take as inputs
- **bottleneck_act** (`Optional[str]`) – if set to a string representation of an activation function, the output of each bottleneck neuron will be passed through the defined activation function before being passed to their associated blocks
- **lookup_init** (`Callable[[str, Optional[int], Optional[int]], Callable[[Tensor], None]]`) – function taking choice of activation function, number of inputs, and number of outputs and returning a function to initialise layer weights.
- **lookup_act** (`Callable[[str], Any]`) – function taking choice of activation function and returning an activation function layer
- **freeze** (`bool`) – whether to start with module parameters set to untrainable

Examples::

```
>>> body = MultiBlock(
...     blocks=[partial(FullyConnected, depth=1, width=50, act='swish'),
...                 partial(FullyConnected, depth=6, width=55, act='swish',
...                           dense=True, growth_rate=-0.1)],
...     feats_per_block=[[f for f in train_feats if 'DER_' in f],
...                       [f for f in train_feats if 'PRI_' in f]])
>>>
>>> body = MultiBlock(
```

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```

...     blocks=[partial(FullyConnected, depth=1, width=50, act='swish'),
...     partial(FullyConnected, depth=6, width=55, act='swish',
...             dense=True, growth_rate=-0.1)],
...     feats_per_block=[[f for f in train_feats if 'DER_' in f],
...                      [f for f in train_feats if 'PRI_' in f]],
...     bottleneck=True)
>>>
>>> body = MultiBlock(
...     blocks=[partial(FullyConnected, depth=1, width=50, act='swish'),
...             partial(FullyConnected, depth=6, width=55, act='swish',
...                     dense=True, growth_rate=-0.1)],
...     feats_per_block=[[f for f in train_feats if 'DER_' in f],
...                      [f for f in train_feats if 'PRI_' in f]],
...     bottleneck=True, bottleneck_act='swish')

```

forward (*x*)

Pass tensor through body

Parameters *x* (Tensor) – incoming tensor**Returns** Resulting tensor**Return type** Tensor**get_out_size** ()

Get size width of output layer

Return type int**Returns** Total number of outputs across all blocks**lumin.nn.models.blocks.endcap** module**class** lumin.nn.models.blocks.endcap.**AbsEndcap** (*model*)

Bases: torch.nn.modules.module.Module

Abstract class for constructing post training layer which performs further calculation on NN outputs. Used when NN was trained to some proxy objective

Parameters *model* (Module) – trained *Model* to wrap**forward** (*x*)

Pass tensor through endcap and compute function

Parameters *x* (Tensor) – model output tensor**Returns** Resulting tensor**Return type** Tensor**abstract func** (*x*)

Transformation function to apply to model outputs

Arguments: *x*: model output tensor**Return type** Tensor

Returns Resulting tensor

predict (*inputs*, *as_np=True*)

Evaluate model on input tensor, and compute function of model outputs

Parameters

- **inputs** (Union[ndarray, DataFrame, Tensor]) – input data as Numpy array, Pandas DataFrame, or tensor on device
- **as_np** (bool) – whether to return predictions as Numpy array (otherwise tensor)

Return type Union[ndarray, Tensor]

Returns model predictions pass through endcap function

lumin.nn.models.blocks.head module

```
class lumin.nn.models.blocks.head.CatEmbHead (cont_feats, do_cont=0,
                                              do_cat=0, cat_embedder=None,
                                              lookup_init=<function>
                                              lookup_normal_init>, freeze=False)
```

Bases: lumin.nn.models.blocks.head.AbsHead

Standard model head for columnar data. Provides inputs for continuous features and embedding matrices for categorical inputs, and uses a dense layer to upscale to width of network body. Designed to be passed as a ‘head’ to *ModelBuilder*. Supports batch normalisation and dropout (at separate rates for continuous features and categorical embeddings). Continuous features are expected to be the first len(cont_feats) columns of input tensors and categorical features the remaining columns. Embedding arguments for categorical features are set using a *CatEmbedder*.

Parameters

- **cont_feats** (List[str]) – list of names of continuous input features
- **do_cont** (float) – if not None will add a dropout layer with dropout rate do acting on the continuous inputs prior to concatenation with the categorical embeddings
- **do_cat** (float) – if not None will add a dropout layer with dropout rate do acting on the categorical embeddings prior to concatenation with the continuous inputs
- **cat_embedder** (Optional[*CatEmbedder*]) – *CatEmbedder* providing details of how to embed categorical inputs
- **lookup_init** (Callable[[str, Optional[int], Optional[int]], Callable[[Tensor], None]]) – function taking choice of activation function, number of inputs, and number of outputs and returning a function to initialise layer weights.
- **freeze** (bool) – whether to start with module parameters set to untrainable

Examples::

```
>>> head = CatEmbHead(cont_feats=cont_feats)
>>>
>>> head = CatEmbHead(cont_feats=cont_feats,
...                   cat_embedder=CatEmbedder.from_fy(train_fy))
>>>
>>> head = CatEmbHead(cont_feats=cont_feats,
...                   cat_embedder=CatEmbedder.from_fy(train_fy),
...                   do_cont=0.1, do_cat=0.05)
```

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```

>>>
>>> head = CatEmbHead(cont_feats=cont_feats,
...                   cat_embedder=CatEmbedder.from_fy(train_fy),
...                   lookup_init=lookup_uniform_init)

```

forward (*x_{in}*)

Pass tensor through head

Parameters *x* – input tensor**Returns** Resulting tensor**Return type** Tensor**get_embeds** ()

Get state_dict for every embedding matrix.

Return type Dict[str, OrderedDict]**Returns** Dictionary mapping categorical features to learned embedding matrix**get_out_size** ()

Get size width of output layer

Return type int**Returns** Width of output layer**plot_embeds** (*savename=None, settings=<lumin.plotting.plot_settings.PlotSettings object>*)

Plot representations of embedding matrices for each categorical feature.

Parameters

- **savename** (Optional[str]) – if not None, will save copy of plot to give path
- **settings** (*PlotSettings*) – *PlotSettings* class to control figure appearance

Return type None**save_embeds** (*path*)

Save learned embeddings to path. Each categorical embedding matrix will be saved as a separate state_dict with name equal to the feature name as set in cat_embedder

Parameters *path* (Path) – path to which to save embedding weights**Return type** None**lumin.nn.models.blocks.tail module**

```

class lumin.nn.models.blocks.tail.ClassRegMulti(n_in, n_out, objective,
                                                y_range=None, bias_init=None,
                                                lookup_init=<function
                                                lookup_normal_init>, freeze=False)

```

Bases: lumin.nn.models.blocks.tail.AbsTail

Output block for (multi(class/label)) classification or regression tasks. Designed to be passed as a ‘tail’ to *ModelBuilder*. Takes output size of network body and scales it to required number of outputs. For regression tasks, *y_range* can be set with per-output minima and maxima. The outputs are then adjusted according to $((y_{\max}-y_{\min})*x)+self.y_{\min}$, where *x* is the output of the network passed through a sigmoid function.

Effectively allowing regression to be performed without normalising and standardising the target values. Note it is safest to allow some leeway in setting the min and max, e.g. $\text{max} = 1.2 * \text{max}$, $\text{min} = 0.8 * \text{min}$. Output activation function is automatically set according to objective and `y_range`.

Parameters

- **n_in** (`int`) – number of inputs to expect
- **n_out** (`int`) – number of outputs required
- **objective** (`str`) – string representation of network objective, i.e. ‘classification’, ‘regression’, ‘multiclass’
- **y_range** (`Union[Tuple, ndarray, None]`) – if not `None`, will apply rescaling to network outputs.
- **lookup_init** (`Callable[[str, Optional[int], Optional[int]], Callable[[Tensor], None]]`) – function taking string representation of activation function, number of inputs, and number of outputs and returning a function to initialise layer weights.

Examples::

```
>>> tail = ClassRegMulti(n_in=100, n_out=1, objective='classification')
>>>
>>> tail = ClassRegMulti(n_in=100, n_out=5, objective='multiclass')
>>>
>>> y_range = (0.8*targets.min(), 1.2*targets.max())
>>> tail = ClassRegMulti(n_in=100, n_out=1, objective='regression',
...                       y_range=y_range)
>>>
>>> min_targs = np.min(targets, axis=0).reshape(targets.shape[1],1)
>>> max_targs = np.max(targets, axis=0).reshape(targets.shape[1],1)
>>> min_targs[min_targs > 0] *=0.8
>>> min_targs[min_targs < 0] *=1.2
>>> max_targs[max_targs > 0] *=1.2
>>> max_targs[max_targs < 0] *=0.8
>>> y_range = np.hstack((min_targs, max_targs))
>>> tail = ClassRegMulti(n_in=100, n_out=6, objective='regression',
...                       y_range=y_range,
...                       lookup_init=lookup_uniform_init)
```

forward (*x*)

Pass tensor through tail

Parameters **x** (`Tensor`) – incoming tensor

Returns Resulting tensor of model outputs

Return type `Tensor`

get_out_size ()

Get size width of output layer

Return type `int`

Returns Width of output layer

Module contents

lumin.nn.models.layers package

Submodules

lumin.nn.models.layers.activations module

`lumin.nn.models.layers.activations.lookup_act` (*act*)

Map activation name to class

Parameters *act* (*str*) – string representation of activation function

Return type *Any*

Returns Class implementing requested activation function

class `lumin.nn.models.layers.activations.Swish` (*inplace=False*)

Bases: `torch.nn.modules.module.Module`

Non-trainable Swish activation function <https://arxiv.org/abs/1710.05941>

Parameters *inplace* – whether to apply activation inplace

Examples::

```
>>> swish = Swish()
```

forward (*x*)

Pass tensor through Swish function

Parameters *x* (*Tensor*) – incoming tensor

Return type *Tensor*

Returns Resulting tensor

Module contents

Submodules

lumin.nn.models.helpers module

class `lumin.nn.models.helpers.CatEmbedder` (*cat_names*, *cat_szs*, *emb_szs=None*,
max_emb_sz=50, *emb_load_path=None*)

Bases: `object`

Helper class for embedding categorical features. Designed to be passed to `ModelBuilder`. Note that the classmethod `from_fy()` may be used to instantiate an `CatEmbedder` from a `FoldYielder`.

Parameters

- **cat_names** (*List[str]*) – list of names of categorical features in order in which they will be passed as inputs columns
- **cat_szs** (*List[int]*) – list of cardinalities (number of unique elements) for each feature
- **emb_szs** (*Optional[List[int]]*) – Optional list of embedding sizes for each feature. If *None*, will use $\min(\text{max_emb_sz}, (1+\text{sz})//2)$

- **max_emb_sz** (int) – Maximum size of embedding if `emb_szs` is None
- **emb_load_path** (Union[Path, str, None]) – if not None, will cause *ModelBuilder* to attempt to load pretrained embeddings from path

Examples::

```
>>> cat_embedder = CatEmbedder(cat_names=['n_jets', 'channel'],
                               cat_szs=[5, 3])
>>>
>>> cat_embedder = CatEmbedder(cat_names=['n_jets', 'channel'],
                               cat_szs=[5, 3], emb_szs=[2, 2])
>>>
>>> cat_embedder = CatEmbedder(cat_names=['n_jets', 'channel'],
                               cat_szs=[5, 3], emb_szs=[2, 2],
                               emb_load_path=Path('weights'))
```

calc_emb_szs ()

Method used to set sizes of embeddings for each categorical feature when no embedding sizes are explicitly passed Uses rule of thumb of $\min(50, (1+\text{cardinality})/2)$

Return type None

classmethod from_fy (fy, emb_szs=None, max_emb_sz=50, emb_load_path=None)

Instantiate an *CatEmbedder* from a *FoldYielder*, i.e. avoid having to pass `cat_names` and `cat_szs`.

Parameters

- **fy** (*FoldYielder*) – *FoldYielder* with training data
- **emb_szs** (Optional[List[int]]) – Optional list of embedding sizes for each feature. If None, will use $\min(\text{max_emb_sz}, (1+\text{sz})/2)$
- **max_emb_sz** (int) – Maximum size of embedding if `emb_szs` is None
- **emb_load_path** (Union[Path, str, None]) – if not None, will cause *ModelBuilder* to attempt to load pretrained embeddings from path

Returns *CatEmbedder*

Examples::

```
>>> cat_embedder = CatEmbedder.from_fy(train_fy)
>>>
>>> cat_embedder = CatEmbedder.from_fy(train_fy, emb_szs=[2, 2])
>>>
>>> cat_embedder = CatEmbedder.from_fy(
    train_fy, emb_szs=[2, 2],
    emb_load_path=Path('weights'))
```

`lumin.nn.models.helpers.Embedder` (`cat_names`, `cat_szs`, `emb_szs=None`, `max_emb_sz=50`, `emb_load_path=None`)

Attention: Depreciated in favour of *CatEmbedder* and will be removed in v0.4.

lumin.nn.models.initialisations module

`lumin.nn.models.initialisations.lookup_normal_init` (*act*, *fan_in=None*,
fan_out=None)

Lookup for weight initialisation using Normal distributions

Parameters

- **act** (*str*) – string representation of activation function
- **fan_in** (*Optional[int]*) – number of inputs to neuron
- **fan_out** (*Optional[int]*) – number of outputs from neuron

Return type *Callable[[Tensor], None]*

Returns Callable to initialise weight tensor

`lumin.nn.models.initialisations.lookup_uniform_init` (*act*, *fan_in=None*,
fan_out=None)

Lookup weight initialisation using Uniform distributions

Parameters

- **act** (*str*) – string representation of activation function
- **fan_in** (*Optional[int]*) – number of inputs to neuron
- **fan_out** (*Optional[int]*) – number of outputs from neuron

Return type *Callable[[Tensor], None]*

Returns Callable to initialise weight tensor

lumin.nn.models.model module

class `lumin.nn.models.model.Model` (*model_builder=None*)

Bases: `lumin.nn.models.abs_model.AbsModel`

Wrapper class to handle training and inference of NNs created via a *ModelBuilder*. Note that saved models can be instantiated directly via *from_save()* classmethod.

Parameters **model_builder** (*Optional[ModelBuilder]*) – *ModelBuilder* which will construct the network, loss, and optimiser

Examples::

```
>>> model = Model(model_builder)
```

evaluate (*inputs, targets, weights=None, callbacks=None, mask_inputs=True*)

Compute loss on provided data.

Parameters

- **inputs** (*Tensor*) – input data as tensor on device
- **targets** (*Tensor*) – targets as tensor on device
- **weights** (*Optional[Tensor]*) – Optional weights as tensor on device
- **callbacks** (*Optional[List[AbsCallback]]*) – list of any callbacks to use during evaluation
- **mask_inputs** (*bool*) – whether to apply input mask if one has been set

Return type float

Returns (weighted) loss of model predictions on provided data

export2onnx (*name*, *bs=1*)

Export network to ONNX format. Note that ONNX expects a fixed batch size (*bs*) which is the number of datapoints your wish to pass through the model concurrently.

Parameters

- **name** (str) – filename for exported file
- **bs** (int) – batch size for exported models

Return type None

export2tfpb (*name*, *bs=1*)

Export network to Tensorflow ProtocolBuffer format, via ONNX. Note that ONNX expects a fixed batch size (*bs*) which is the number of datapoints your wish to pass through the model concurrently.

Parameters

- **name** (str) – filename for exported file
- **bs** (int) – batch size for exported models

Return type None

fit (*batch_yielder*, *callbacks=None*)

Fit network for one complete iteration of a *BatchYielder*, i.e. one (sub-)epoch

Parameters

- **batch_yielder** (*BatchYielder*) – *BatchYielder* providing training data in form of tuple of inputs, targets, and weights as tensors on device
- **callbacks** (Optional[List[AbsCallback]]) – list of *AbsCallback* to be used during training

Return type float

Returns Loss on training data averaged across all minibatches

classmethod from_save (*name*, *model_builder*)

Instantiated a *Model* and load saved state from file.

Parameters

- **name** (str) – name of file containing saved state
- **model_builder** (*ModelBuilder*) – *ModelBuilder* which was used to construct the network

Return type AbsModel

Returns Instantiated *Model* with network weights, optimiser state, and input mask loaded from saved state

Examples::

```
>>> model = Model.from_save('weights/model.h5', model_builder)
```

get_feat_importance (*fy*, *eval_metric=None*)

Call *get_nn_feat_importance()* passing this *Model* and provided arguments

Parameters

- **fy** (*FoldYielder*) – *FoldYielder* interfacing to data on which to evaluate importance
- **eval_metric** (*Optional[EvalMetric]*) – Optional *EvalMetric* to use for quantifying performance

Return type *DataFrame*

get_lr()

Get learning rate of optimiser

Return type *float*

Returns learning rate of optimiser

get_mom()

Get momentum/beta_1 of optimiser

Return type *float*

Returns momentum/beta_1 of optimiser

get_out_size()

Get number of outputs of model

Return type *int*

Returns Number of outputs of model

get_param_count (*trainable=True*)

Return number of parameters in model.

Parameters **trainable** (*bool*) – if true (default) only count trainable parameters

Return type *int*

Returns Number of (trainable) parameters in model

get_weights()

Get state_dict of weights for network

Return type *OrderedDict*

Returns state_dict of weights for network

load (*name, model_builder=None*)

Load model, optimiser, and input mask states from file

Parameters

- **name** (*str*) – name of save file
- **model_builder** (*Optional[ModelBuilder]*) – if *Model* was not initialised with a *ModelBuilder*, you will need to pass one here

Return type *None*

predict (*inputs, as_np=True, pred_name='pred'*)

Apply model to inputted data and compute predictions. A compatibility method to call *predict_array()* or meth:~*lumin.nn.models.model.Model.predict_folds*, depending on input type.

Parameters

- **inputs** (*Union[ndarray, DataFrame, Tensor, FoldYielder]*) – input data as Numpy array, Pandas DataFrame, or tensor on device, or *FoldYielder* interfacing to data

- **as_np** (bool) – whether to return predictions as Numpy array (otherwise tensor) if inputs are a Numpy array, Pandas DataFrame, or tensor
- **pred_name** (str) – name of group to which to save predictions if inputs are a *FoldYielder*

Return type Union[ndarray, Tensor, None]

Returns if inputs are a Numpy array, Pandas DataFrame, or tensor, will return predictions as either array or tensor

predict_array (*inputs*, *as_np=True*, *mask_inputs=True*)

Pass inputs through network and obtain predictions.

Parameters

- **inputs** (Union[ndarray, DataFrame, Tensor]) – input data as Numpy array, Pandas DataFrame, or tensor on device
- **as_np** (bool) – whether to return predictions as Numpy array (otherwise tensor)
- **mask_inputs** (bool) – whether to apply input mask if one has been set

Return type Union[ndarray, Tensor]

Returns Model prediction(s) per datapoint

predict_folds (*fy*, *pred_name='pred'*)

Apply model to all data accessed by a *FoldYielder* and save predictions as new group in fold file

Parameters

- **fy** (*FoldYielder*) – *FoldYielder* interfacing to data
- **pred_name** (str) – name of group to which to save predictions

Return type None

save (*name*)

Save model, optimiser, and input mask states to file

Parameters **name** (str) – name of save file

Return type None

set_input_mask (*mask*)

Mask input columns by only using input columns whose indices are listed in mask

Parameters **mask** (ndarray) – array of column indices to use from all input columns

Return type None

set_lr (*lr*)

set learning rate of optimiser

Parameters **lr** (float) – learning rate of optimiser

Return type None

set_mom (*mom*)

Set momentum/beta_1 of optimiser

Parameters **mom** (float) – momentum/beta_1 of optimiser

Return type None

set_weights (*weights*)

Set state_dict of weights for network

Parameters `weights` (OrderedDict) – state_dict of weights for network

Return type None

`lumin.nn.models.model_builder` module

```
class lumin.nn.models.model_builder.ModelBuilder (objective, n_out, cont_feats=None,
model_args=None, opt_args=None,
cat_embedder=None,
loss='auto', head=<class 'lumin.nn.models.blocks.head.CatEmbHead'>,
body=<class 'lumin.nn.models.blocks.body.FullyConnected'>,
tail=<class 'lumin.nn.models.blocks.tail.ClassRegMulti'>,
lookup_init=<function
lookup_normal_init>,
lookup_act=<function
lookup_act>, pretrain_file=None,
freeze_head=False,
freeze_body=False,
freeze_tail=False, cat_args=None,
n_cont_in=None)
```

Bases: object

Class to build models to specified architecture on demand along with an optimiser.

Attention: `cat_args` is now depreciated in favour of `cat_embedder` and will be removed in *v0.4*

Attention: `n_cont_in` is now depreciated in favour of `cont_feats` and will be removed in *v0.4*

Parameters

- **objective** (str) – string representation of network objective, i.e. ‘classification’, ‘regression’, ‘multiclass’
- **n_out** (int) – number of outputs required
- **cont_feats** (Optional[List[str]]) – list of names of continuous input features
- **model_args** (Optional[Dict[str, Dict[str, Any]]]) – dictionary of dictionaries of keyword arguments to pass to head, body, and tail to control architecture
- **opt_args** (Optional[Dict[str, Any]]) – dictionary of arguments to pass to optimiser. Missing kargs will be filled with default values. Currently, only ADAM (default), RAdam, Ranger, and SGD are available.
- **cat_embedder** (Optional[CatEmbedder]) – *CatEmbedder* for embedding categorical inputs
- **loss** (Any) – either and uninstantiated loss class, or leave as ‘auto’ to select loss according to objective
- **head** (AbsHead) – uninstantiated class which can receive input data and upscale it to model width

- **body** (AbsBody) – uninstantiated class which implements the main bulk of the model's hidden layers
- **tail** (AbsTail) – uninstantiated class which scales the body to the required number of outputs and implements any final activation function and output scaling
- **lookup_init** (Callable[[str, Optional[int], Optional[int]], Callable[[Tensor], None]]) – function taking choice of activation function, number of inputs, and number of outputs and returning a function to initialise layer weights.
- **lookup_act** (Callable[[str], Module]) – function taking choice of activation function and returning an activation function layer
- **pretrain_file** (Optional[str]) – if set, will load saved parameters for entire network from saved model
- **freeze_head** (bool) – whether to start with the head parameters set to untrainable
- **freeze_body** (bool) – whether to start with the body parameters set to untrainable
- **cat_args** (Optional[Dict[str, Any]]) – deprecated in place of cat_embedder
- **n_cont_in** (Optional[int]) – deprecated in favour of cont_feats

Examples::

```

>>> model_builder = ModelBuilder(objective='classifier',
>>>                               cont_feats=cont_feats, n_out=1,
>>>                               model_args={'body':{'depth':4,
>>>                                               'width':100}})
>>>
>>> min_targs = np.min(targets, axis=0).reshape(targets.shape[1],1)
>>> max_targs = np.max(targets, axis=0).reshape(targets.shape[1],1)
>>> min_targs[min_targs > 0] *=0.8
>>> min_targs[min_targs < 0] *=1.2
>>> max_targs[max_targs > 0] *=1.2
>>> max_targs[max_targs < 0] *=0.8
>>> y_range = np.hstack((min_targs, max_targs))
>>> model_builder = ModelBuilder(
>>>     objective='regression', cont_feats=cont_feats, n_out=6,
>>>     cat_embedder=CatEmbedder.from_fy(train_fy),
>>>     model_args={'body':{'depth':4, 'width':100},
>>>                 'tail':{'y_range=y_range}})
>>>
>>> model_builder = ModelBuilder(objective='multiclassifier',
>>>                               cont_feats=cont_feats, n_out=5,
>>>                               model_args={'body':{'width':100,
>>>                                               'depth':6,
>>>                                               'do':0.1,
>>>                                               'res':True}})
>>>
>>> model_builder = ModelBuilder(objective='classifier',
>>>                               cont_feats=cont_feats, n_out=1,
>>>                               model_args={'body':{'depth':4,
>>>                                               'width':100}},
>>>                               opt_args={'opt':'sgd',
>>>                                           'momentum':0.8,
>>>                                           'weight_decay':1e-5},
>>>                               loss=partial(SignificanceLoss,
>>>                                             sig_weight=sig_weight,

```

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```
>>>                                     bkg_weight=bkg_weight,
>>>                                     func=calc_ams_torch))
```

build_model()

Construct entire network module

Return type Module**Returns** Instantiated nn.Module

classmethod from_model_builder(*model_builder*, *pretrain_file=None*, *freeze_head=False*, *freeze_body=False*, *freeze_tail=False*, *loss=None*, *opt_args=None*)

Instantiate a *ModelBuilder* from an existing *ModelBuilder*, but with options to adjust loss, optimiser, pretraining, and module freezing

Parameters

- **model_builder** – existing *ModelBuilder* or filename for a pickled *ModelBuilder*
- **pretrain_file** (Optional[str]) – if set, will load saved parameters for entire network from saved model
- **freeze_head** (bool) – whether to start with the head parameters set to untrainable
- **freeze_body** (bool) – whether to start with the body parameters set to untrainable
- **freeze_tail** (bool) – whether to start with the tail parameters set to untrainable
- **loss** (Optional[Any]) – either and uninstantiated loss class, or leave as ‘auto’ to select loss according to objective
- **opt_args** (Optional[Dict[str, Any]]) – dictionary of arguments to pass to optimiser. Missing kargs will be filled with default values. Choice of optimiser (‘opt’) keyword can either be set by passing the string name (e.g. ‘adam’), but only ADAM and SGD are available this way, or by passing an uninstantiated optimiser (e.g. torch.optim.Adam). If no optimiser is set, then it defaults to ADAM. Additional keyword arguments can be set, and these will be passed to the optimiser during instantiation

Returns Instantiated *ModelBuilder***Examples::**

```
>>> new_model_builder = ModelBuilder.from_model_builder(
>>>     ModelBuilder)
>>>
>>> new_model_builder = ModelBuilder.from_model_builder(
>>>     ModelBuilder, loss=partial(
>>>         SignificanceLoss, sig_weight=sig_weight,
>>>         bkg_weight=bkg_weight, func=calc_ams_torch))
>>>
>>> new_model_builder = ModelBuilder.from_model_builder(
>>>     'weights/model_builder.pkl',
>>>     opt_args={'opt':'sgd', 'momentum':0.8, 'weight_decay':1e-5})
>>>
>>> new_model_builder = ModelBuilder.from_model_builder(
>>>     'weights/model_builder.pkl',
>>>     opt_args={'opt':torch.optim.Adam,
```

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```

...         'momentum':0.8,
...         'weight_decay':1e-5})

```

get_body (*n_in*, *feat_map*)

Construct body module

Return type `AbsBody`**Returns** Instantiated body `nn.Module`**get_head** ()

Construct head module

Return type `AbsHead`**Returns** Instantiated head `nn.Module`**get_model** ()

Construct model, loss, and optimiser, optionally loading pretrained weights

Return type `Tuple[Module, Optimizer, Any]`**Returns** Instantiated network, optimiser linked to model parameters, and uninstantiated loss**get_out_size** ()

Get number of outputs of model

Return type `int`**Returns** number of outputs of network**get_tail** (*n_in*)

Construct tail module

Return type `Module`**Returns** Instantiated tail `nn.Module`**load_pretrained** (*model*)

Load model weights from pretrained file

Parameters **model** (`Module`) – instantiated model, i.e. return of `build_model()`**Returns** model with weights loaded**set_lr** (*lr*)

Set learning rate for all model parameters

Parameters **lr** (`float`) – learning rate**Return type** `None`

Module contents

4.1.8 lumin.nn.training package

Submodules

lumin.nn.training.fold_train module

```
lumin.nn.training.fold_train.fold_train_ensemble (fy, n_models, bs, model_builder,
                                                  callback_partials=None,
                                                  eval_metrics=None,
                                                  train_on_weights=True,
                                                  eval_on_weights=True, patience=10,
                                                  max_epochs=200, plots=['history',
                                                  'realtime'], shuffle_fold=True,
                                                  shuffle_folds=True, bulk_move=True,
                                                  savepath=PosixPath('train_weights'),
                                                  verbose=False, log_output=False,
                                                  plot_settings=<lumin.plotting.plot_settings.PlotSettings
                                                  object>, callback_args=None)
```

Main training method for *Model*. Trains a specified number of models created by a *ModelBuilder* on data provided by a *FoldYielder*, and save them to *savepath*. Note, this does not return trained models, instead they are saved and must be loaded later. Instead this method returns results of model training. Each *Model* is trained on N-1 folds, for a *FoldYielder* with N folds, and the remaining fold is used as validation data. Training folds are loaded iteratively, and model evaluation takes place after each fold use (a sub-epoch), rather than after every use of all folds (epoch). Training continues until:

- All of the training folds are used *max_epoch* number of times;
- Or validation loss does not decrease for *patience* number of training folds; (or cycles, if using an *AbsCyclicCallback*);
- Or a callback triggers training to stop, e.g. *OneCycle*

Once training is finished, the state with the lowest validation loss is loaded, evaluated, and saved.

Attention: `callback_args` is now deprecated in favour of `callback_partials` and will be removed in v0.4

Parameters

- **fy** (*FoldYielder*) – *FoldYielder* interfacing of training data
- **n_models** (*int*) – number of models to train
- **bs** (*int*) – batch size. Number of data points per iteration
- **model_builder** (*ModelBuilder*) – *ModelBuilder* creating the networks to train
- **callback_partials** (Optional[*List*[*partial*]]) – optional list of functions, each of which will instantiate *Callback* when called
- **eval_metrics** (Optional[*Dict*[*str*, *EvalMetric*]]) – list of instantiated *EvalMetric*. At the end of training, validation data and model predictions will be passed to each, and the results printed and saved
- **train_on_weights** (*bool*) – If weights are present in training data, whether to pass them to the loss function during training
- **eval_on_weights** (*bool*) – If weights are present in validation data, whether to pass them to the loss function during validation
- **patience** (*int*) – number of folds (sub-epochs) or cycles to train without decrease in validation loss before ending training (early stopping)
- **max_epochs** (*int*) – maximum number of epochs for which to train

- **plots** (`List[str]`) – list of string representation of plots to produce. currently: ‘history’: loss history of all models after all training has finished ‘realtime’: live loss evolution during training ‘cycle’: call the plot method of the last (if any) *AbsCyclicCallback* listed in `callback_partials` after every complete model training.
- **shuffle_fold** (`bool`) – whether to tell *BatchYielder* to shuffle data
- **shuffle_folds** (`bool`) – whether to shuffle the order of the trainign folds
- **bulk_move** (`bool`) – whether to pass all training data to device at once, or by minibatch. Bulk moving will be quicker, but may not fit in memory.
- **savepath** (`Path`) – path to to which to save model weights and results
- **verbose** (`bool`) – whether to print out extra information during training
- **log_output** (`bool`) – whether to save printed results to a log file rather than printing them
- **plot_settings** (*PlotSettings*) – *PlotSettings* class to control figure appearance
- **callback_args** (`Optional[List[Dict[str, Any]]]`) – deprecated in favour of `callback_partials`

Return type `Tuple[List[Dict[str, float]], List[Dict[str, List[float]]], List[Dict[str, float]]]`

Returns

- results list of validation losses and other `eval_metrics` results, ordered by model training. Can be used to create an *Ensemble*.
- histories list of loss histories, ordered by model training
- `cycle_losses` if an *AbsCyclicCallback* was passed, list of validation losses at the end of each cycle, ordered by model training. Can be passed to *Ensemble*.

Module contents

4.2 Module contents

LUMIN.OPTIMISATION PACKAGE

5.1 Submodules

5.2 `lumin.optimisation.features` module

`lumin.optimisation.features.get_rf_feat_importance` (*rf*, *inputs*, *targets*, *weights=None*)
Compute feature importance for a Random Forest model using `rfpimp`.

Parameters

- **rf** (`ForestRegressor`) – trained Random Forest model
- **inputs** (`DataFrame`) – input data as Pandas `DataFrame`
- **targets** (`ndarray`) – target data as Numpy array
- **weights** (`Optional[ndarray]`) – Optional data weights as Numpy array

Return type `DataFrame`

`lumin.optimisation.features.rf_rank_features` (*train_df*, *val_df*, *objective*,
train_feats, *targ_name='gen_target'*,
wgt_name=None, *importance_cut=0.0*, *n_estimators=40*,
n_rfs=1, *savename=None*,
plot_settings=<lumin.plotting.plot_settings.PlotSettings object>)

Compute relative permutation importance of input features via using Random Forests. A reduced set of ‘important features’ is obtained by cutting on relative importance and a new model is trained and evaluated on this reduced set. RFs will have their hyper-parameters roughly optimised, both when training on all features and once when training on important features. Relative importances may be computed multiple times (via `n_rfs`) and averaged. In which case the standard error is also computed.

Parameters

- **train_df** (`DataFrame`) – training data as Pandas `DataFrame`
- **val_df** (`DataFrame`) – validation data as Pandas `DataFrame`
- **objective** (`str`) – string representation of objective: either ‘classification’ or ‘regression’
- **train_feats** (`List[str]`) – complete list of training features
- **targ_name** (`str`) – name of column containing target data

- **wgt_name** (Optional[str]) – name of column containing weight data. If set, will use weights for training and evaluation, otherwise will not
- **importance_cut** (float) – minimum importance required to be considered an ‘important feature’
- **n_estimators** (int) – number of trees to use in each forest
- **n_rfs** (int) – number of trainings to perform on all training features in order to compute importances
- **savename** (Optional[str]) – Optional name of file to which to save the plot of feature importances
- **plot_settings** (*PlotSettings*) – *PlotSettings* class to control figure appearance

Return type List[str]

Returns List of features passing importance_cut, ordered by importance

5.3 lumin.optimisation.hyper_param module

```
lumin.optimisation.hyper_param.get_opt_rf_params(x_trn, y_trn, x_val, y_val, objective, w_trn=None, w_val=None,
params={'max_features': [0.3, 0.5, 0.7, 0.9], 'min_samples_leaf': [1, 3, 5, 10, 25, 50, 100]},
n_estimators=40, verbose=True)
```

Use an ordered parameter-scan to roughly optimise Random Forest hyper-parameters.

Parameters

- **x_trn** (ndarray) – training input data
- **y_trn** (ndarray) – training target data
- **x_val** (ndarray) – validation input data
- **y_val** (ndarray) – validation target data
- **objective** (str) – string representation of objective: either ‘classification’ or ‘regression’
- **w_trn** (Optional[ndarray]) – training weights
- **w_val** (Optional[ndarray]) – validation weights
- **params** (OrderedDict) – ordered dictionary mapping parameters to optimise to list of values to consider
- **n_estimators** (int) – number of trees to use in each forest
- **verbose** – Print extra information and show a live plot of model performance

Returns dictionary mapping parameters to their optimised values rf: best performing Random Forest

Return type params

```
lumin.optimisation.hyper_param.fold_lr_find(fy, model_builder, bs,
                                             train_on_weights=True, shuffle_fold=True, n_folds=-1, lr_bounds=[1e-05, 10], callback_partials=None,
                                             plot_settings=<lumin.plotting.plot_settings.PlotSettings object>)
```

Wrapper function for training using *LRFinder* which runs a Smith LR range test (<https://arxiv.org/abs/1803.09820>) using folds in *FoldYielder*. Trains models for 1 fold, interpolating LR between set bounds. This repeats for each fold in *FoldYielder*, and loss evolution is averaged.

Parameters

- **fy** (*FoldYielder*) – *FoldYielder* providing training data
- **model_builder** (*ModelBuilder*) – *ModelBuilder* providing networks and optimisers
- **bs** (*int*) – batch size
- **train_on_weights** (*bool*) – If weights are present, whether to use them for training
- **shuffle_fold** (*bool*) – whether to shuffle data in folds
- **n_folds** (*int*) – if ≥ 1 , will only train *n_folds* number of models, otherwise will train one model per fold
- **lr_bounds** (*Tuple*[*float*, *float*]) – starting and ending LR values
- **callback_partials** (*Optional*[*List*[*partial*]]) – optional list of *func-tools.partial*, each of which will instantiate *Callback* when called
- **plot_settings** (*PlotSettings*) – *PlotSettings* class to control figure appearance

Return type *List*[*LRFinder*]

Returns List of *LRFinder* which were used for each model trained

5.4 lumin.optimisation.threshold module

```
lumin.optimisation.threshold.binary_class_cut(df, top_perc=5.0, min_pred=0.9,
                                              wgt_factor=1.0, br=0.0,
                                              syst_unc_b=0.0, pred_name='pred',
                                              targ_name='gen_target',
                                              wgt_name='gen_weight',
                                              plot_settings=<lumin.plotting.plot_settings.PlotSettings object>)
```

Attention: Deprecated as renamed to *binary_class_cut_by_ams()*. Will be removed in v0.4.

Return type *Tuple*[*float*, *float*, *float*]

```
lumin.optimisation.threshold.binary_class_cut_by_ams (df, top_perc=5.0,
min_pred=0.9,
wgt_factor=1.0, br=0.0,
syst_unc_b=0.0,
pred_name='pred',
targ_name='gen_target',
wgt_name='gen_weight',
plot_settings=<lumin.plotting.plot_settings.PlotSettings
object>)
```

Optimise a cut on a signal-background classifier prediction by the Approximate Median Significance Cut which should generalise better by taking the mean class prediction of the top `top_perc` percentage of points as ranked by AMS

Parameters

- **df** (`DataFrame`) – Pandas `DataFrame` containing data
- **top_perc** (`float`) – top percentage of events to consider as ranked by AMS
- **min_pred** (`float`) – minimum prediction to consider
- **wgt_factor** (`float`) – single multiplicative coefficient for rescaling signal and background weights before computing AMS
- **br** (`float`) – background offset bias
- **syst_unc_b** (`float`) – fractional systematic uncertainty on background
- **pred_name** (`str`) – column to use as predictions
- **targ_name** (`str`) – column to use as truth labels for signal and background
- **wgt_name** (`str`) – column to use as weights for signal and background events
- **plot_settings** (`PlotSettings`) – `PlotSettings` class to control figure appearance

Return type `Tuple[float, float, float]`

Returns Optimised cut AMS at cut Maximum AMS

5.5 Module contents

LUMIN.PLOTTING PACKAGE

6.1 Submodules

6.2 `lumin.plotting.data_viewing` module

```
lumin.plotting.data_viewing.plot_feat(df, feat, wgt_name=None, cuts=None, labels="", plot_bulk=True, n_samples=100000, plot_params=None, size='mid', show_moments=True, ax_labels={'x': None, 'y': 'Density'}, savename=None, settings=<lumin.plotting.plot_settings.PlotSettings object>)
```

A flexible function to provide indicative information about the 1D distribution of a feature. By default it will produce a weighted KDE+histogram for the [1,99] percentile of the data, as well as compute the mean and standard deviation of the data in this region. Distributions are weighted by sampling with replacement the data with probabilities proportional to the sample weights. By passing a list of cuts and labels, it will plot multiple distributions of the same feature for different cuts. Since it is designed to provide quick, indicative information, more specific functions (such as `plot_kdes_from_bs`) should be used to provide final results.

Parameters

- **df** (`DataFrame`) – Pandas DataFrame containing data
- **feat** (`str`) – column name to plot
- **wgt_name** (`Optional[str]`) – if set, will use column to weight data
- **cuts** (`Optional[List[Series]]`) – optional list of cuts to apply to feature. Will add one KDE+hist for each cut listed on the same plot
- **labels** (`Optional[List[str]]`) – optional list of labels for each KDE+hist
- **plot_bulk** (`bool`) – whether to plot the [1,99] percentile of the data, or all of it
- **n_samples** (`int`) – if plotting weighted distributions, how many samples to use
- **plot_params** (`Union[Dict[str, Any], List[Dict[str, Any]], None]`) – optional list of arguments to pass to Seaborn Distplot for each KDE+hist
- **size** (`str`) – string to pass to `str2sz()` to determine size of plot
- **show_moments** (`bool`) – whether to compute and display the mean and standard deviation
- **ax_labels** (`Dict[str, Any]`) – dictionary of x and y axes labels

- **savename** (Optional[str]) – Optional name of file to which to save the plot of feature importances
- **settings** (*PlotSettings*) – *PlotSettings* class to control figure appearance

Return type None

`lumin.plotting.data_viewing.compare_events(events)`

Plot at least two events side by side in their transverse and longitudinal projections

Parameters **events** (list) – list of DataFrames containing vector coordinates for 3 momenta

Return type None

`lumin.plotting.data_viewing.plot_rank_order_dendrogram(df, savename=None, settings=<lumin.plotting.plot_settings.PlotSettings object>)`

Plot dendrogram of features in df clustered via Spearman's rank correlation coefficient

Parameters

- **df** (DataFrame) – Pandas DataFrame containing data
- **savename** (Optional[str]) – Optional name of file to which to save the plot of feature importances
- **settings** (*PlotSettings*) – *PlotSettings* class to control figure appearance

Return type None

`lumin.plotting.data_viewing.plot_kdes_from_bs(x, bs_stats, name2args, feat, units=None, moments=True, savename=None, settings=<lumin.plotting.plot_settings.PlotSettings object>)`

Plot KDEs computed via `bootstrap_stats()`

Parameters

- **bs_stats** (Dict[str, Any]) – (filtered) dictionary returned by `bootstrap_stats()`
- **name2args** (Dict[str, Dict[str, Any]]) – Dictionary mapping names of different distributions to arguments to pass to seaborn tsplot
- **feat** (str) – Name of feature being plotted (for axis labels)
- **units** (Optional[str]) – Optional units to show on axes
- **moments** – whether to display mean and standard deviation of each distribution
- **savename** (Optional[str]) – Optional name of file to which to save the plot of feature importances
- **settings** (*PlotSettings*) – *PlotSettings* class to control figure appearance

Return type None

6.3 lumin.plotting.interpretation module

```
lumin.plotting.interpretation.plot_importance(df, feat_name='Feature',
                                              imp_name='Importance',
                                              unc_name='Uncertainty',
                                              savename=None, settings=
                                              <lumin.plotting.plot_settings.PlotSettings
                                              object>)
```

Plot feature importances as computed via `get_nn_feat_importance`, `get_ensemble_feat_importance`, or `rf_rank_features`

Parameters

- **df** (DataFrame) – DataFrame containing columns of features, importances and, optionally, uncertainties
- **feat_name** (str) – column name for features
- **imp_name** (str) – column name for importances
- **unc_name** (str) – column name for uncertainties (if present)
- **savename** (Optional[str]) – Optional name of file to which to save the plot of feature importances
- **settings** (*PlotSettings*) – *PlotSettings* class to control figure appearance

Return type None

```
lumin.plotting.interpretation.plot_embedding(embed, feat, savename=None, settings=
                                              <lumin.plotting.plot_settings.PlotSettings
                                              object>)
```

Visualise weights in provided categorical entity-embedding matrix

Parameters

- **embed** (OrderedDict) – `state_dict` of trained `nn.Embedding`
- **feat** (str) – name of feature embedded
- **savename** (Optional[str]) – Optional name of file to which to save the plot of feature importances
- **settings** (*PlotSettings*) – *PlotSettings* class to control figure appearance

Return type None

```
lumin.plotting.interpretation.plot_1d_partial_dependence(model, df, feat,
                                                         train_feats, ignore_feats=None,
                                                         input_pipe=None,
                                                         sample_sz=None,
                                                         wgt_name=None,
                                                         n_clusters=10,
                                                         n_points=20,
                                                         pdp_isolate_kargs=None,
                                                         pdp_plot_kargs=None,
                                                         savename=None, settings=
                                                         <lumin.plotting.plot_settings.PlotSettings
                                                         object>)
```

Wrapper for PDPbox to plot 1D dependence of specified feature using provided NN or RF. If features have

been preprocessed using an SK-Learn Pipeline, then that can be passed in order to rescale the x-axis back to its original values.

Parameters

- **model** (*Any*) – any trained model with a `.predict` method
- **df** (*DataFrame*) – *DataFrame* containing training data
- **feat** (*str*) – feature for which to evaluate the partial dependence of the model
- **train_feats** (*List[str]*) – list of all training features including ones which were later ignored, i.e. input features considered when `input_pipe` was fitted
- **ignore_feats** (*Optional[List[str]*) – features present in training data which were not used to train the model (necessary to correctly preprocess feature using `input_pipe`)
- **input_pipe** (*Optional[Pipeline]*) – SK-Learn Pipeline which was used to process the training data
- **sample_sz** (*Optional[int]*) – if set, will only compute partial dependence on a random sample with replacement of the training data, sampled according to weights (if set). Speeds up computation and allows weighted partial dependencies to be computed.
- **wgt_name** (*Optional[str]*) – Optional column name to use as sampling weights
- **n_points** (*int*) – number of points at which to evaluate the model output, passed to `pdp_isolate` as `num_grid_points`
- **n_clusters** (*Optional[int]*) – number of clusters in which to group dependency lines. Set to `None` to show all lines
- **pdp_isolate_kargs** (*Optional[Dict[str, Any]]*) – optional dictionary of keyword arguments to pass to `pdp_isolate`
- **pdp_plot_kargs** (*Optional[Dict[str, Any]]*) – optional dictionary of keyword arguments to pass to `pdp_plot`
- **savename** (*Optional[str]*) – Optional name of file to which to save the plot of feature importances
- **settings** (*PlotSettings*) – *PlotSettings* class to control figure appearance

Return type `None`

```
lumin.plotting.interpretation.plot_2d_partial_dependence(model, df, feats,
                                                         train_feats, ignore_feats=None,
                                                         input_pipe=None,
                                                         sample_sz=None,
                                                         wgt_name=None,
                                                         n_points=[20, 20],
                                                         pdp_interact_kargs=None,
                                                         pdp_interact_plot_kargs=None,
                                                         savename=None, settings=<lumin.plotting.plot_settings.PlotSettings
                                                         object>)
```

Wrapper for `PDPbox` to plot 2D dependence of specified pair of features using provided NN or RF. If features have been preprocessed using an SK-Learn Pipeline, then that can be passed in order to rescale them back to their original values.

Parameters

- **model** (*Any*) – any trained model with a `.predict` method

- **df** (`DataFrame`) – `DataFrame` containing training data
- **feats** (`Tuple[str, str]`) – pair of features for which to evaluate the partial dependence of the model
- **train_feats** (`List[str]`) – list of all training features including ones which were later ignored, i.e. input features considered when `input_pipe` was fitted
- **ignore_feats** (`Optional[List[str]]`) – features present in training data which were not used to train the model (necessary to correctly preprocess feature using `input_pipe`)
- **input_pipe** (`Optional[Pipeline]`) – SK-Learn Pipeline which was used to process the training data
- **sample_sz** (`Optional[int]`) – if set, will only compute partial dependence on a random sample with replacement of the training data, sampled according to weights (if set). Speeds up computation and allows weighted partial dependencies to be computed.
- **wgt_name** (`Optional[str]`) – Optional column name to use as sampling weights
- **n_points** (`Tuple[int, int]`) – pair of numbers of points at which to evaluate the model output, passed to `pdp_interact` as `num_grid_points`
- **n_clusters** – number of clusters in which to group dependency lines. Set to `None` to show all lines
- **pdp_isolate_kargs** – optional dictionary of keyword arguments to pass to `pdp_isolate`
- **pdp_plot_kargs** – optional dictionary of keyword arguments to pass to `pdp_plot`
- **savename** (`Optional[str]`) – Optional name of file to which to save the plot of feature importances
- **settings** (`PlotSettings`) – `PlotSettings` class to control figure appearance

Return type `None`

```
lumin.plotting.interpretation.plot_multibody_weighted_outputs(model, inputs,
                                                             block_names=None,
                                                             use_mean=False,
                                                             save-
                                                             name=None, set-
                                                             tings=<lumin.plotting.plot_settings.PlotSe
                                                             object>)
```

Interpret how a model relies on the outputs of each block in a `:class:MultiBlock` by plotting the outputs of each block as weighted by the tail block. This function currently only supports models whose tail block contains a single neuron in the first dense layer. Input data is passed through the model and the absolute sums of the weighted block outputs are computed per datum, and optionally averaged over the number of block outputs.

Parameters

- **model** (`AbsModel`) – model to interpret
- **inputs** (`Union[ndarray, Tensor]`) – input data to use for interpretation
- **block_names** (`Optional[List[str]]`) – names for each block to use when plotting
- **use_mean** (`bool`) – if `True`, will average the weighted outputs over the number of output neurons in each block
- **savename** (`Optional[str]`) – Optional name of file to which to save the plot of feature importances
- **settings** (`PlotSettings`) – `PlotSettings` class to control figure appearance

Return type None

```
lumin.plotting.interpretation.plot_bottleneck_weighted_inputs(model, bottleneck_idx, inputs, log_y=True, save_name=None, settings=<lumin.plotting.plot_settings.PlotSettings object>)
```

Interpret how a single-neuron bottleneck in a :class:MultiBlock relies on input features by plotting the absolute values of the features times their associated weight for a given set of input data.

Parameters

- **model** (AbsModel) – model to interpret
- **bottleneck_idx** (int) – index of the bottleneck to interpret, i.e. `model.body.bottleneck_blocks[bottleneck_idx]`
- **inputs** (Union[ndarray, Tensor]) – input data to use for interpretation
- **log_y** (bool) – whether to plot a log scale for the y-axis
- **savename** (Optional[str]) – Optional name of file to which to save the plot of feature importances
- **settings** (PlotSettings) – *PlotSettings* class to control figure appearance

Return type None

6.4 lumin.plotting.plot_settings module

```
class lumin.plotting.plot_settings.PlotSettings(**kwargs)
```

Bases: object

Class to provide control over plot appearances. Default parameters are set automatically, and can be adjusted by passing values as keyword arguments during initialisation (or changed after instantiation)

Parameters arguments (*keyword*) – used to set relevant plotting parameters

str2sz (*sz, ax*)

Used to map requested plot sizes to actual dimensions

Parameters

- **sz** (str) – string representation of size
- **ax** (str) – axis dimension requested

Return type float

Returns width of plot dimension

6.5 lumin.plotting.results module

```
lumin.plotting.results.plot_roc(data,      pred_name='pred',      targ_name='gen_target',
                               wgt_name=None,                  labels=None,
                               plot_params=None,               n_bootstrap=0,        log_x=False,
                               plot_baseline=True,              savename=None,        settings=<lumin.plotting.plot_settings.PlotSettings object>)
```

Plot receiver operating characteristic curve(s), optionally using bootstrap resampling

Parameters

- **data** (Union[DataFrame, List[DataFrame]]) – (list of) DataFrame(s) from which to draw predictions and targets
- **pred_name** (str) – name of column to use as predictions
- **targ_name** (str) – name of column to use as targets
- **wgt_name** (Optional[str]) – optional name of column to use as sample weights
- **labels** (Union[str, List[str], None]) – (list of) label(s) for plot legend
- **plot_params** (Union[Dict[str, Any], List[Dict[str, Any]], None]) – (list of) dictionary[ies] of argument(s) to pass to line plot
- **n_bootstrap** (int) – if greater than 0, will bootstrap resample the data that many times when computing the ROC AUC. Currently, this does not affect the shape of the lines, which are based on computing the ROC for the entire dataset as is.
- **log_x** (bool) – whether to use a log scale for plotting the x-axis, useful for high AUC line
- **plot_baseline** (bool) – whether to plot a dotted line for AUC=0.5. Currently incompatible with log_x=True
- **savename** (Optional[str]) – Optional name of file to which to save the plot of feature importances
- **settings** (*PlotSettings*) – *PlotSettings* class to control figure appearance

Return type Dict[str, Union[float, Tuple[float, float]]]

Returns Dictionary mapping data labels to aucs (and uncertainties if n_bootstrap > 0)

```
lumin.plotting.results.plot_binary_class_pred(df,      pred_name='pred',
                                              targ_name='gen_target',
                                              wgt_name=None,      wgt_scale=1,
                                              log_y=False,   lim_x=(0, 1),  den-
                                              sity=True,     savename=None,  set-
                                              tings=<lumin.plotting.plot_settings.PlotSettings
                                              object>)
```

Basic plotter for prediction distribution in a binary classification problem. Note that labels are set using the settings.targ2class dictionary, which by default is {0: 'Background', 1: 'Signal'}.

Parameters

- **df** (DataFrame) – DataFrame with targets and predictions
- **pred_name** (str) – name of column to use as predictions
- **targ_name** (str) – name of column to use as targets
- **wgt_name** (Optional[str]) – optional name of column to use as sample weights

- **wgt_scale** (float) – applies a global multiplicative rescaling to sample weights. Default 1 = no rescaling
- **log_y** (bool) – whether to use a log scale for the y-axis
- **lim_x** (Tuple[float, float]) – limit for plotting on the x-axis
- **density** – whether to normalise each distribution to one, or keep set to sum of weights / datapoints
- **savename** (Optional[str]) – Optional name of file to which to save the plot of feature importances
- **settings** (*PlotSettings*) – *PlotSettings* class to control figure appearance

Return type None

```
lumin.plotting.results.plot_sample_pred(df, pred_name='pred', targ_name='gen_target',
                                       wgt_name='gen_weight', sample_name='gen_sample', wgt_scale=1, bins=35,
                                       log_y=True, lim_x=(0, 1), density=False, zoom_args=None, savename=None,
                                       settings=<lumin.plotting.plot_settings.PlotSettings object>)
```

More advanced plotter for prediction distribution in a binary class problem with stacked distributions for backgrounds and user-defined binning. Can also zoom in to specified parts of plot. Note that plotting colours can be controlled by setting the `settings.sample2col` dictionary.

Parameters

- **df** (DataFrame) – DataFrame with targets and predictions
- **pred_name** (str) – name of column to use as predictions
- **targ_name** (str) – name of column to use as targets
- **wgt_name** (str) – name of column to use as sample weights
- **sample_name** (str) – name of column to use as process names
- **wgt_scale** (float) – applies a global multiplicative rescaling to sample weights. Default 1 = no rescaling
- **bins** (Union[int, List[int]]) – either the number of bins to use for a uniform binning, or a list of bin edges for a variable-width binning
- **log_y** (bool) – whether to use a log scale for the y-axis
- **lim_x** (Tuple[float, float]) – limit for plotting on the x-axis
- **density** – whether to normalise each distribution to one, or keep set to sum of weights / datapoints
- **zoom_args** (Optional[Dict[str, Any]]) – arguments to control the optional zoomed in section, e.g. {'x':(0.4,0.45), 'y':(0.2, 1500), 'anchor':(0,0.25,0.95,1), 'width_scale':1, 'width_zoom':4, 'height_zoom':3}
- **savename** (Optional[str]) – Optional name of file to which to save the plot of feature importances
- **settings** (*PlotSettings*) – *PlotSettings* class to control figure appearance

Return type None

6.6 lumin.plotting.training module

```
lumin.plotting.training.plot_train_history (histories, savename=None,
                                           ignore_trn=True, settings=
                                           <lumin.plotting.plot_settings.PlotSettings
                                           object>)
```

Plot histories object returned by `fold_train_ensemble()` showing the loss evolution over time per model trained.

Parameters

- **histories** (List[Dict[str, List[float]]]) – list of dictionaries mapping loss type to values at each (sub)-epoch
- **savename** (Optional[str]) – Optional name of file to which to save the plot of feature importances
- **ignore_trn** – whether to ignore training loss
- **settings** (*PlotSettings*) – *PlotSettings* class to control figure appearance

Return type None

```
lumin.plotting.training.plot_lr_finders (lr_finders, loss='loss', cut=-10,
                                         settings=
                                         <lumin.plotting.plot_settings.PlotSettings
                                         object>)
```

Plot mean loss evolution against learning rate for several `fold_lr_find`.

Parameters

- **lr_finders** (List[*LRFinder*]) – list of `fold_lr_find`
- **loss** – string representation of loss to plot
- **cut** – number of final iterations to cut
- **settings** (*PlotSettings*) – *PlotSettings* class to control figure appearance

Return type None

6.7 Module contents

LUMIN.UTILS PACKAGE

7.1 Submodules

7.2 `lumin.utils.data` module

`lumin.utils.data.check_val_set` (*train*, *val*, *test=None*, *n_folds=None*)

Method to check validation set suitability by seeing whether Random Forests can predict whether events belong to one dataset or another. If a *FoldYielder* is passed, then trainings are run once per fold and averaged. Will compute the ROC AUC for set discrimination (should be close to 0.5) and compute the feature importances to aid removal of discriminating features.

Parameters

- **train** (Union[DataFrame, ndarray, *FoldYielder*]) – training data
- **val** (Union[DataFrame, ndarray, *FoldYielder*]) – validation data
- **test** (Union[DataFrame, ndarray, *FoldYielder*, None]) – optional testing data
- **n_folds** (Optional[int]) – if set and if passed a *FoldYielder*, will only use the first *n_folds* folds

Return type None

7.3 `lumin.utils.misc` module

`lumin.utils.misc.to_np` (*x*)

Convert Tensor *x* to a Numpy array

Parameters **x** (Tensor) – Tensor to convert

Return type ndarray

Returns *x* as a Numpy array

`lumin.utils.misc.to_device` (*x*, *device=device(type='cpu')*)

Recursively place Tensor(s) onto device

Parameters **x** (Union[Tensor, List[Tensor]]) – Tensor(s) to place on device

Return type Union[Tensor, List[Tensor]]

Returns Tensor(s) on device

`lumin.utils.misc.to_tensor` (*x*)

Convert Numpy array to Tensor with possibility of a None being passed

Parameters `x` (Optional[ndarray]) – Numpy array or None

Return type Optional[Tensor]

Returns `x` as Tensor or None

`lumin.utils.misc.str2bool` (*string*)

Convert string representation of Boolean to bool

Parameters `string` (Union[str, bool]) – string representation of Boolean (or a Boolean)

Return type bool

Returns bool if bool was passed else, True if lowercase string matches is in (“yes”, “true”, “t”, “1”)

`lumin.utils.misc.to_binary_class` (*df, zero_preds, one_preds*)

Map class precitions back to a binary prediction. The maximum prediction for features listed in `zero_preds` is treated as the prediction for class 0, vice versa for `one_preds`. The binary prediction is added to `df` in place as column ‘pred’

Parameters

- `df` (DataFrame) – DataFrame containing prediction features
- `zero_preds` (List[str]) – list of column names for predictions associated with class 0
- `one_preds` (List[str]) – list of column names for predictions associated with class 0

Return type None

`lumin.utils.misc.ids2unique` (*ids*)

Map a permutaion of integers to a unique number, or a 2D array of integers to unique numbers by row. Returned numbers are unique for a given permutation of integers. This is achieved by computing the product of primes raised to powers equal to the integers. Beacause of this, it can be easy to produce numbers which are too large to be stored if many (large) integers are passed.

Parameters `ids` (Union[List[int], ndarray]) – (array of) permutation(s) of integers to map

Return type ndarray

Returns (Array of) unique id(s) for given permutation(s)

`class lumin.utils.misc.FowardHook` (*module, hook_fn=None*)

Bases: object

Create a hook for performing an action based on the forward pass thourgh a `nn.Module`

Parameters

- `module` – `nn.Module` to hook
- `hook_fn` – Optional function to perform. Default is to record input and output of module

Examples::

```
>>> hook = ForwardHook(model.tail.dense)
>>> model.predict(inputs)
>>> print(hook.inputs)
```

`hook_fn` (*module, input, output*)

Default hook function records inputs and outputs of module

Parameters

- `module` (Module) – `nn.Module` to hook

- **input** (Union[Tensor, Tuple[Tensor]]) – input tensor
- **output** (Union[Tensor, Tuple[Tensor]]) – output tensor of module

Return type None

remove ()

Call when finished to remove hook

Return type None

7.4 lumin.utils.multiprocessing module

`lumin.utils.multiprocessing.mp_run (args, func)`

Run multiple instances of function simultaneously by using a list of argument dictionaries Runs given function once per entry in args list.

Important: Function should put a dictionary of results into the `mp.Queue` and each result key should be unique otherwise they will overwrite one another.

Parameters

- **args** (List[Dict[Any, Any]]) – list of dictionaries of arguments
- **func** (Callable[[Any], Any]) – function to which to pass dictionary arguments

Return type Dict[Any, Any]

Returns Dictionary of results

7.5 lumin.utils.statistics module

`lumin.utils.statistics.bootstrap_stats (args, out_q=None)`

Computes statistics and KDEs of data via sampling with replacement

Parameters

- **args** (Dict[str, Any]) – dictionary of arguments. Possible keys are: name - name prepended to returned keys in result dict weights - array of weights matching length of data to use for weighted resampling n - number of times to resample data x - points at which to compute the kde values of resample data kde - whether to compute the kde values at x-points for resampled data mean - whether to compute the means of the resampled data std - whether to compute standard deviation of resampled data c68 - whether to compute the width of the absolute central 68.2 percentile of the resampled data
- **out_q** (Optional[<bound method BaseContext.Queue of <multiprocessing.context.DefaultContext object at 0x7f5b729d75c0>>]) – if using multiprocessing can place result dictionary in provided queue

Return type Union[None, Dict[str, Any]]

Returns Result dictionary if `out_q` is `None` else `None`.

`lumin.utils.statistics.get_moments (arr)`

Computes mean and std of data, and their associated uncertainties

Parameters **arr** (ndarray) – univariate data

Return type Tuple[float, float, float, float]

Returns

- mean
- statistical uncertainty of mean
- standard deviation
- statistical uncertainty of standard deviation

`lumin.utils.statistics.uncert_round` (*value*, *uncert*)

Round value according to given uncertainty using one significant figure of the uncertainty

Parameters

- **value** (float) – value to round
- **uncert** (float) – uncertainty of value

Return type Tuple[float, float]

Returns

- rounded value
- rounded uncertainty

7.6 Module contents

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