Chapter_3_A_Tour_of_Machine_Learning_Classifiers_Using_Scikit-Learn

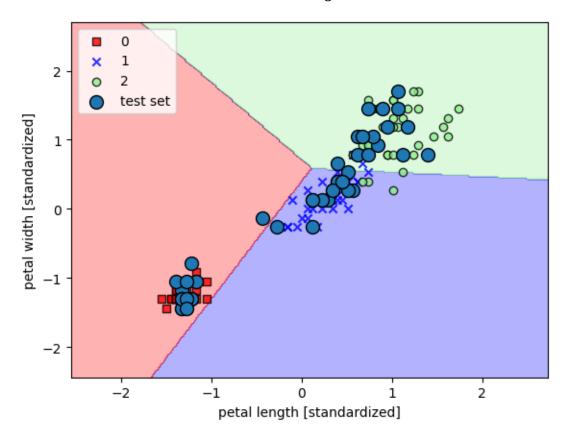
March 17, 2024

0.1 Implementing Perceptron using scikit-learn on iris dataset

```
[34]: from sklearn import datasets
      import numpy as np
      iris = datasets.load_iris()
      X = iris.data[:, [2, 3]]
      y = iris.target
      print('Class labels:', np.unique(y))
     Class labels: [0 1 2]
[35]: # iris
[36]: from sklearn.model_selection import train_test_split
      X train, X test, y train, y test = train_test_split(X, y, test_size=0.3,_
       →random_state=1, stratify=y)
[37]: print('Labels counts in y:', np.bincount(y))
      print('Labels counts in y_train:', np.bincount(y_train))
      print('Labels counts in y_test:', np.bincount(y_test))
     Labels counts in y: [50 50 50]
     Labels counts in y_train: [35 35 35]
     Labels counts in y_test: [15 15 15]
[38]: from sklearn.preprocessing import StandardScaler
      sc = StandardScaler()
      sc.fit(X_train)
      X_train_std = sc.transform(X_train)
      X_test_std = sc.transform(X_test)
[39]: from sklearn.linear model import Perceptron
      ppn = Perceptron(max_iter=40, eta0=0.1, random_state=1)
      ppn.fit(X_train_std, y_train)
[39]: Perceptron(eta0=0.1, max_iter=40, random_state=1)
```

```
[40]: y_pred = ppn.predict(X_test_std)
      print('Misclassified samples: %d' % (y_test != y_pred).sum())
     Misclassified samples: 1
[41]: from sklearn.metrics import accuracy_score
      print('Accuracy: %.2f' % accuracy_score(y_test, y_pred))
     Accuracy: 0.98
[42]: print('Accuracy: %.2f' % ppn.score(X_test_std, y_test))
     Accuracy: 0.98
[43]: from matplotlib.colors import ListedColormap
      import matplotlib.pyplot as plt
      def plot_decision_regions(X, y, classifier, test_idx=None,resolution=0.02):
          # setup marker generator and color map
         markers = ('s', 'x', 'o', '^', 'v')
          colors = ('red', 'blue', 'lightgreen', 'gray', 'cyan')
          cmap = ListedColormap(colors[:len(np.unique(y))])
          # plot the decision surface
         x1 \text{ min}, x1 \text{ max} = X[:, 0].min() - 1, X[:, 0].max() + 1
         x2_{min}, x2_{max} = X[:, 1].min() - 1, X[:, 1].max() + 1
         xx1, xx2 = np.meshgrid(np.arange(x1_min, x1_max, resolution),
         np.arange(x2_min, x2_max, resolution))
         Z = classifier.predict(np.array([xx1.ravel(), xx2.ravel()]).T)
         Z = Z.reshape(xx1.shape)
         plt.contourf(xx1, xx2, Z, alpha=0.3, cmap=cmap)
         plt.xlim(xx1.min(), xx1.max())
         plt.ylim(xx2.min(), xx2.max())
         for idx, cl in enumerate(np.unique(y)):
             plt.scatter(x=X[y == cl, 0], y=X[y == cl, 1],alpha=0.8, 
       →c=colors[idx], marker=markers[idx], label=cl,edgecolor='black')
          # highlight test samples
         if test idx:
              # plot all samples
             X_test, y_test = X[test_idx, :], y[test_idx]
             plt.scatter(X_test[:, 0], X_test[:, 1], edgecolor='black', alpha=1.
       [44]: X_combined_std = np.vstack((X_train_std, X_test_std))
      y_combined = np.hstack((y_train, y_test))
      plot_decision_regions(X=X_combined_std,y=y_combined,classifier=ppn,test_idx=range(105,__
       →150))
      plt.xlabel('petal length [standardized]')
      plt.ylabel('petal width [standardized]')
      plt.legend(loc='upper left')
      plt.show()
```

plt.scatter(x=X[y == cl, 0], y=X[y == cl, 1],alpha=0.8,
c=colors[idx],marker=markers[idx], label=cl,edgecolor='black')



[45]: $\# \# \ decision \ boundary \ is \ not \ perfectly \ separable \ because \ data \ classes \ are \ not_{\sqcup}$ $\hookrightarrow linearly \ separable. \ So, \ use \ logistic \ algorithm$

[46]: help(Perceptron)

Help on class Perceptron in module sklearn.linear_model._perceptron:

class Perceptron(sklearn.linear_model._stochastic_gradient.BaseSGDClassifier)
 | Perceptron(*, penalty=None, alpha=0.0001, l1_ratio=0.15, fit_intercept=True,
max_iter=1000, tol=0.001, shuffle=True, verbose=0, eta0=1.0, n_jobs=None,
random_state=0, early_stopping=False, validation_fraction=0.1,
n_iter_no_change=5, class_weight=None, warm_start=False)
 |
 | Linear perceptron classifier.

```
| The implementation is a wrapper around
:class:`~sklearn.linear_model.SGDClassifier`
by fixing the `loss` and `learning_rate` parameters as::
       SGDClassifier(loss="perceptron", learning_rate="constant")
   Other available parameters are described below and are forwarded to
   :class:`~sklearn.linear_model.SGDClassifier`.
  Read more in the :ref:`User Guide <perceptron>`.
  Parameters
   _____
   penalty : {'12','11','elasticnet'}, default=None
       The penalty (aka regularization term) to be used.
   alpha: float, default=0.0001
       Constant that multiplies the regularization term if regularization is
       used.
   l1_ratio : float, default=0.15
       The Elastic Net mixing parameter, with `0 <= 11_ratio <= 1`.
       `l1_ratio=0` corresponds to L2 penalty, `l1_ratio=1` to L1.
       Only used if `penalty='elasticnet'`.
       .. versionadded:: 0.24
   fit_intercept : bool, default=True
       Whether the intercept should be estimated or not. If False, the
       data is assumed to be already centered.
  max_iter : int, default=1000
       The maximum number of passes over the training data (aka epochs).
       It only impacts the behavior in the ``fit`` method, and not the
       :meth:`partial_fit` method.
       .. versionadded:: 0.19
   tol : float or None, default=1e-3
       The stopping criterion. If it is not None, the iterations will stop
       when (loss > previous_loss - tol).
       .. versionadded:: 0.19
   shuffle : bool, default=True
       Whether or not the training data should be shuffled after each epoch.
```

```
verbose : int, default=0
     The verbosity level.
eta0 : float, default=1
     Constant by which the updates are multiplied.
n_jobs : int, default=None
     The number of CPUs to use to do the OVA (One Versus All, for
     multi-class problems) computation.
     ``None`` means 1 unless in a :obj:`joblib.parallel_backend` context.
     ``-1`` means using all processors. See :term:`Glossary <n_jobs>`
     for more details.
 random_state : int, RandomState instance or None, default=0
     Used to shuffle the training data, when ``shuffle`` is set to
     ``True``. Pass an int for reproducible output across multiple
     function calls.
     See :term:`Glossary <random_state>`.
 early_stopping : bool, default=False
     Whether to use early stopping to terminate training when validation
     score is not improving. If set to True, it will automatically set aside
     a stratified fraction of training data as validation and terminate
     training when validation score is not improving by at least `tol` for
     `n_iter_no_change` consecutive epochs.
     .. versionadded:: 0.20
 validation_fraction : float, default=0.1
     The proportion of training data to set aside as validation set for
     early stopping. Must be between 0 and 1.
     Only used if early_stopping is True.
     .. versionadded:: 0.20
 n_iter_no_change : int, default=5
     Number of iterations with no improvement to wait before early stopping.
     .. versionadded:: 0.20
class_weight : dict, {class_label: weight} or "balanced", default=None
     Preset for the class_weight fit parameter.
     Weights associated with classes. If not given, all classes
     are supposed to have weight one.
     The "balanced" mode uses the values of y to automatically adjust
     weights inversely proportional to class frequencies in the input data
```

```
as ``n_samples / (n_classes * np.bincount(y))``.
  warm_start : bool, default=False
       When set to True, reuse the solution of the previous call to fit as
       initialization, otherwise, just erase the previous solution. See
       :term:`the Glossary <warm_start>`.
 Attributes
  -----
  classes_ : ndarray of shape (n_classes,)
       The unique classes labels.
| coef_ : ndarray of shape (1, n_features) if n_classes == 2 else
(n_classes, n_features)
       Weights assigned to the features.
   intercept_ : ndarray of shape (1,) if n_classes == 2 else (n_classes,)
       Constants in decision function.
   loss_function_ : concrete LossFunction
       The function that determines the loss, or difference between the
       output of the algorithm and the target values.
  n_features_in_ : int
       Number of features seen during :term:`fit`.
       .. versionadded:: 0.24
   feature_names_in_ : ndarray of shape (`n_features_in_`,)
       Names of features seen during :term:`fit`. Defined only when `X`
       has feature names that are all strings.
       .. versionadded:: 1.0
  n iter : int
       The actual number of iterations to reach the stopping criterion.
       For multiclass fits, it is the maximum over every binary fit.
  t_ : int
       Number of weight updates performed during training.
       Same as ``(n_iter_ * n_samples + 1)``.
| See Also
   sklearn.linear_model.SGDClassifier : Linear classifiers
       (SVM, logistic regression, etc.) with SGD training.
   Notes
```

```
``Perceptron`` is a classification algorithm which shares the same
 underlying implementation with ``SGDClassifier``. In fact,
   ``Perceptron()`` is equivalent to `SGDClassifier(loss="perceptron",
   eta0=1, learning_rate="constant", penalty=None)`.
 References
   _____
 https://en.wikipedia.org/wiki/Perceptron and references therein.
 | Examples
  -----
 | >>> from sklearn.linear_model import Perceptron
 >>> X, y = load_digits(return_X_y=True)
 >>> clf = Perceptron(tol=1e-3, random_state=0)
 | >>> clf.fit(X, y)
 | Perceptron()
 | >>> clf.score(X, y)
 I 0.939...
   Method resolution order:
       Perceptron
       sklearn.linear_model._stochastic_gradient.BaseSGDClassifier
       sklearn.linear_model._base.LinearClassifierMixin
       sklearn.base.ClassifierMixin
       sklearn.linear_model._stochastic_gradient.BaseSGD
       sklearn.linear_model._base.SparseCoefMixin
       sklearn.base.BaseEstimator
       sklearn.utils._estimator_html_repr._HTMLDocumentationLinkMixin
       sklearn.utils._metadata_requests._MetadataRequester
       builtins.object
  Methods defined here:
 __init__(self, *, penalty=None, alpha=0.0001, l1_ratio=0.15,
fit_intercept=True, max_iter=1000, tol=0.001, shuffle=True, verbose=0, eta0=1.0,
n_jobs=None, random_state=0, early_stopping=False, validation_fraction=0.1,
n_iter_no_change=5, class_weight=None, warm_start=False)
       Initialize self. See help(type(self)) for accurate signature.
 | set_fit_request(self: sklearn.linear_model._perceptron.Perceptron, *,
coef_init: Union[bool, NoneType, str] = '$UNCHANGED$', intercept_init:
Union[bool, NoneType, str] = '$UNCHANGED$', sample_weight: Union[bool, NoneType,
str] = '$UNCHANGED$') -> sklearn.linear_model._perceptron.Perceptron
       Request metadata passed to the ``fit`` method.
       Note that this method is only relevant if
```

```
``enable_metadata_routing=True`` (see :func:`sklearn.set_config`).
        Please see :ref: `User Guide <metadata_routing>` on how the routing
        mechanism works.
       The options for each parameter are:
        - ``True``: metadata is requested, and passed to ``fit`` if provided.
The request is ignored if metadata is not provided.
        - ``False``: metadata is not requested and the meta-estimator will not
pass it to ``fit``.
        - ``None``: metadata is not requested, and the meta-estimator will raise
an error if the user provides it.
        - ``str``: metadata should be passed to the meta-estimator with this
given alias instead of the original name.
        The default (``sklearn.utils.metadata_routing.UNCHANGED``) retains the
        existing request. This allows you to change the request for some
        parameters and not others.
        .. versionadded:: 1.3
        .. note::
            This method is only relevant if this estimator is used as a
            sub-estimator of a meta-estimator, e.g. used inside a
            :class:`~sklearn.pipeline.Pipeline`. Otherwise it has no effect.
        Parameters
        coef_init : str, True, False, or None,
default=sklearn.utils.metadata_routing.UNCHANGED
           Metadata routing for ``coef_init`` parameter in ``fit``.
        intercept_init : str, True, False, or None,
default=sklearn.utils.metadata_routing.UNCHANGED
           Metadata routing for ``intercept_init`` parameter in ``fit``.
        sample_weight : str, True, False, or None,
default=sklearn.utils.metadata_routing.UNCHANGED
            Metadata routing for ``sample_weight`` parameter in ``fit``.
       Returns
        -----
        self : object
            The updated object.
```

```
| set_partial_fit_request(self: sklearn.linear_model._perceptron.Perceptron,
*, classes: Union[bool, NoneType, str] = '$UNCHANGED$', sample_weight:
Union[bool, NoneType, str] = '$UNCHANGED$') ->
sklearn.linear_model._perceptron.Perceptron
        Request metadata passed to the ``partial_fit`` method.
        Note that this method is only relevant if
        ``enable_metadata_routing=True`` (see :func:`sklearn.set_config`).
       Please see :ref:`User Guide <metadata_routing>` on how the routing
       mechanism works.
       The options for each parameter are:
        - ``True``: metadata is requested, and passed to ``partial_fit`` if
provided. The request is ignored if metadata is not provided.
        - ``False``: metadata is not requested and the meta-estimator will not
pass it to ``partial_fit``.
        - ``None``: metadata is not requested, and the meta-estimator will raise
an error if the user provides it.
        - ``str``: metadata should be passed to the meta-estimator with this
given alias instead of the original name.
        The default (``sklearn.utils.metadata_routing.UNCHANGED``) retains the
        existing request. This allows you to change the request for some
        parameters and not others.
        .. versionadded:: 1.3
        .. note::
            This method is only relevant if this estimator is used as a
            sub-estimator of a meta-estimator, e.g. used inside a
            :class:`~sklearn.pipeline.Pipeline`. Otherwise it has no effect.
       Parameters
        classes : str, True, False, or None,
default=sklearn.utils.metadata_routing.UNCHANGED
           Metadata routing for ``classes`` parameter in ``partial_fit``.
        sample_weight : str, True, False, or None,
default=sklearn.utils.metadata_routing.UNCHANGED
           Metadata routing for ``sample_weight`` parameter in ``partial_fit``.
       Returns
```

```
self : object
            The updated object.
 | set_score_request(self: sklearn.linear_model._perceptron.Perceptron, *,
sample weight: Union[bool, NoneType, str] = '$UNCHANGED$') ->
sklearn.linear_model._perceptron.Perceptron
        Request metadata passed to the ``score`` method.
        Note that this method is only relevant if
        ``enable_metadata_routing=True`` (see :func:`sklearn.set_config`).
        Please see :ref:`User Guide <metadata routing>` on how the routing
        mechanism works.
       The options for each parameter are:
        - ``True``: metadata is requested, and passed to ``score`` if provided.
The request is ignored if metadata is not provided.
        - ``False``: metadata is not requested and the meta-estimator will not
pass it to ``score``.
        - ``None``: metadata is not requested, and the meta-estimator will raise
an error if the user provides it.
        - ``str``: metadata should be passed to the meta-estimator with this
given alias instead of the original name.
        The default (``sklearn.utils.metadata_routing.UNCHANGED``) retains the
        existing request. This allows you to change the request for some
       parameters and not others.
        .. versionadded:: 1.3
        .. note::
            This method is only relevant if this estimator is used as a
            sub-estimator of a meta-estimator, e.g. used inside a
            :class:`~sklearn.pipeline.Pipeline`. Otherwise it has no effect.
       Parameters
        sample_weight : str, True, False, or None,
default=sklearn.utils.metadata_routing.UNCHANGED
            Metadata routing for ``sample_weight`` parameter in ``score``.
       Returns
        -----
        self : object
            The updated object.
```

```
Data and other attributes defined here:
   __abstractmethods__ = frozenset()
    __annotations__ = {'_parameter_constraints': <class 'dict'>}
   Methods inherited from
sklearn.linear_model._stochastic_gradient.BaseSGDClassifier:
   fit(self, X, y, coef_init=None, intercept_init=None, sample_weight=None)
       Fit linear model with Stochastic Gradient Descent.
       Parameters
       X : {array-like, sparse matrix}, shape (n_samples, n_features)
            Training data.
       y : ndarray of shape (n_samples,)
            Target values.
       coef_init : ndarray of shape (n_classes, n_features), default=None
            The initial coefficients to warm-start the optimization.
        intercept_init : ndarray of shape (n_classes,), default=None
            The initial intercept to warm-start the optimization.
        sample_weight : array-like, shape (n_samples,), default=None
            Weights applied to individual samples.
            If not provided, uniform weights are assumed. These weights will
            be multiplied with class_weight (passed through the
            constructor) if class_weight is specified.
       Returns
        self : object
           Returns an instance of self.
   partial_fit(self, X, y, classes=None, sample_weight=None)
       Perform one epoch of stochastic gradient descent on given samples.
       Internally, this method uses ``max_iter = 1``. Therefore, it is not
       guaranteed that a minimum of the cost function is reached after calling
       it once. Matters such as objective convergence, early stopping, and
       learning rate adjustments should be handled by the user.
```

```
Parameters
       _____
       X : {array-like, sparse matrix}, shape (n_samples, n_features)
           Subset of the training data.
       y : ndarray of shape (n_samples,)
           Subset of the target values.
       classes : ndarray of shape (n_classes,), default=None
           Classes across all calls to partial_fit.
           Can be obtained by via `np.unique(y_all)`, where y_all is the
           target vector of the entire dataset.
           This argument is required for the first call to partial_fit
           and can be omitted in the subsequent calls.
           Note that y doesn't need to contain all labels in `classes`.
       sample_weight : array-like, shape (n_samples,), default=None
           Weights applied to individual samples.
           If not provided, uniform weights are assumed.
       Returns
       _____
       self : object
           Returns an instance of self.
   Data and other attributes inherited from
sklearn.linear_model._stochastic_gradient.BaseSGDClassifier:
   loss_functions = {'epsilon_insensitive': (<class 'sklearn.linear_model...
   _____
  Methods inherited from sklearn.linear_model._base.LinearClassifierMixin:
   decision_function(self, X)
       Predict confidence scores for samples.
       The confidence score for a sample is proportional to the signed
       distance of that sample to the hyperplane.
       Parameters
       X : {array-like, sparse matrix} of shape (n_samples, n_features)
           The data matrix for which we want to get the confidence scores.
       Returns
       _____
       scores : ndarray of shape (n_samples,) or (n_samples, n_classes)
```

```
Confidence scores per `(n_samples, n_classes)` combination. In the
         binary case, confidence score for `self.classes_[1]` where >0 means
         this class would be predicted.
predict(self, X)
     Predict class labels for samples in X.
     Parameters
     X : {array-like, sparse matrix} of shape (n_samples, n_features)
         The data matrix for which we want to get the predictions.
     Returns
     _____
     y_pred : ndarray of shape (n_samples,)
         Vector containing the class labels for each sample.
Methods inherited from sklearn.base.ClassifierMixin:
 score(self, X, y, sample_weight=None)
     Return the mean accuracy on the given test data and labels.
     In multi-label classification, this is the subset accuracy
     which is a harsh metric since you require for each sample that
     each label set be correctly predicted.
     Parameters
     X : array-like of shape (n_samples, n_features)
         Test samples.
     y : array-like of shape (n_samples,) or (n_samples, n_outputs)
         True labels for `X`.
     sample_weight : array-like of shape (n_samples,), default=None
         Sample weights.
    Returns
     -----
     score : float
         Mean accuracy of ``self.predict(X)`` w.r.t. `y`.
Data descriptors inherited from sklearn.base.ClassifierMixin:
__dict__
     dictionary for instance variables (if defined)
```

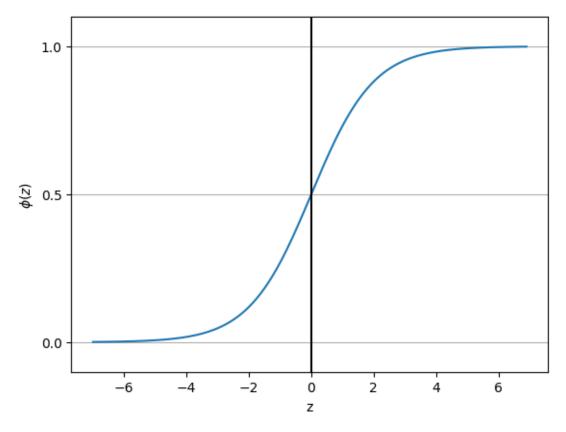
```
__weakref__
        list of weak references to the object (if defined)
   Readonly properties inherited from
sklearn.linear_model._stochastic_gradient.BaseSGD:
   loss_function_
  Methods inherited from sklearn.linear_model._base.SparseCoefMixin:
   densify(self)
        Convert coefficient matrix to dense array format.
        Converts the ``coef_`` member (back) to a numpy.ndarray. This is the
        default format of ``coef_`` and is required for fitting, so calling
        this method is only required on models that have previously been
        sparsified; otherwise, it is a no-op.
       Returns
        ____
        self
           Fitted estimator.
   sparsify(self)
        Convert coefficient matrix to sparse format.
        Converts the ``coef_`` member to a scipy.sparse matrix, which for
        L1-regularized models can be much more memory- and storage-efficient
        than the usual numpy.ndarray representation.
       The ``intercept_`` member is not converted.
       Returns
        self
           Fitted estimator.
       Notes
       For non-sparse models, i.e. when there are not many zeros in ``coef_``,
       this may actually *increase* memory usage, so use this method with
        care. A rule of thumb is that the number of zero elements, which can
        be computed with ``(coef_ == 0).sum()``, must be more than 50% for this
        to provide significant benefits.
```

```
After calling this method, further fitting with the partial_fit
     method (if any) will not work until you call densify.
Methods inherited from sklearn.base.BaseEstimator:
 __getstate__(self)
 __repr__(self, N_CHAR_MAX=700)
    Return repr(self).
 __setstate__(self, state)
 __sklearn_clone__(self)
get_params(self, deep=True)
     Get parameters for this estimator.
     Parameters
     _____
     deep : bool, default=True
         If True, will return the parameters for this estimator and
         contained subobjects that are estimators.
     Returns
     _____
     params : dict
         Parameter names mapped to their values.
 set_params(self, **params)
     Set the parameters of this estimator.
     The method works on simple estimators as well as on nested objects
     (such as :class:`~sklearn.pipeline.Pipeline`). The latter have
     parameters of the form ``<component>__<parameter>`` so that it's
     possible to update each component of a nested object.
     Parameters
     **params : dict
         Estimator parameters.
     Returns
     self : estimator instance
         Estimator instance.
```

```
Methods inherited from sklearn.utils._metadata_requests._MetadataRequester:
   get_metadata_routing(self)
        Get metadata routing of this object.
        Please check :ref:`User Guide <metadata_routing>` on how the routing
       mechanism works.
       Returns
       routing : MetadataRequest
            A :class: `~sklearn.utils.metadata_routing.MetadataRequest`
encapsulating
           routing information.
                              _____
   Class methods inherited from
sklearn.utils._metadata_requests._MetadataRequester:
    __init_subclass__(**kwargs) from abc.ABCMeta
        Set the ``set_{method}_request`` methods.
       This uses PEP-487 [1] to set the ``set_{method}_request`` methods. It
        looks for the information available in the set default values which are
        set using ``__metadata_request__*`` class attributes, or inferred
        from method signatures.
       The ``__metadata_request__*`` class attributes are used when a method
        does not explicitly accept a metadata through its arguments or if the
        developer would like to specify a request value for those metadata
        which are different from the default ``None``.
       References
        .. [1] https://www.python.org/dev/peps/pep-0487
```

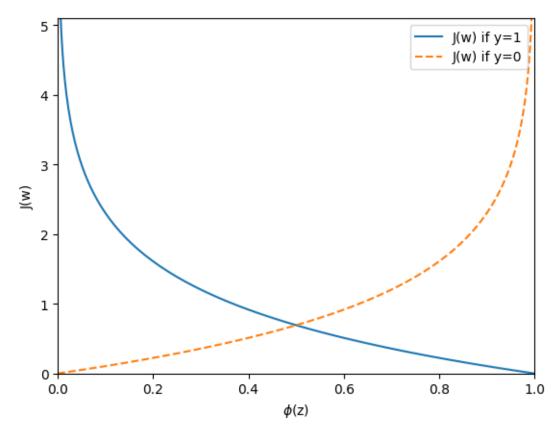
0.2 Implementing Logistic Regression using scikit-learn on iris dataset

```
phi_z = sigmoid(z)
plt.plot(z, phi_z)
plt.axvline(0.0, color='k')
plt.ylim(-0.1, 1.1)
plt.xlabel('z')
plt.ylabel('$\phi (z)$')
# y axis ticks and gridline
plt.yticks([0.0, 0.5, 1.0])
ax = plt.gca()
ax.yaxis.grid(True)
plt.show()
```



```
[48]: def cost_1(z):
    return - np.log(sigmoid(z))
def cost_0(z):
    return - np.log(1 - sigmoid(z))
z = np.arange(-10, 10, 0.1)
phi_z = sigmoid(z)
c1 = [cost_1(x) for x in z]
plt.plot(phi_z, c1, label='J(w) if y=1')
c0 = [cost_0(x) for x in z]
```

```
plt.plot(phi_z, c0, linestyle='--', label='J(w) if y=0')
plt.ylim(0.0, 5.1)
plt.xlim([0, 1])
plt.xlabel('$\phi$(z)')
plt.ylabel('J(w)')
plt.legend(loc='best')
plt.show()
```



0.2.1 Implementing Logistic Regression by modifying Adaline

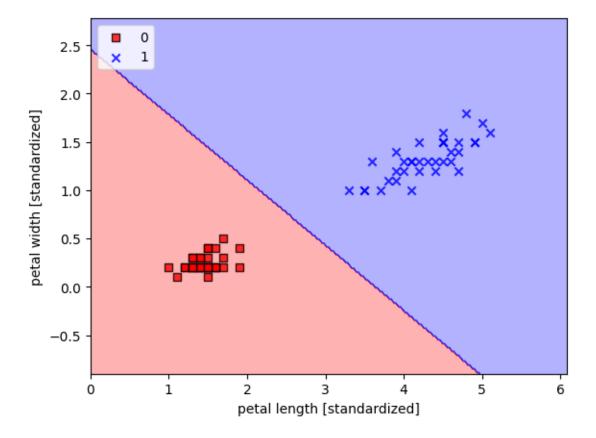
```
[49]: class Logistic Regression GD (object):
    """Logistic Regression Classifier using gradient descent.
    Parameters
    ------
    eta: float
    Learning rate (between 0.0 and 1.0)
    n_iter: int
    Passes over the training dataset.
    random_state: int
    Random number generator seed for random weight
```

```
initialization.
Attributes
_____
w_{\perp}: 1d-array
Weights after fitting.
cost_{-} : list
Sum-of-squares cost function value in each epoch.
def __init__(self, eta=0.05, n_iter=100, random_state=1):
    self.eta = eta
    self.n_iter = n_iter
    self.random_state = random_state
def fit(self, X, y):
    """ Fit training data.
    Parameters
    X : {array-like}, shape = [n_samples, n_features]
    Training vectors, where n_samples is the number of
    samples and
    n_features is the number of features.
    y : array-like, shape = [n_samples]
    Target values.
    Returns
    self : object
    rgen = np.random.RandomState(self.random_state)
    self.w_ = rgen.normal(loc=0.0, scale=0.01, size=1 + X.shape[1])
    self.cost_ = []
    for i in range(self.n_iter):
        net_input = self.net_input(X)
        output = self.activation(net_input)
        errors = (y - output)
        self.w_[1:] += self.eta * X.T.dot(errors)
        self.w_[0] += self.eta * errors.sum()
        # note that we compute the logistic `cost` now
        # instead of the sum of squared errors cost
        cost = (-y.dot(np.log(output)) -
        ((1 - y).dot(np.log(1 - output))))
        self.cost_.append(cost)
    return self
def net_input(self, X):
    """Calculate net input"""
    return np.dot(X, self.w_[1:]) + self.w_[0]
def activation(self, z):
    """Compute logistic sigmoid activation"""
    return 1. / (1. + np.exp(-np.clip(z, -250, 250)))
```

```
def predict(self, X):
    """Return class label after unit step"""
    return np.where(self.net_input(X) >= 0.0, 1, 0)
    # equivalent to:
    # return np.where(self.activation(self.net_input(X))
# >= 0.5, 1, 0)
```

```
[50]: X_train_01_subset = X_train[(y_train == 0) | (y_train == 1)]
    y_train_01_subset = y_train[(y_train == 0) | (y_train == 1)]
    lrgd = LogisticRegressionGD(eta=0.05,n_iter=1000,random_state=1)
    lrgd.fit(X_train_01_subset,y_train_01_subset)
    plot_decision_regions(X=X_train_01_subset,y=y_train_01_subset,classifier=lrgd)
    plt.xlabel('petal length [standardized]')
    plt.ylabel('petal width [standardized]')
    plt.legend(loc='upper left')
    plt.show()
```

plt.scatter(x=X[y == cl, 0], y=X[y == cl, 1],alpha=0.8,
c=colors[idx],marker=markers[idx], label=cl,edgecolor='black')

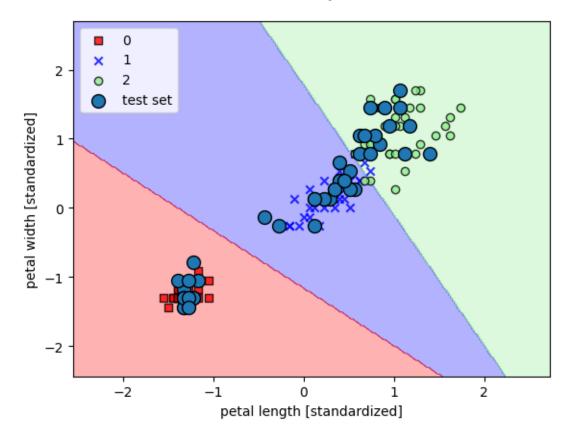


0.3 Training a logistic regression model with scikit-learn

```
[51]: from sklearn.linear_model import LogisticRegression
    lr = LogisticRegression(C=100.0, random_state=1)
    lr.fit(X_train_std, y_train)
    plot_decision_regions(X_combined_std,y_combined,classifier=lr,test_idx=range(105,u=150))
    plt.xlabel('petal length [standardized]')
    plt.ylabel('petal width [standardized]')
    plt.legend(loc='upper left')
    plt.show()
```

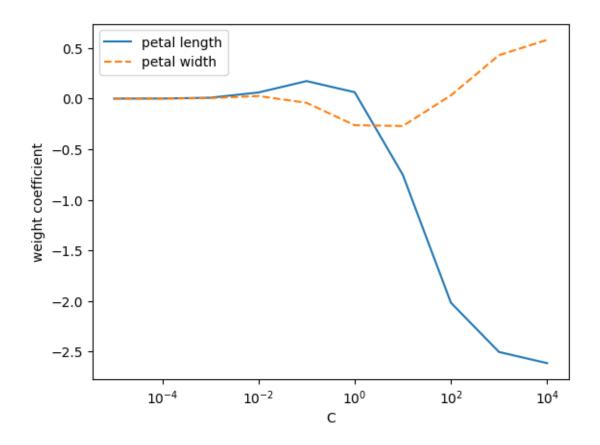
C:\Users\ankit19.gupta\AppData\Local\Temp\ipykernel_15304\1174551119.py:19:
UserWarning: You passed a edgecolor/edgecolors ('black') for an unfilled marker ('x'). Matplotlib is ignoring the edgecolor in favor of the facecolor. This behavior may change in the future.

plt.scatter(x=X[y == cl, 0], y=X[y == cl, 1],alpha=0.8,
c=colors[idx],marker=markers[idx], label=cl,edgecolor='black')



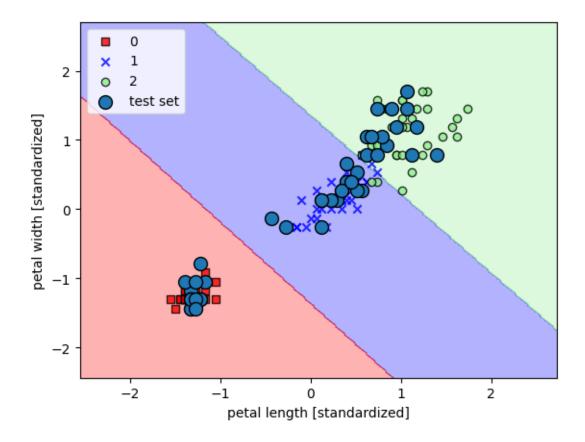
0.4 Tackling Overfitting and Regularization

```
[57]: weights, params = [], []
for c in np.arange(-5, 5):
    lr = LogisticRegression(C=10.**c, random_state=1)
    lr.fit(X_train_std, y_train)
    weights.append(lr.coef_[1])
    params.append(10.**c)
weights = np.array(weights)
plt.plot(params, weights[:, 0],label='petal length')
plt.plot(params, weights[:, 1], linestyle='--',label='petal width')
plt.ylabel('weight coefficient')
plt.xlabel('C')
plt.legend(loc='upper left')
plt.xscale('log')
plt.show()
```



0.5 support vector machines

```
plt.scatter(x=X[y == cl, 0], y=X[y == cl, 1],alpha=0.8,
c=colors[idx],marker=markers[idx], label=cl,edgecolor='black')
```



0.6 Alternative implementations in scikit-learn

```
[59]: # for large datasets

from sklearn.linear_model import SGDClassifier

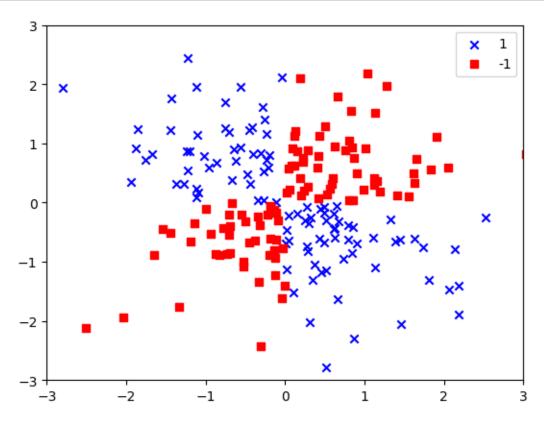
ppn = SGDClassifier(loss='perceptron')

lr = SGDClassifier(loss='log')

svm = SGDClassifier(loss='hinge')
```

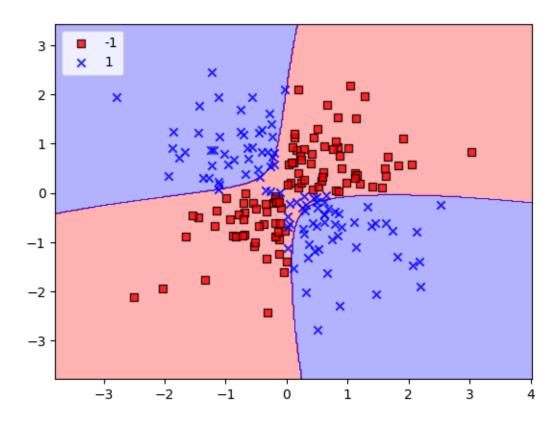
0.7 Solving nonlinear problems using a kernel SVM

0.7.1 Kernel methods for linearly inseparable data

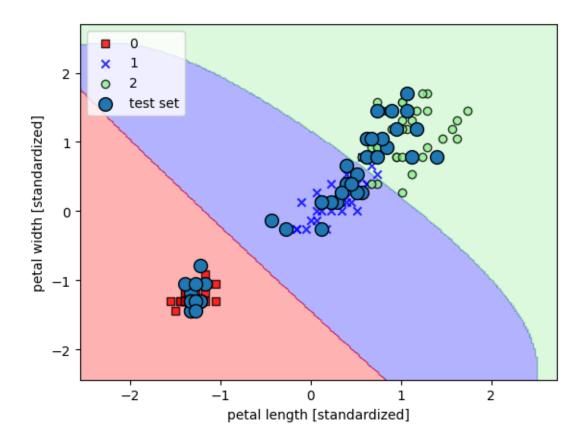


```
[61]: svm = SVC(kernel='rbf', random_state=1, gamma=0.10, C=10.0)
svm.fit(X_xor, y_xor)
plot_decision_regions(X_xor, y_xor, classifier=svm)
plt.legend(loc='upper left')
plt.show()
```

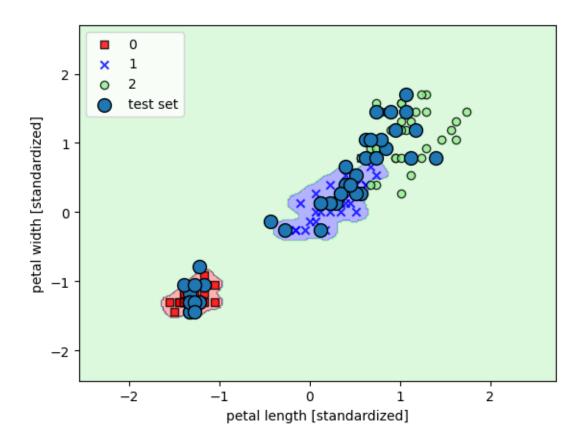
```
plt.scatter(x=X[y == cl, 0], y=X[y == cl, 1],alpha=0.8,
c=colors[idx],marker=markers[idx], label=cl,edgecolor='black')
```



```
plt.scatter(x=X[y == cl, 0], y=X[y == cl, 1],alpha=0.8,
c=colors[idx],marker=markers[idx], label=cl,edgecolor='black')
```

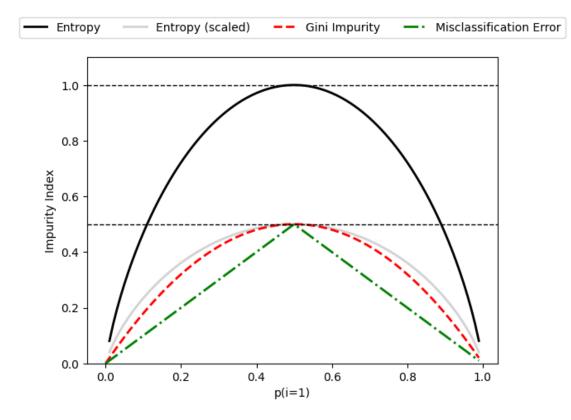


```
plt.scatter(x=X[y == cl, 0], y=X[y == cl, 1],alpha=0.8,
c=colors[idx],marker=markers[idx], label=cl,edgecolor='black')
```



0.8 Decision tree learning

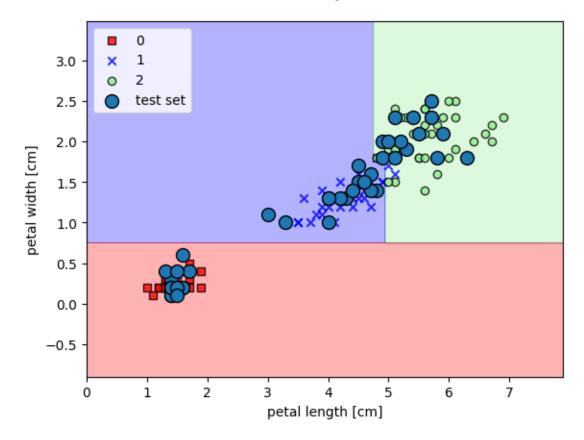
```
[64]: import matplotlib.pyplot as plt
     import numpy as np
     def gini(p):
         return (p)*(1 - (p)) + (1 - p)*(1 - (1-p))
     def entropy(p):
         return - p*np.log2(p) - (1 - p)*np.log2((1 - p))
     def error(p):
         return 1 - np.max([p, 1 - p])
     x = np.arange(0.0, 1.0, 0.01)
     ent = [entropy(p) if p != 0 else None for p in x]
     sc ent = [e*0.5 if e else None for e in ent]
     err = [error(i) for i in x]
     fig = plt.figure()
     ax = plt.subplot(111)
     for i, lab, ls, c, in zip([ent, sc_ent, gini(x), err],['Entropy', 'Entropy_
      ⇒(scaled)','Gini Impurity','Misclassification Error'],['-', '--', '--', '-.
      line = ax.plot(x, i, label=lab,linestyle=ls, lw=2, color=c)
```



0.9 Building a decision tree

```
plt.show()
```

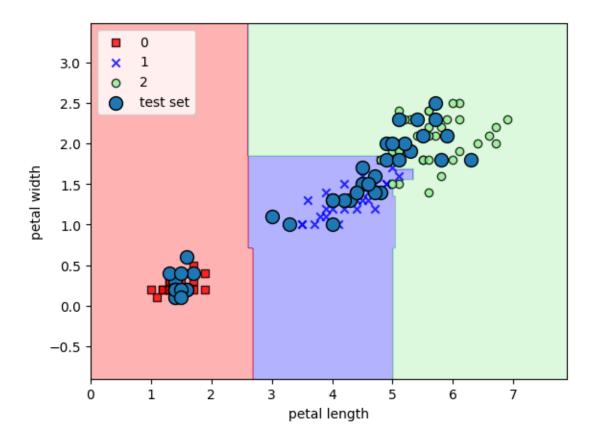
```
plt.scatter(x=X[y == cl, 0], y=X[y == cl, 1],alpha=0.8,
c=colors[idx],marker=markers[idx], label=cl,edgecolor='black')
```



[69]: True

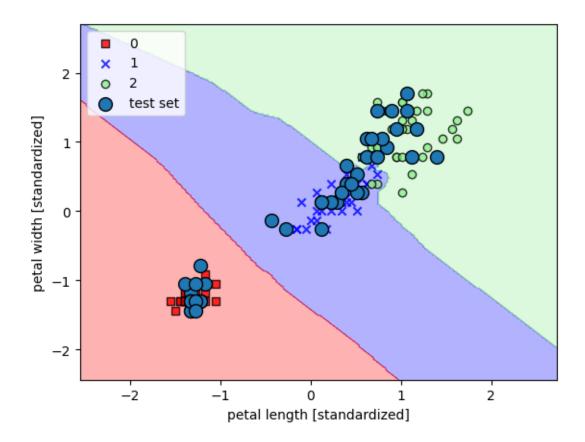
0.10 Combining multiple decision trees via random forests

```
plt.scatter(x=X[y == cl, 0], y=X[y == cl, 1],alpha=0.8,
c=colors[idx],marker=markers[idx], label=cl,edgecolor='black')
```



0.11 KNN - A Lazy Learning Algorithm

```
plt.scatter(x=X[y == cl, 0], y=X[y == cl, 1],alpha=0.8,
c=colors[idx],marker=markers[idx], label=cl,edgecolor='black')
```



[]: