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三维点云处理第三次作业



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K-Means

- fit 函数 E-step

- 循环计算所有数据与每个聚类中心的距离，存储在 `dist_mat`
- 每笔数据的标签：与该笔数据最近的聚类中心的 `id`

```
def fit(self, data):  
    # 作业1  
    # 屏蔽开始  
  
    m = data.shape[0]  
    self.centers_ = data[np.random.choice(m, self.k_, replace = False),:]  
  
    for _ in range(self.max_iter_):  
        dist_mat = np.empty((m, self.k_))  
        for ci, center in enumerate(self.centers_):  
            dist_mat[:,ci] = np.linalg.norm(data - center, axis=1)  
  
        labels = np.argmin(dist_mat, axis=1)
```

K-Means

● fit函数M-step

- distortion: 所有数据点到所属聚类中心距离的平方和
- 更新聚类中心: 属于该聚类的所有点的坐标平均值

```
distortion = (np.min(dist_mat, axis=1)**2).sum()

if distortion < self.tolerance_:
    break

for ci in range(self.k_):
    ci_points = data[labels == ci]
    self.centers_[ci] = np.mean(ci_points, axis=0)

# 屏蔽结束
```

K-Means

● predict函数

- 循环计算所有数据与每个聚类中心的距离，存储在dist_mat
- 每笔数据的标签：与该笔数据最近的聚类中心的id

```
def predict(self, p_datas):  
    result = []  
    # 作业2  
    # 屏蔽开始  
  
    dist_mat = np.empty((p_datas.shape[0], self.k_))  
    for ci, center in enumerate(self.centers_):  
        dist_mat[:,ci] = np.linalg.norm(p_datas - center, axis=1)  
    result = np.argmin(dist_mat, axis=1).tolist()  
  
    # 屏蔽结束  
    return result
```

●初始化

```
def __init__(self, n_clusters, max_iter=50, tol=0.001):  
    self.n_clusters = n_clusters  
    self.max_iter = max_iter  
  
    self.means = None  
    self.covs = None  
    self.weights = np.ones(n_clusters)/n_clusters  
    self.tol = tol
```

● fit函数E-step

- 计算数据点属于k个高斯模型的概率，即
gamma

```
def _gauss(self, j, datum):  
    dim = datum.shape[-1]  
    return 1/pow(2*np.pi, dim/2) * \  
        (1/pow(np.linalg.det(self.covs[j]), 0.5)) * \  
        np.exp(-1/2*np.dot(np.dot((datum-self.means[j]).T,  
                                np.linalg.inv(self.covs[j])),  
                            (datum-self.means[j])))
```

```
def fit(self, data):  
    # 作业3  
    # 屏蔽开始  
  
    N = data.shape[0]  
    dim = data.shape[1]  
    self.means = np.random.random((self.n_clusters, dim))  
    self.covs = np.array(self.n_clusters * [np.identity(dim)])  
    last_nll = float("inf")  
  
    for _ in range(self.max_iter):  
        # E-step  
  
        # posterior probability  
        # (N, k)  
        gamma = np.zeros((N, self.n_clusters))  
  
        for n, datum in enumerate(data):  
            for k in range(self.n_clusters):  
                # gamma[n][k] = self.weights[k]*self._gauss(k, datum)  
                gamma[n][k] = self.weights[k]*multivariate_normal.pdf(datum,  
                                mean=self.means[k], cov=self.covs[k])  
            gamma[n] /= gamma[n].sum()
```

GMM

- fit函数M-step

- 更新k个高斯模型的mean, covariance matrix及weight

- 计算negative log likelihood, 判断是否该跳出循环

```
# M-step
for k in range(self.n_clusters):
    N_k = gamma[:,k].sum()

    self.means[k] = gamma[:,k].dot(data) / N_k

    self.covs[k] = np.zeros((dim, dim))
    for n in range(N):
        diff = np.array([data[n] - self.means[k]])
        self.covs[k] += gamma[n][k] * \
            np.matmul(diff.T, diff)
    self.covs[k] /= N_k

    self.weights[k] = N_k/N
```

```
# negative log likelihood
nll = 0
for n in range(N):
    tmp = 0
    for k in range(self.n_clusters):
        tmp += self.weights[k] * self._gauss(k, data[n])
    nll += np.log(tmp)
nll *= -1

if last_nll - nll < self.tol:
    break
last_nll = nll
```

●predict函数

```
def predict(self, data):  
    # 屏蔽开始  
    N = data.shape[0]  
  
    gamma = np.zeros((N, self.n_clusters))  
  
    for n, datum in enumerate(data):  
        for k in range(self.n_clusters):  
            gamma[n][k] = self.weights[k]*self._gauss(k, datum)  
  
        gamma[n] /= gamma[n].sum()  
  
    return np.argmax(gamma, axis=1)
```


● fit函数E-step

- Gaussian函数：计算“在k个高斯模型里出现这N笔数据的概率”
- 计算gamma

```
def Gaussian(self, data):  
    res = np.empty((self.n_clusters, data.shape[0]))  
    for i in range(self.n_clusters):  
        res[i, :] = multivariate_normal.pdf(data,  
            mean=self.means[i], cov=self.covs[i])  
    return res
```

```
def fit(self, data):  
    # 作业3  
    # 屏蔽开始  
  
    """  
    N: data.shape[0]  
    dim: data.shape[1]  
    k: self.n_clusters  
    """  
  
    dim = data.shape[1]  
    self.means = np.random.random((self.n_clusters, dim))  
    self.covs = np.array(self.n_clusters * [np.identity(dim)])  
    last_nll = float("inf")  
  
    for _ in range(self.max_iter):  
        # E-step  
        # (k, N)  
        gauss_dist = self.Gaussian(data)  
  
        # (k, N)  
        gamma = gauss_dist * self.weights  
        gamma = gamma / np.sum(gamma, axis=0, keepdims=True)
```

GMM-向量化

- fit函数M-step
 - 更新n_clusters个高斯模型的mean, covariance matrix及weight
- 计算negative log likelihood

```
# M-step
# (k, 1)
N_ks = np.sum(gamma, axis=1, keepdims=True)
# (k, 1, dim)
self.means = (gamma.dot(data)/N_ks)
# (k, N, dim, 1)
demeaned_data = (data - self.means[:,np.newaxis,:])[...,np.newaxis]
# (k, N, dim, dim)
demeaned_data2 = np.matmul(demeaned_data,
                             np.transpose(demeaned_data, (0,1,3,2)))
# (k, N, 1, 1) * (k, N, dim, dim) --sum--> (k, dim, dim)
# (k, dim, dim) / (k, 1, dim) -> (k, dim, dim)
self.covs = np.sum(gamma[...,np.newaxis,np.newaxis] * demeaned_data2,
                    axis=1)/N_ks[...np.newaxis]
# (k, 1)
self.weights = N_ks / data.shape[0]
```

```
# negative log likelihood
nll = -np.sum(np.log(np.sum(self.weights * gauss_dist, axis=0)))
if last_nll - nll < self.tol:
    break
last_nll = nll
```

- predict函数
 - 计算gamma
 - 把数据分配到可能性最大的类别里

```
def predict(self, data):  
    # 屏蔽开始  
    gauss_dist = self.Gaussian(data)  
    gamma = gauss_dist * self.weights  
    gamma = gamma / np.sum(gamma, axis=0, keepdims=True)  
  
    return np.argmax(gamma, axis=0)  
    # 屏蔽结束
```

Spectral Clustering

●初始化

- normalized_: Laplacian是否做 normalization
- use_radius_nn_: 使用radius nn 或knn
- nnk_, nnradius_: 用于寻找最近邻
- use_gauss_dist: 使用距离的倒数或高斯函数计算权重
- gauss_sigma: 高斯函数的标准差

```
def __init__(self, n_clusters=2, nnk = 3, nnradius = 1,
             normalized = True, use_radius_nn = False,
             use_gauss_dist = False, gauss_sigma = 5e-1):
    self.n_clusters = n_clusters
    # 屏蔽开始
    # the k for KNN
    self.nnk_ = nnk
    self.nnradius_ = nnradius
    self.labels_ = np.empty(0)
    self.normalized_ = normalized
    self.use_radius_nn_ = use_radius_nn
    self.use_gauss_dist_ = use_gauss_dist
    self.gauss_sigma_ = gauss_sigma
    # 屏蔽结束
```

Spectral Clustering

● fit 函数

- 近邻查找方式: radius nn 或 knn

- 建构 adjacency matrix W

- 将权重设为

- 距离的倒数 或是

- 将距离代入高斯函数得到的值

- 计算 degree matrix D

```
def gauss(self, x):  
    return np.exp(-x*x/(2*self.gauss_sigma*self.gauss_sigma))
```

```
def fit(self, data):  
    # 屏蔽开始  
  
    m = data.shape[0]  
  
    tree = KDTree(data)  
  
    W = np.zeros((m, m))  
  
    for di, datum in enumerate(data):  
        if self.use_radius_nn:  
            nis, ndists = tree.query_radius([datum], self.nnradius_,  
                                           return_distance=True)  
        else:  
            ndists, nis = tree.query([datum], self.nnk+1,  
                                     return_distance=True)  
  
        nis = nis[0]  
        ndists = ndists[0]  
  
        for ni, ndist in zip(nis, ndists):  
            # the point itself will be one of its knn, need to skip it  
            if ni == di: continue  
            if self.use_gauss_dist:  
                W[di][ni] = W[ni][di] = self.gauss(ndist)  
            else:  
                W[di][ni] = W[ni][di] = 1/ndist  
  
    D = np.diag(W.sum(axis=1))
```

Spectral Clustering

● fit函数

- 由D及W计算L

- L_{rw} 有两种计算方式 $L_{rw} = D^{-1}L = I - D^{-1}W$

- 选取最小的k个特征值对应的特征向量建构V矩阵

- 注意numpy.linalg.eig返回的特征值及特征向量是未排序的

- (import numpy.linalg as LA)

```
# unnormalized Laplacian
L = D - W

if self.normalized_:
    L = np.matmul(LA.inv(D), L)
# L = np.identity(m) - np.matmul(LA.inv(D), W)
```

```
eigvals, eigvecs = LA.eig(L)
sorted_idx = np.argsort(eigvals)
V = eigvecs[:, sorted_idx[:self.n_clusters]]
```

Spectral Clustering

- fit 函数

- 使用 KMeans 对 V 矩阵的各 row 进行分类
- 将结果储存于 self.labels_

```
self.labels_ = KMeans(n_clusters=self.n_clusters).fit_predict(V)
```

Spectral Clustering

- predict函数

- 谱聚类算法较难对新数据进行分类
- 参考sklearn的API，同样没有predict函数：<https://scikit-learn.org/stable/modules/generated/sklearn.cluster.SpectralClustering.html>

```
def predict(self, p_datas):  
    pass
```




感谢各位聆听 !
Thanks for Listening

