Applications of Machine Learning in Remote Sensing

Homework 2

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https://github.com/np1140/MLRemoteSensing

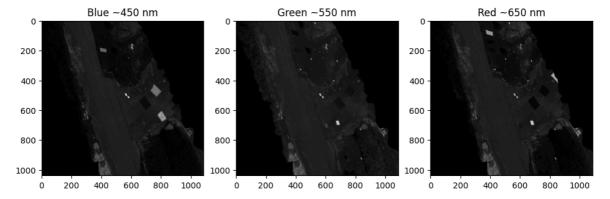
Problem 1

```
In [1]:
        import spectral as spy
        import spectral.io.envi as envi
        import numpy as np
        import re
        import matplotlib.pyplot as plt
        from PIL import Image
        from sklearn.preprocessing import StandardScaler
In [2]: data = spy.envi.open('tait labsphere.hdr', 'tait labsphere') # data loadi
        print(data)
        print(data.shape) # (rows, cols, bands)
              Data Source: './tait labsphere'
              # Rows:
                                 1039
                                1087
              # Samples:
              # Bands:
                                 272
              Interleave:
                                  BS<sub>0</sub>
              Quantization: 32 bits
              Data format: float32
      (1039, 1087, 272)
In [3]: wavelengths = np.array([
            float(re.search(r"\(([\d.]+) nm", name).group(1))
            for name in data.metadata["band names"]
        ])
        # print(wavelengths) # list all the wavelengths
        print(f"Max Wavelength: {wavelengths.max()}, Min Wavelength:{wavelengths.
        # Find band indices
        blue band = np.argmin(abs(wavelengths - 450))
        green_band = np.argmin(abs(wavelengths - 550))
        red_band = np.argmin(abs(wavelengths - 650))
        nir band = np.argmin(abs(wavelengths - 850))
        print(f"Bands indices; Blue:{blue band}, Green: {green band}, Red: {red b
      Max Wavelength: 1001.81, Min Wavelength: 398.573
      Bands indices; Blue:23, Green: 68, Red: 113, Nir: 203
```

Select blue, green, red bands

```
In [4]: # Extract bands
blue = data.read_band(blue_band)
green = data.read_band(green_band)
red = data.read_band(red_band)
nir = data.read_band(nir_band)
```

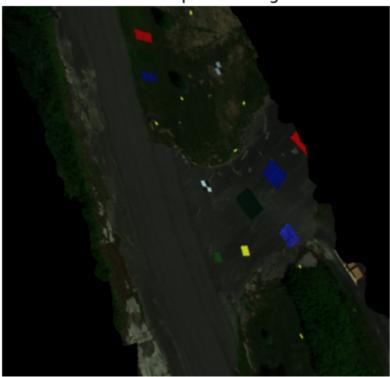
```
In [5]: fig, axs = plt.subplots(1, 3, figsize=(12, 4))
   axs[0].imshow(blue, cmap='gray'); axs[0].set_title("Blue ~450 nm")
   axs[1].imshow(green, cmap='gray'); axs[1].set_title("Green ~550 nm")
   axs[2].imshow(red, cmap='gray'); axs[2].set_title("Red ~650 nm")
   plt.show()
```



```
In [6]: rgb= np.dstack((red, green, blue))
    rgb_norm = rgb / np.max(rgb) # Normalize to [0, 1], max normalization
    plt.imshow(rgb_norm)
    plt.title("RGB Composite Image")
    plt.axis('off')
```

Out[6]: (np.float64(-0.5), np.float64(1086.5), np.float64(1038.5), np.float64(-0.5))





Create pseudocolor BGR (using G, R, NIR)

```
In [7]: pseudo_bgr = np.dstack((
          data.read_band(nir_band), #red - > nir
          data.read_band(red_band), #green - > red
          data.read_band(green_band) # blue - > green
          ))
    pseudo_bgr = pseudo_bgr.astype(float)
    pseudo_bgr /= pseudo_bgr.max()

plt.imshow(pseudo_bgr)
    plt.title("Pseudo-color BGR (Green, Red, NIR)")
    plt.show()
```

Pseudo-color BGR (Green, Red, NIR) 200 400 800 1000 200 400 600 800 1000

In the pseudo-color image, the red channel is assigned to NIR values. The strong red appearance of vegetation indicates that vegetation reflects significantly more NIR than green wavelengths (seen in the plot above). Similarly, blue objects also appear red, leading to the conclusion that they reflect more strongly in the NIR region as well.

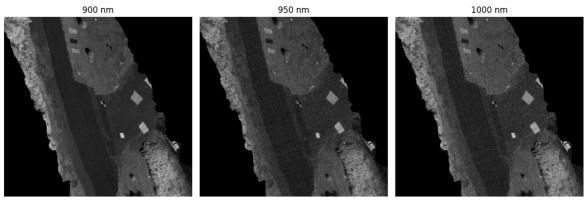
Plot bands at 900, 950, 1000 nm

```
In [8]: wls = [900, 950, 1000]

fig, axs = plt.subplots(1, len(wls), figsize=(12, 5))

for i, wl in enumerate(wls):
    band = np.argmin(abs(wavelengths - wl))
    axs[i].imshow(data.read_band(band), cmap='gray')
    axs[i].set_title(f"{wl} nm")
```

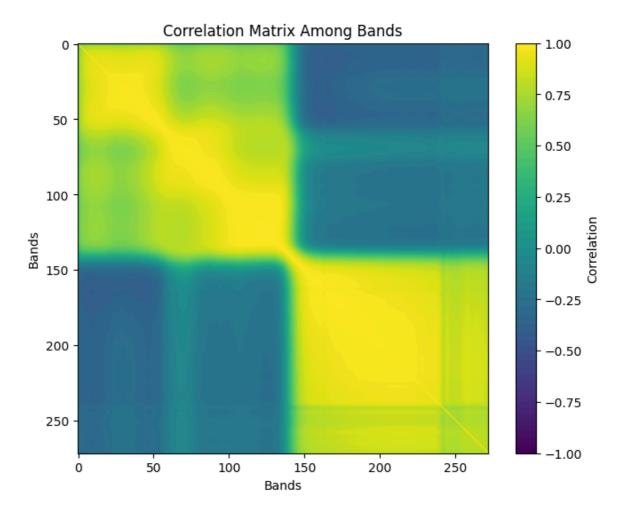
```
axs[i].axis('off')
plt.tight_layout()
```



In these plots, the objects are more distinctly visible compared to the visible range.

Display the correlation matrix as an image

```
In [9]: data = data.load()
         h, w, b = data.shape
         X = data.reshape(-1, b) # shape = (h*w, b)
         mask = np.any(X > 0, axis=1) # Mask no-data pixels
         X mask = X[mask] # keep only valid pixels
         corr_matrix = np.corrcoef(X_mask, rowvar=False) # correlation across ban
         print(corr_matrix.shape) # (272, 272)
        (272, 272)
         plt.figure(figsize=(8, 6))
In [10]:
         plt.imshow(corr matrix, cmap='viridis', vmin=-1, vmax=1)
         plt.colorbar(label="Correlation")
         plt.title("Correlation Matrix Among Bands")
         plt.xlabel("Bands")
         plt.ylabel("Bands")
         plt.show()
```



We observe that bands 0–150 and 150–272 exhibit high intra-group correlation. This is likely because hyperspectral sensors capture data at very fine wavelength intervals, causing adjacent bands to appear nearly identical.

Problem 2: Principal Component Analysis (PCA)

```
In [ ]: def principal_component_analysis(X):
              # Standardize the data
              mean = np.mean(X, axis=0)
              std = np.std(X, axis=0)
              X \text{ std} = (X - \text{mean}) / \text{std}
              # Compute SVD
              U, S, Vt = np.linalg.svd(X_std, full_matrices=False)
              pcs = Vt.T
              m = X std.shape[0]
              eigenvalues = (S**2) / (m - 1)
              return pcs, eigenvalues, X std
In [12]:
         pcs, eigenvalues, X_std = principal_component_analysis(X)
          print(f"PCs shape: {pcs.shape}, Eigenvalues shape: {eigenvalues.shape}")
        PCs shape: (272, 272), Eigenvalues shape: (272,)
 In [ ]: h, w, f = data.shape
          K = 10
```

```
top_pcs = pcs[:, :K]

# Project standardized data to top 10 PCs
X_reduced = np.dot(X_std, top_pcs)

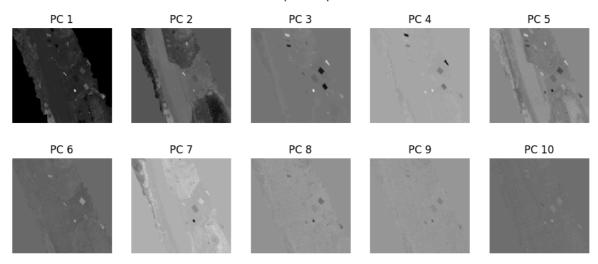
# Reshape
pca_top10_images = X_reduced.reshape(h, w, K)

In [14]: pca_top10_images.shape

Out[14]: (1039, 1087, 10)

In [15]: fig, axs = plt.subplots(2, 5, figsize=(12, 5))
for i in range(K):
    ax = axs[i//5, i%5]
    im = ax.imshow(pca_top10_images[:, :, i], cmap='gray')
    ax.set_title(f"PC {i+1}")
    ax.axis('off')
plt.suptitle("First 10 Principal Components")
plt.show()
```

First 10 Principal Components

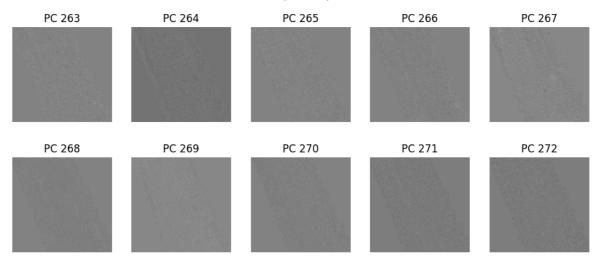


The first principal components (PCs) capture the largest variance, highlighting the main regions such as areas with no data, regions with data, and various objects. The second PC differentiates between vegetation and road. PCs 3, 4, and 5 capture objects of different colors. The last five PCs represent progressively smaller variations, with PCs 8–10 contributing very little additional information.

```
In [16]: last_pcs = pcs[:, -10:]
X_last10 = np.dot(X_std, last_pcs)
pca_last10_images = X_last10.reshape(h, w, 10)

fig, axs = plt.subplots(2, 5, figsize=(12, 5))
for i in range(10):
    ax = axs[i//5, i%5]
    im = ax.imshow(pca_last10_images[:, :, i], cmap='gray')
    ax.set_title(f"PC {f-10+i+1}")
    ax.axis('off')
plt.suptitle("Last 10 Principal Components")
plt.show()
```

Last 10 Principal Components

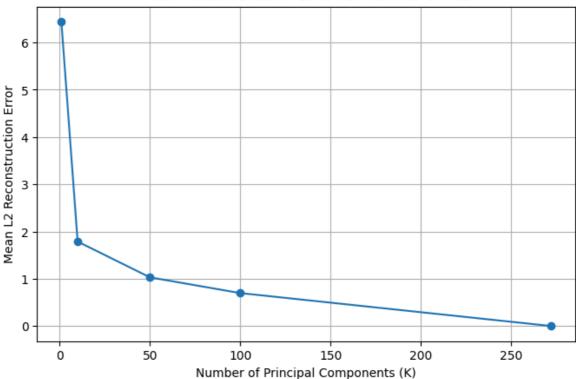


The last PCs do not highlight any meaningful features in the image and can be considered as primarily representing noise.

Error visualization for a selected number of principal components (1, 10, 50, 100)

```
In [17]: def reconstruct pca(X std, pcs, K):
                                                   # K principal components
             top pcs = pcs[:, :K]
             X_reduced = X_std @ top_pcs # Project data to K PCs
             X reconstructed = np.dot(X reduced, top pcs.T) # Reconstruct data
             return X reconstructed
In [18]:
         pc list = [1, 10, 50, 100, 272] # selected number of PCs
         errors = []
         for K in pc list:
             X_rec = reconstruct_pca(X_std, pcs, K)
             error = np.linalg.norm(X_std - X_rec, axis=1) # L2 per pixel
             mean error = error.mean()
             errors.append(mean error)
             print(f"PCs={K}, Mean reconstruction L2 error: {mean_error:.4f}")
        PCs=1, Mean reconstruction L2 error: 6.4400
        PCs=10, Mean reconstruction L2 error: 1.7872
        PCs=50, Mean reconstruction L2 error: 1.0302
        PCs=100, Mean reconstruction L2 error: 0.6965
        PCs=272, Mean reconstruction L2 error: 0.0000
In [19]: plt.figure(figsize=(8, 5))
         plt.plot(pc list, errors, marker='o')
         plt.title("Mean Reconstruction Error vs. Number of PCs")
         plt.xlabel("Number of Principal Components (K)")
         plt.ylabel("Mean L2 Reconstruction Error")
         plt.grid(True)
         plt.show()
```





The reconstruction error is high when only one principal component is used. As more PCs are included, the reconstruction error gradually decreases. When all 272 PCs are used, the reconstruction error becomes zero, as all the information is incorporated; however, using fewer PCs omits certain details.

```
In [20]: explained_var_ratio = eigenvalues / eigenvalues.sum()
    cumulative_var = np.cumsum(explained_var_ratio)
    K = np.argmax(cumulative_var >= 0.99) + 1
    print(f"Number of PCs explaining ≥99% variance: {K}")
```

Number of PCs explaining ≥99% variance: 36

```
In [21]: X_pc = np.dot(X_std, pcs)
X_mean = X_mask.mean(axis=0)
X_pc_cpy1 = X_pc.copy()
# K= 1
# Zero out contributions after K
X_pc_cpy1[:, K:] = 0

# Backward transform
X_reconstructed = np.dot(X_pc_cpy1, pcs.T) + X_mean
X_reconstructed = X_reconstructed.reshape(h, w, f)
```

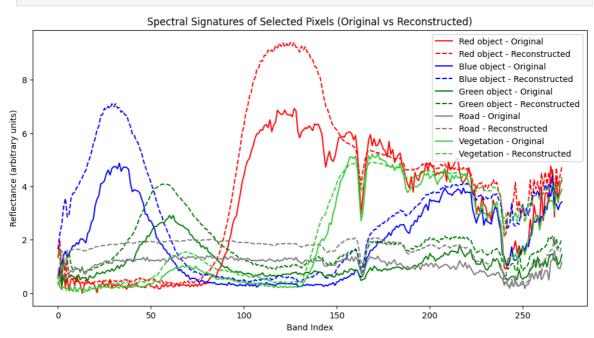
```
In [22]: X_reconstructed.shape
```

Out[22]: (1039, 1087, 272)

```
In [23]: import plotly.express as px
fig = px.imshow(
    rgb_norm,
    title="Normalized RGB Image"
)
fig.update_layout(
```

```
width=800,
height=500,
dragmode=False
)
fig.show()
```

```
five_interesting_pixels = {
In [24]:
             "Red object": (370, 825),
             "Blue object": (467, 767),
             "Green object": (704, 608),
             "Road": (594, 423),
             "Vegetation": (909, 949),
         }
         # Assign fixed colors
         pixel colors = {
             "Red object": "red",
             "Blue object": "blue",
             "Green object": "green",
             "Road": "gray",
             "Vegetation": "limegreen",
         }
         plt.figure(figsize=(12, 6))
         for label, (r, c) in five interesting pixels.items():
             orig_spectrum = np.array(data[r, c, :])
                                                           # original
             recon spectrum = X reconstructed[r, c, :] # reconstructed
             color = pixel colors[label]
             plt.plot(orig_spectrum[0][0], color=color, label=f"{label} - Original
             plt.plot(recon spectrum, "--", color=color, label=f"{label} - Reconst
         plt.title("Spectral Signatures of Selected Pixels (Original vs Reconstruc
         plt.xlabel("Band Index")
         plt.ylabel("Reflectance (arbitrary units)")
         plt.legend()
         plt.show()
```



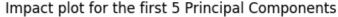
When utilizing all the PC that represent 99% variability in total, the spectral signatures of five different pixels representing various objects closely resembles to that of original data spectral signature.

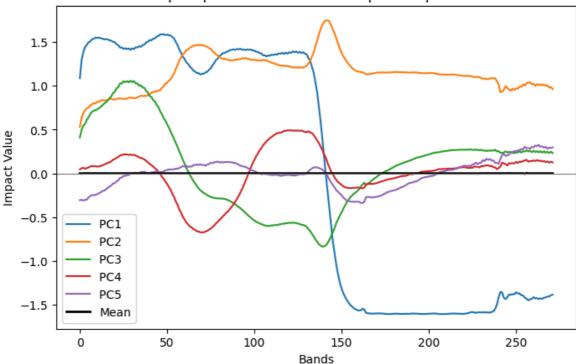
SNR of Original and Reconstructed Signals

```
In [25]: data = data.load()
         data array = np.array(data)
         data array[data array == 0] = np.nan
         mu = np.nanmean(data array) # mean per band
         sigma = np.nanstd(data array) # std per band
         SNR org = mu / sigma
         print("Mean SNR (original):", SNR org)
       Mean SNR (original): 1.5260783
In [26]: X reconstructed = X reconstructed.astype(np.float32)
         valid reconstructed = X reconstructed
         valid reconstructed[valid reconstructed == 0] = np.nan
         mu rec = np.nanmean(valid reconstructed)
         sigma rec = np.nanstd(valid reconstructed)
         snr per band rec = mu rec / sigma rec
         SNR_rec = snr_per_band_rec.mean()
         print("Mean SNR (reconstructed):", SNR rec)
         improvement = SNR rec / SNR org
         print(f"SNR Improvement factor: {improvement:.2f}x")
       Mean SNR (reconstructed): 1.1749474
       SNR Improvement factor: 0.77x
         Impact plot
In [27]: data = spy.envi.open('tait labsphere.hdr', 'tait labsphere') # data loadi
         data = data.load()
         h, w, f = data.shape
         print(data.shape)
        (1039, 1087, 272)
 In []: X = data.reshape(-1, f)
         mask = np.any(X != 0, axis=1)
         X_{valid} = X[mask]
         scaler = StandardScaler()
         X std = scaler.fit transform(X valid) # (num valid pixels, 272)
         X std.shape
 Out[]: (729978, 272)
```

In []: pcs, eigenvalues, X std = principal component analysis(X valid)

```
In []: num pcs = 5
        impacts = []
        for i in range(num pcs):
            pc = pcs[:, i]
                                           # PC
                                           # Eigenvalue
            ev = eigenvalues[i]
            impact = 2 * np.sqrt(ev) * pc # Scale by 2*sqrt(eigenvalue), reflec
            impacts.append(impact)
        impacts = np.array(impacts)
        print("Impacts shape:", impacts.shape) # (5, 272)
       Impacts shape: (5, 272)
In [ ]: import numpy as np
        import matplotlib.pyplot as plt
        mean features = X std.mean(axis=0)
        features = np.arange(X_std.shape[1])
        # Plot mean and impacts for the first 5 PCs
        plt.figure(figsize=(8, 5))
        for i in range(5):
            impact = impacts[i]
            plt.plot(features, impact, label=f'PC{i+1} ')
        plt.plot(features, mean features, color='black', linewidth=2, label='Mean
        plt.xlabel("Bands")
        plt.ylabel("Impact Value")
        plt.title("Impact plot for the first 5 Principal Components")
        plt.axhline(0, color='gray', linewidth=0.8)
        plt.legend()
        plt.show()
```





The impact plot for the first five principal components illustrates how each component influences the spectral bands. PC1 shows a strong positive impact on the first ~150 bands and a negative impact on the remaining bands, indicating a contrast between the lower and higher wavelength regions. PC2 is nearly flat across all bands, suggesting it captures minor uniform variation rather than specific spectral differences. PC3 primarily affects bands 0–165, highlighting variability in the lower-to-mid wavelength range, while PC4 exhibits significant impact on bands 50–150, reflecting localized spectral variability within this region. PC5 shows minimal impact across all bands, contributing little to the overall spectral variance. Overall, these observations indicate that the first few principal components capture the most meaningful spectral variations, with PC1 dominating global contrast and subsequent PCs highlighting finer, localized differences.

Problem 3: K-Means Clustering

```
In []: img = Image.open("jellybeans.tiff") # Load TIFF image
    plt.imshow(img)
    plt.axis("off")
    plt.show()
```



In []: def kmeans(X, K, max iters=100, tol=1e-4, seed=42):

```
np.random.seed(seed)
            # Initialize cluster centers randomly
            rand indices = np.random.choice(X.shape[0], K, replace=False)
            centroids = X[rand indices]
            for i in range(max iters):
                # Assign points to nearest centroid
                distances = np.linalg.norm(X[:, None] - centroids[None, :], axis=
                labels = np.argmin(distances, axis=1)
                # Compute new centroids
                new_centroids = np.array([X[labels == k].mean(axis=0) for k in ra
                # Check convergence criteria
                if np.linalg.norm(new centroids - centroids) < tol:</pre>
                    print(f"Converged at iteration {i}")
                    break
                centroids = new centroids
            return labels, centroids
In [ ]: data = np.array(img) # shape: (H, W, 3)
        # Reshape
        X = data.reshape(-1, 3).astype(np.float64)
        # Standardize
        X_{std} = (X - X.mean(axis=0)) / X.std(axis=0)
In []: K = 6 # adjusted based on colors in an image
        labels, centroids = kmeans(X_std, K)
        # Map labels back to image space using original color scale
```

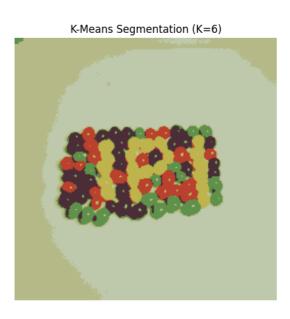
```
segmented = centroids[labels].reshape(data.shape)
# Undo standardization
segmented = (segmented * X.std(axis=0) + X.mean(axis=0)).astype(np.uint8)

# Display original and clustered image
plt.figure(figsize=(12, 6))
plt.subplot(1, 2, 1)
plt.title("Original Image")
plt.imshow(img)
plt.axis("off")

plt.subplot(1, 2, 2)
plt.title(f"K-Means Segmentation (K={K})")
plt.imshow(segmented)
plt.axis("off")
```

Converged at iteration 54





Experimentally it is determined that K =6 works well for jellybeans image.

Problem 3(b)

```
In []: sentinal_data = np.load("sentinel2_rochester.npy")
    rows, cols, bands = sentinal_data.shape # (rows, cols, bands)
    X = sentinal_data.reshape(-1, bands)
    mask = np.any(X != 0, axis=1) # Avoid no data pixels
    X_valid = X[mask]
    print("Original pixels:", X.shape[0])
    print("Valid pixels:", X_valid.shape[0])

Original pixels: 683064
    Valid pixels: 630024

In []: pcs, eigenvalues, X_std = principal_component_analysis(X_valid)

In []: pc_representations = {}
    for k in [3, 4, 5, 6]:
        # Selecting the first k principal components
        top_pcs = pcs[:, :k] # (bands, k)
```

```
# Project the data
            X reduced = np.dot(X std, top pcs) # (N samples, k)
            # Store for later use
            pc representations[k] = X reduced
            print(f"Shape with {k} PCs:", X reduced.shape)
       Shape with 3 PCs: (630024, 3)
       Shape with 4 PCs: (630024, 4)
       Shape with 5 PCs: (630024, 5)
       Shape with 6 PCs: (630024, 6)
In [ ]: cluster_results = {}
        for k in [3, 4, 5, 6]:
            X_reduced = pc_representations[k] # (rows*cols, k)
            X reduced standard = (X reduced - X reduced.mean(axis=0)) / X reduced
            # number of clusters
            K = 5
            labels, centroids = kmeans(X reduced standard, K)
            full labels = np.full(X.shape[0], -1, dtype=int)
            full labels[mask] = labels
            cluster results[k] = full labels.reshape(rows, cols)
            print(f"K-Means done for {k} PCs -> output shape {cluster results[k].
       Converged at iteration 36
       K-Means done for 3 PCs -> output shape (954, 716)
       Converged at iteration 49
       K-Means done for 4 PCs -> output shape (954, 716)
       K-Means done for 5 PCs -> output shape (954, 716)
       Converged at iteration 57
       K-Means done for 6 PCs -> output shape (954, 716)
In [ ]: fig, axes = plt.subplots(1, 4, figsize=(12, 4))
        for idx, k in enumerate([3, 4, 5, 6]):
            axes[idx].imshow(cluster results[k])
            axes[idx].set title(f"K-Means on {k} PCs")
            axes[idx].axis("off")
        plt.show()
         K-Means on 3 PCs
                             K-Means on 4 PCs
                                                 K-Means on 5 PCs
                                                                    K-Means on 6 PCs
```

Here, we found that instead of manually selecting relevant features, the earlier PCs that capture high variability can be used to transform the data into a lower-dimensional

representation, enabling effective clustering.

problem 3(c)

```
In []: data = spy.envi.open('tait_labsphere.hdr', 'tait_labsphere') # data loadi
    data = data.load()
    data_patch = data[650:900, 650:900, :]

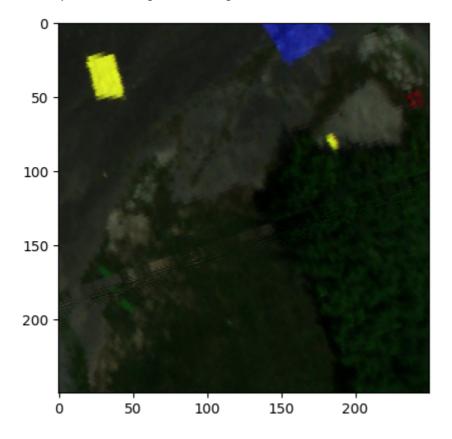
    rgb_bands = [red_band, green_band, blue_band]

# Extract RGB bands
    rgb_data_patch = data_patch[:, :, rgb_bands]

    rgb_data_patch_norm = (rgb_data_patch - np.min(rgb_data_patch)) / (np.max
    plt.imshow(rgb_data_patch_norm)

/tmp/ipykernel_105721/1400234148.py:11: DeprecationWarning:
    _array_wrap__ must accept context and return_scalar arguments (positional
    ly) in the future. (Deprecated NumPy 2.0)
```

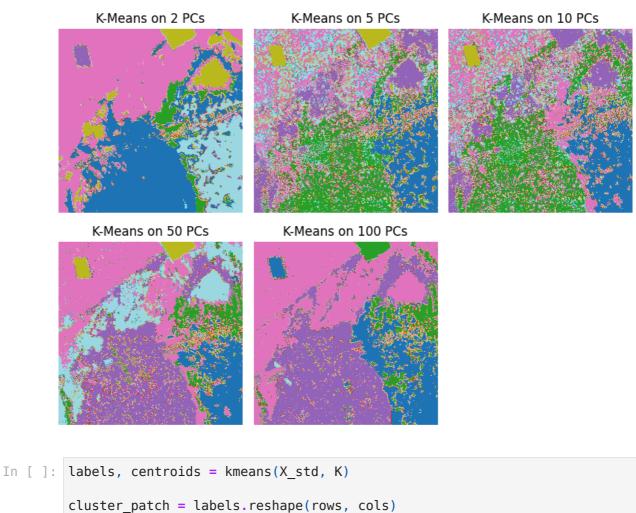
Out[]: <matplotlib.image.AxesImage at 0x76880c02b640>



```
In [ ]: rows, cols, bands = data_patch.shape
    X_data_patch = data_patch.reshape(-1, bands)
    pcs, eigenvalues, X_std = principal_component_analysis(X_data_patch)

In [ ]: pc_representations = {}
    for k in [2, 5, 10, 50, 100]:
        top_pcs = pcs[:, :k]
```

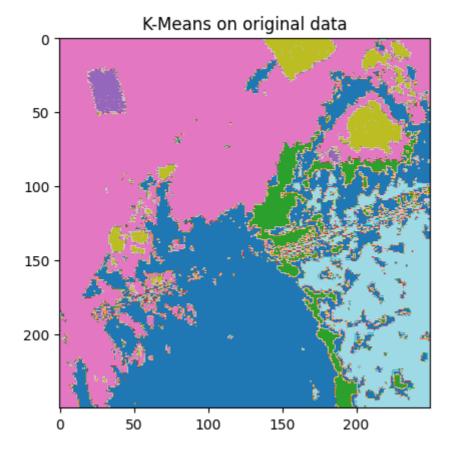
```
X \text{ reduced} = \text{np.dot}(X \text{ std, top pcs}) \# (N \text{ valid, } k)
            pc representations[k] = X reduced
In [ ]: cluster results = {}
        K = 6 # number of clusters
        for k in [2, 5, 10, 50, 100]:
            X reduced = pc representations[k]
            X_reduced_standard = (X_reduced - X_reduced.mean(axis=0)) / X reduced
            labels, centroids = kmeans(X reduced standard, K)
            cluster_results[k] = labels.reshape(rows, cols)
            print(f"K-Means done for {k} PCs -> output shape {cluster results[k].
       Converged at iteration 16
       K-Means done for 2 PCs -> output shape (250, 250)
       Converged at iteration 41
       K-Means done for 5 PCs -> output shape (250, 250)
       Converged at iteration 49
       K-Means done for 10 PCs -> output shape (250, 250)
       Converged at iteration 63
       K-Means done for 50 PCs -> output shape (250, 250)
       Converged at iteration 35
       K-Means done for 100 PCs -> output shape (250, 250)
In [ ]: fig, axes = plt.subplots(2, 3, figsize=(8, 6))
        # Flatten axes
        axes = axes.ravel()
        for idx, k in enumerate([2, 5, 10, 50, 100]):
            axes[idx].imshow(cluster_results[k], cmap="tab20")
            axes[idx].set title(f"K-Means on {k} PCs")
            axes[idx].axis("off")
        axes[-1].axis("off")
        plt.tight_layout()
        plt.show()
```



```
print(f"K-Means done for {k} PCs -> output shape {cluster_patch.shape}")

Converged at iteration 20
K-Means done for 100 PCs -> output shape (250, 250)

In []: plt.imshow(cluster_patch, cmap='tab20')
plt.title("K-Means on original data")
plt.show()
```



We observe that good clusters can be formed using fewer principal components rather than relying on the full-dimensional data. This reduces the computational burden associated with calculating Euclidean distances in the k-means algorithm. We also observe that as more PCs are included, increasingly subtle features are distinguished, which may act as noise and provide limited benefit.

In []: