

### Prob 1

let's assume we have an eigenvector  $\underline{v}$  of

$$\text{Matrix } \underline{S} = \frac{1}{N} \underline{X} \underline{X}^T$$

$\therefore$  It will satisfy  $\underline{S} \underline{v} = \lambda \underline{v}$  eigenvalue.

$$\Rightarrow \frac{1}{N} (\underline{X} \underline{X}^T) \underline{v} = \lambda \underline{v}$$

$$\Rightarrow \frac{1}{N} (\underline{X}^T \underline{X}) (\underline{X}^T \underline{v}) = \lambda (\underline{X}^T \underline{v}) \quad \left[ \begin{array}{l} \text{pre multiply } \underline{X}^T \\ \text{both side} \end{array} \right]$$

$$\text{Substitute } \underline{u} = \underline{X}^T \underline{v}$$

$$\Rightarrow \frac{1}{N} (\underline{X}^T \underline{X}) \underline{u} = \lambda \underline{u}$$

$\therefore \underline{u} = \underline{X}^T \underline{v}$  is nothing else but an eigenvector for  $\frac{1}{N} (\underline{X}^T \underline{X})$

In normal case to compute  $K$  eigenvectors for  $\frac{1}{N} (\underline{X}^T \underline{X})$  complexity will be  $O(KD^2)$  but in this case order will be  $O(KN^2) + O(KND)$

$\therefore$  overall complexity  $O(KND)$  which is less than  $O(KD^2)$

$\downarrow$  for decomp. of  $\frac{1}{N} (\underline{X} \underline{X}^T)$        $\downarrow$  for multi matrix

$$\text{as } \boxed{N < D}$$

Prob 2

$$h(x) = x \sigma(\beta x)$$

# if we choose  $\beta = 0$

$$\Rightarrow \sigma(0) = 1$$

$$\Rightarrow h(x) = \frac{x}{2}$$

~~a linear~~  
activation  
function.

# if we choose  $\beta \rightarrow \infty$

$$\Rightarrow \text{for } x < 0 \Rightarrow \sigma(\beta x) = \sigma(-\infty) = 0$$

$$\text{or for } x > 0 \Rightarrow \sigma(\beta x) = \sigma(\infty) = 1$$

$$\Rightarrow h(x) = \begin{cases} x & ; x > 0 \\ 0 & ; x < 0 \end{cases}$$

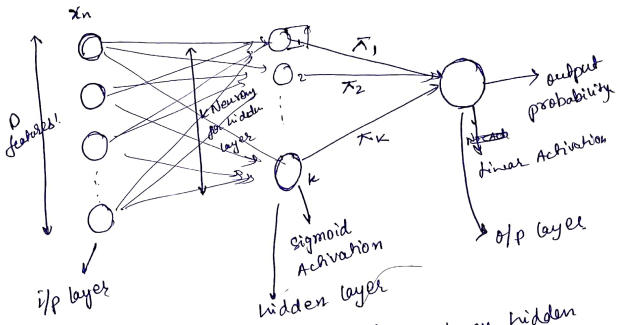
} ReLU  
↓  
Activation  
f.n

Prob 3

$$P(y_n=1 | x_n) = \prod_{k=1}^K P(y_n=1 | z_n=k, x_n) P(z_n=k | x_n)$$

∴ Above equation can be modelled as a multiplication of two probabilistic fns.

⇒ 2 layer NN with one hidden layer.



∴ basically  $\sigma(w_n^T x_n) \rightarrow$  acts as an hidden neurons

Activation is sigmoid -  
∴ Weights for first layer will be  $(w_{z1}, w_{z2}, \dots, w_{zk})$  for  $k$  neurons.

∴ Weight for outputs will be  $[\pi_1, \pi_2, \dots, \pi_k]^T$   
multiplying the probability of  $P(z_n | x_n)$

Prob 4

$$P(u_n, v_m, \theta_n, \phi_m) \propto X$$

$$= \prod_{n=1, m=1}^{N, N} P(x_{nm} | u_n, v_m, \theta_n, \phi_m) P(u_n) P(v_m)$$

Q

$$= \prod_{n,m} \sqrt{\frac{\lambda_v}{2\pi}} \exp \left\{ -\frac{\lambda_v}{2} (x_{nm} - \theta_n - \phi_m - u_n^T v_m)^2 \right\} \\ \times \sqrt{\frac{\lambda_u}{2\pi}} \left\{ \exp \left( -\frac{\lambda_u}{2} \|u_n - w_u a_n\|^2 \right) \right\} \\ \times \sqrt{\frac{\lambda_v}{2\pi}} \exp \left( -\frac{\lambda_v}{2} \|v_m - w_v a_m\|^2 \right)$$

$$\Rightarrow NLL = \sum_{n,m} \left\{ \frac{\lambda_v}{2} (x_{nm} - \theta_n - \phi_m - u_n^T v_m)^2 \right\} \\ + \sum_n \left\{ \frac{\lambda_u}{2} \|u_n - w_u a_n\|^2 \right\} \\ + \sum_m \left\{ \frac{\lambda_v}{2} \|v_m - w_v a_m\|^2 \right\}$$

NLL  $\rightarrow$  cost function

# Now, for  $N_u, W_u \rightarrow$  minimization will be similar to linear regression model.

which will give :-

$$W_u^T = (A^T A)^{-1} A^T U$$

where;  $A \rightarrow N \times K$  matrix of all  $a_n$ 's

$U \rightarrow N \times K$  matrix of all  $u_n$ 's

$$\Rightarrow W_u = U^T A (A^T A)^{-1}$$

Similarly,  $W_b = V^T B (B^T B)^{-1}$

# for  $\theta_n$  &  $\phi_m$  taking the derivatives yield.

$$\sum_m -\lambda_n (x_{nm} - \theta_n - \phi_m - u_n^T v_m) = 0$$

$$\Rightarrow \theta_n = \frac{1}{n} \sum_m (x_{nm} - \phi_m - u_n^T v_m)$$

$$\text{Similarly } \phi_m = \frac{1}{N} \sum_n (x_{nm} - \theta_n - u_n^T v_m)$$

# Now taking the derivative w.r.t.  $\mathbf{v}_m$  we get.

$$\sum_n -\lambda_n (x_{nm} - \theta_n - \phi_m - \mathbf{u}_n^T \mathbf{v}_m) \mathbf{u}_n + \lambda_\nu (\mathbf{v}_m - \mathbf{W}_\nu \mathbf{b}_m) = 0$$

$$\Rightarrow \left( \sum_n \lambda_n \mathbf{u}_n \mathbf{u}_n^T + \lambda_\nu \mathbf{I}_K \right) \mathbf{v}_m = \lambda_\nu \mathbf{W}_\nu \mathbf{b}_m + \sum_n \lambda_n (x_{nm} - \theta_n - \phi_m) \mathbf{u}_n$$

We can write  $\sum_n \mathbf{u}_n \mathbf{u}_n^T = \mathbf{U}^T \mathbf{U}$

$$\Rightarrow \mathbf{v}_m = \left( \lambda_n \mathbf{U}^T \mathbf{U} + \lambda_\nu \mathbf{I}_K \right)^{-1} \left[ \sum_n \lambda_n (x_{nm} - \theta_n - \phi_m) \mathbf{u}_n + \lambda_\nu \mathbf{W}_\nu \mathbf{b}_m \right]$$

Similarly, for  $\mathbf{u}_m$

$$\mathbf{u}_m = \left( \lambda_n \mathbf{V}^T \mathbf{V} + \lambda_u \mathbf{I}_K \right)^{-1} \left[ \sum_m \lambda_n (x_{nm} - \theta_n - \phi_m) \mathbf{v}_m + \lambda_u \mathbf{W}_u \mathbf{a}_n \right]$$

$\therefore$  Using the equation shown above. ALI-OPT algo will be :-

• Initial

PTO

### ALT - OPT

- latent variables  $U_n = \hat{U}_n; V_m = \hat{V}_m$
- ① Initialize
  - ② Solve for  $\hat{W}_u, \hat{W}_v, \hat{P}_n, \hat{Q}_m$  using the eq<sup>n</sup>.
  - ③ Solve  $\hat{U}_n, \hat{V}_m$  using the eq<sup>n</sup>
  - ④ Go to step ② if not converge.

**Introduction to ML (CS771), Autumn 2018**  
**Indian Institute of Technology Kanpur**  
**Homework Assignment Number 5**

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*Roll Number:* 160645

*Date:* November 17, 2018

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**QUESTION**

**5**

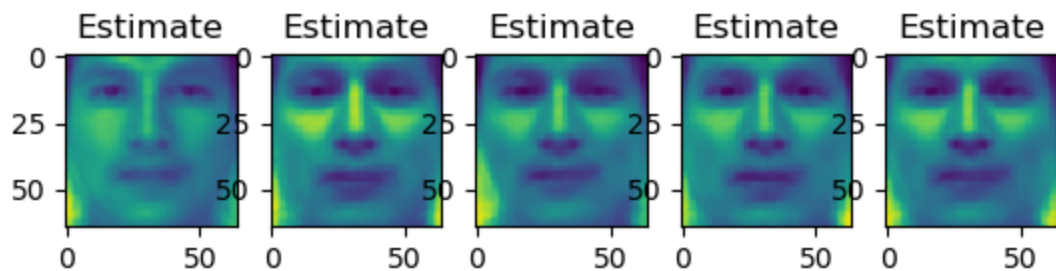
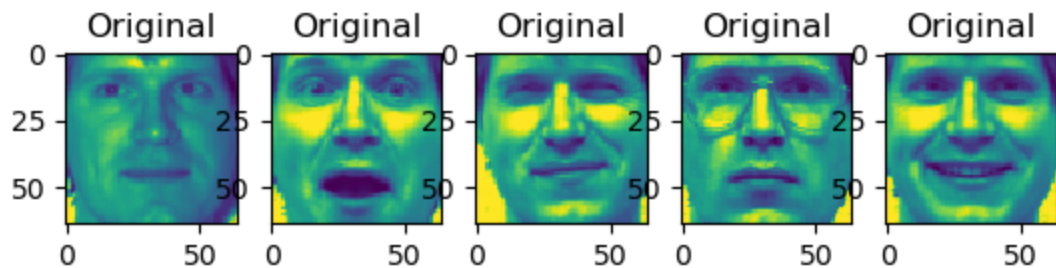
**Programming Part 1:**

On increasing  $k$  there is improvement in the reconstruction of images! Moreover the images appear to be more clearer. Explanation is simple higher the value of  $k$  more the no of features retained in  $z$ .

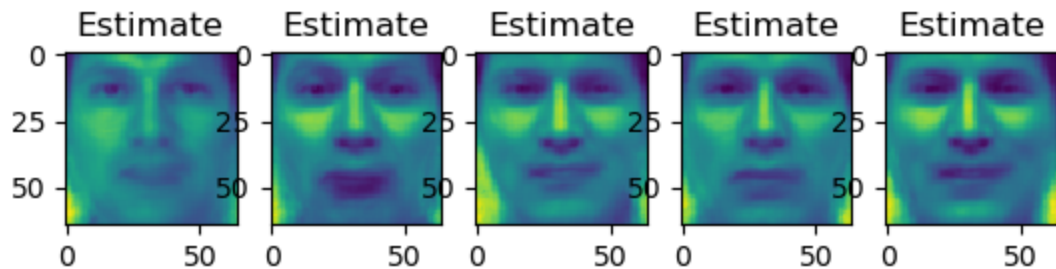
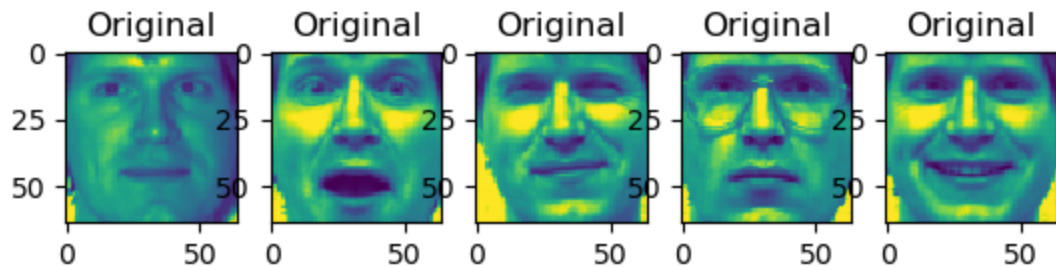
Please find the plots from next page.



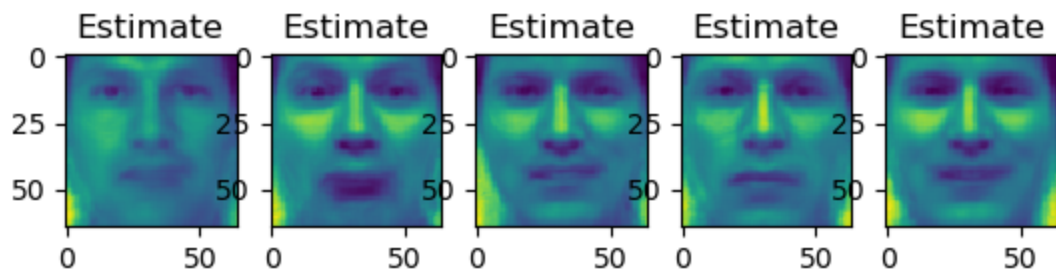
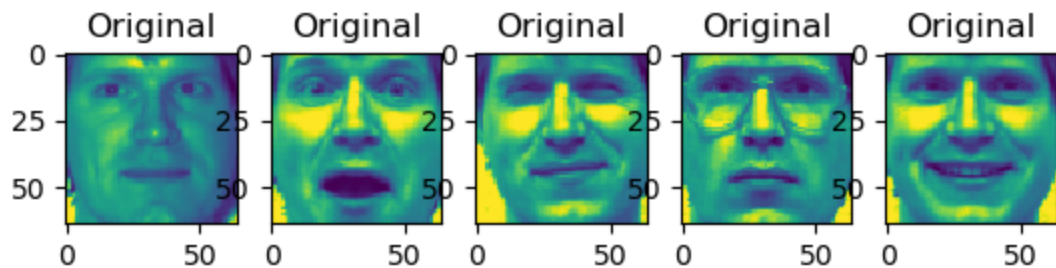
For  $K = 10$



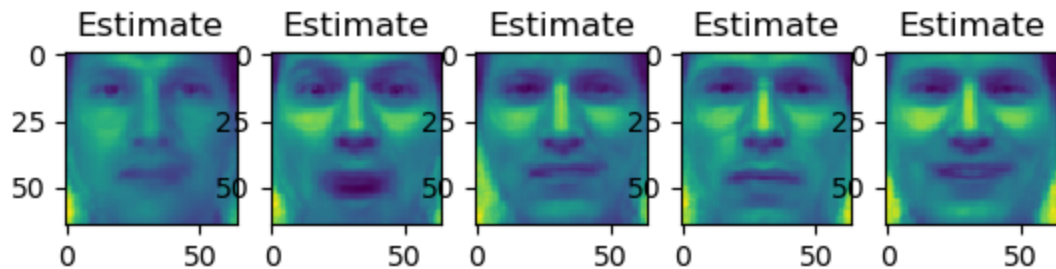
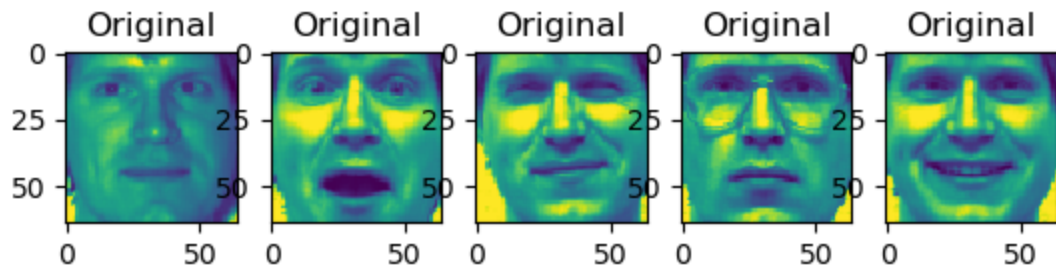
For  $K = 20$



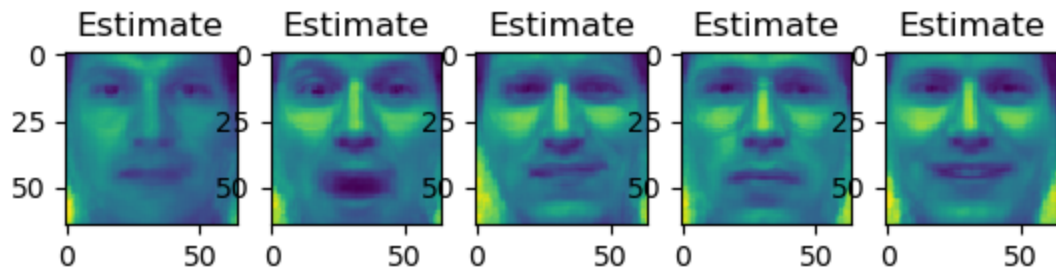
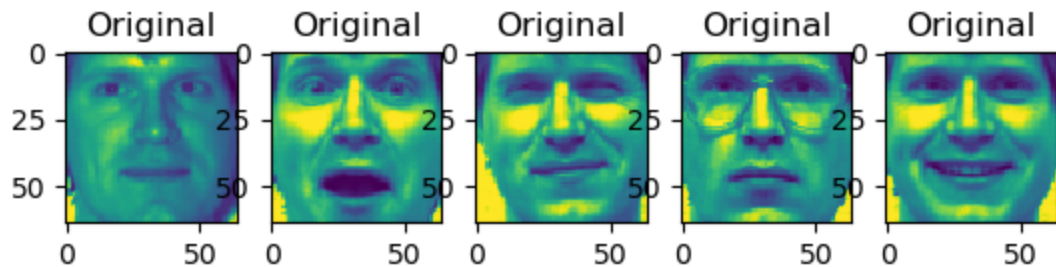
For  $K = 30$



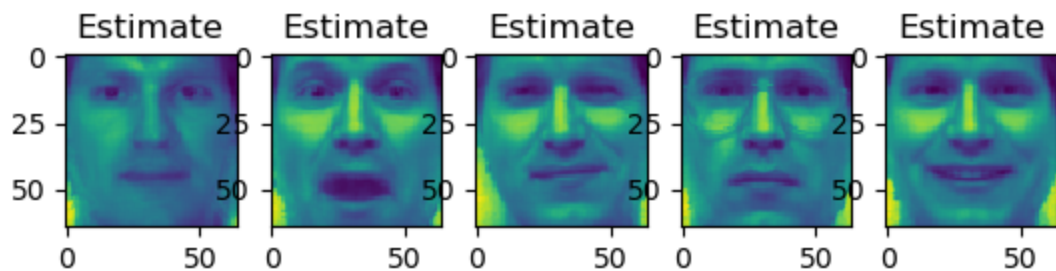
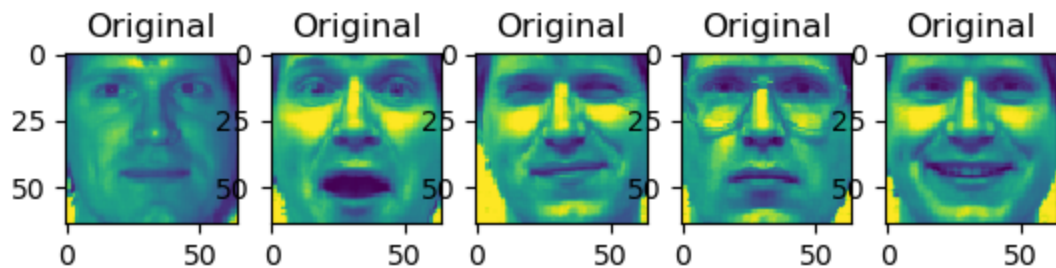
For  $K = 40$



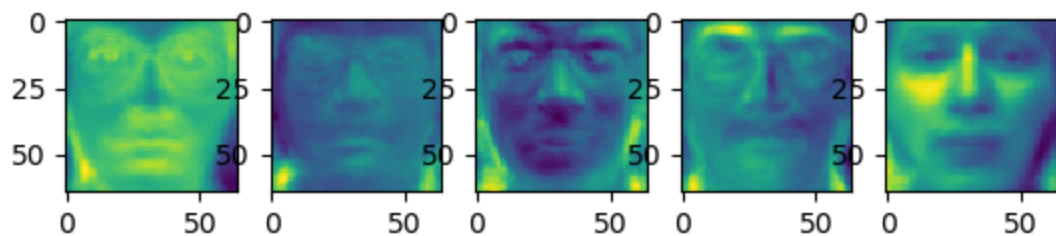
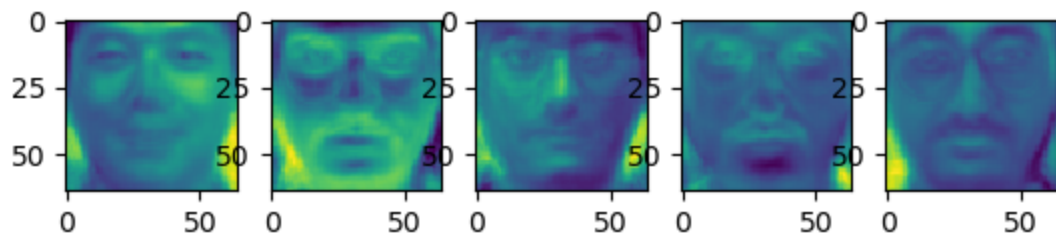
For  $K = 50$



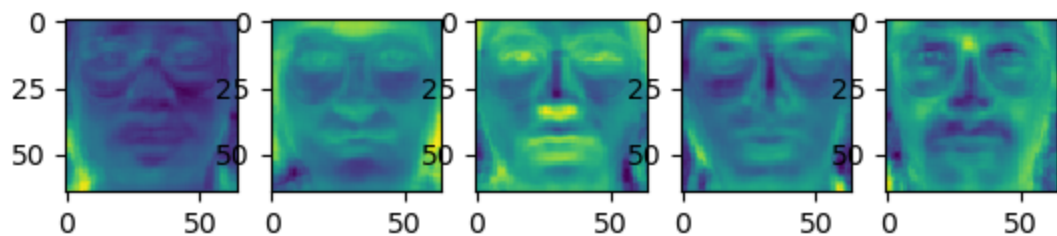
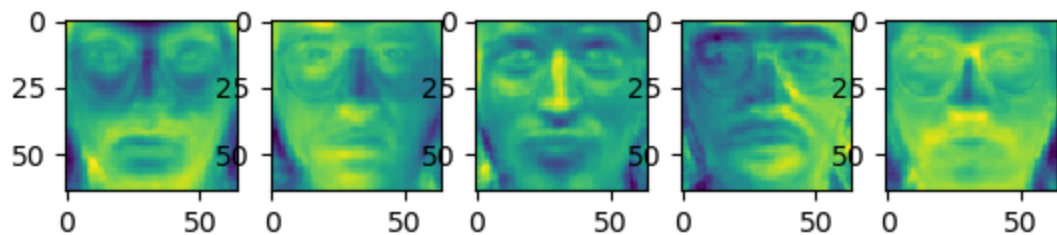
For  $K = 100$



For  $K = 10$

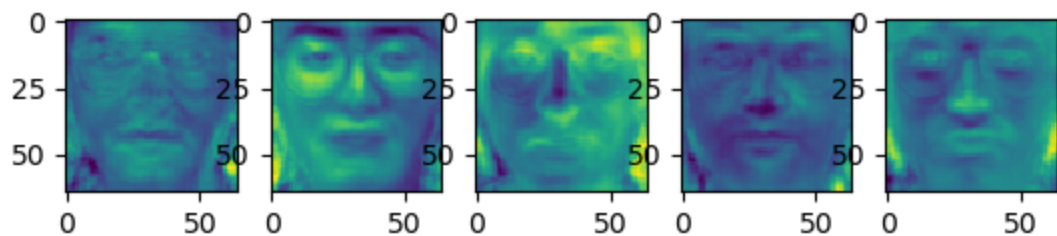
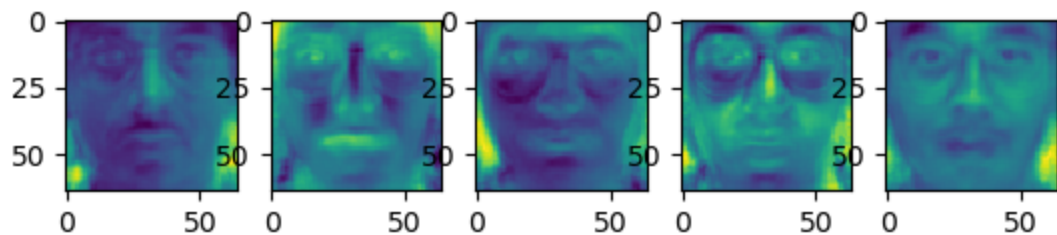


For  $K = 20$

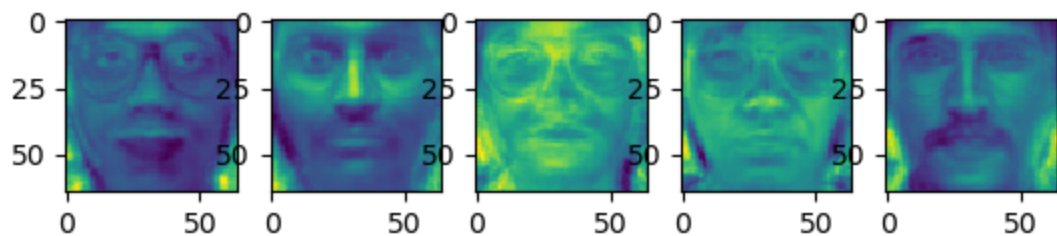
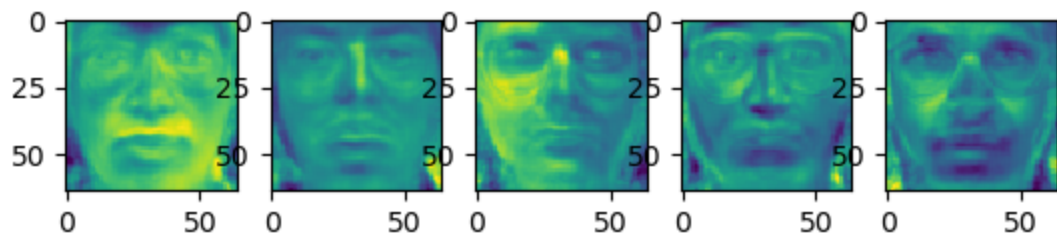




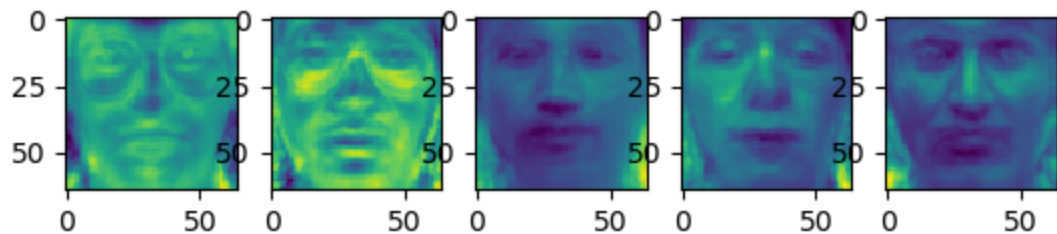
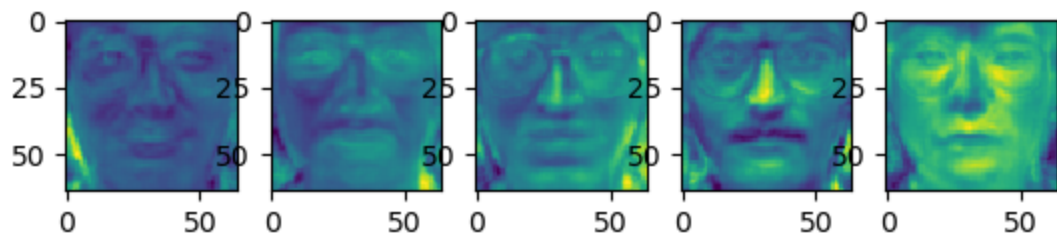
For  $K = 30$



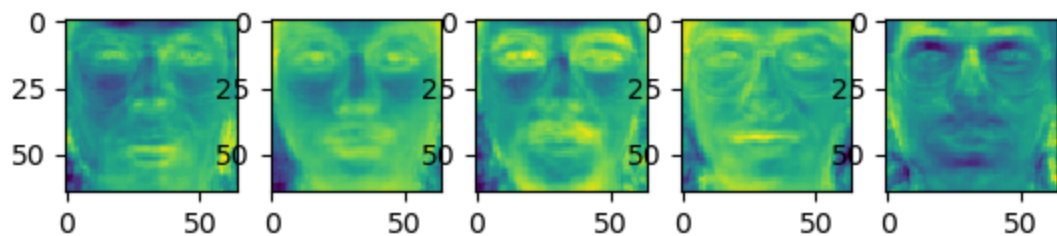
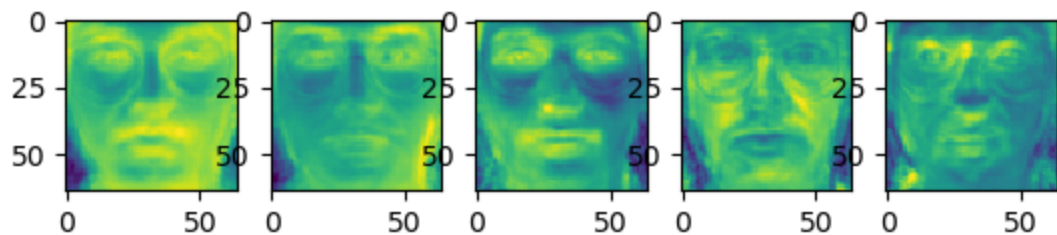
For  $K = 40$



For  $K = 50$



For  $K = 100$

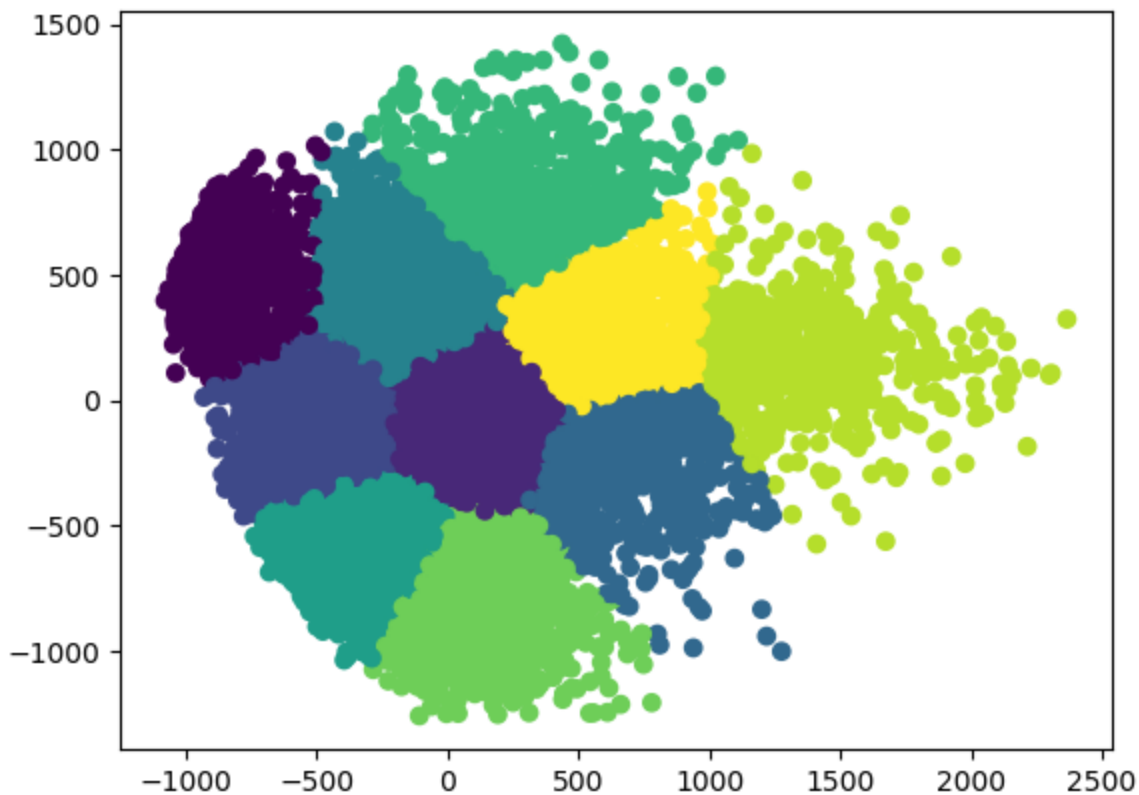


**Programming Part 2:**

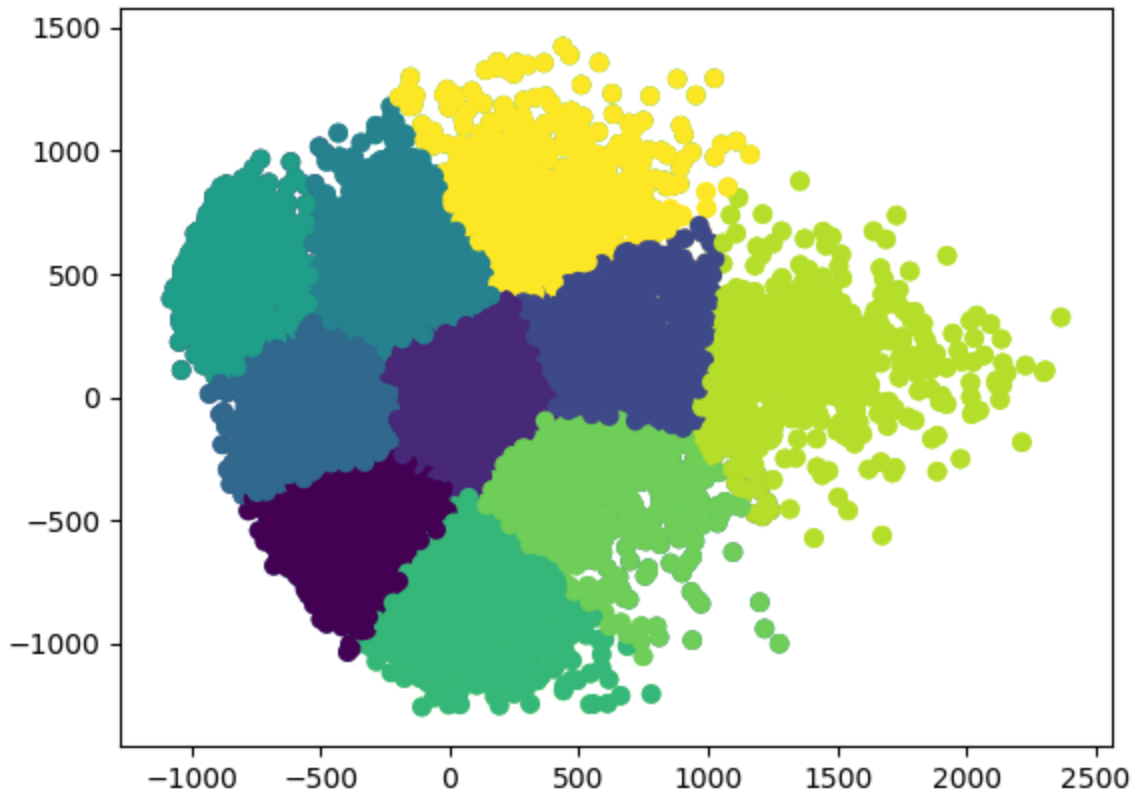
Visually tSNE appears to be better, clusters are more separated as well as on running 10 different initialisations errors in tSNE seems to be less than that of PCA!

Please find the plots from next page.

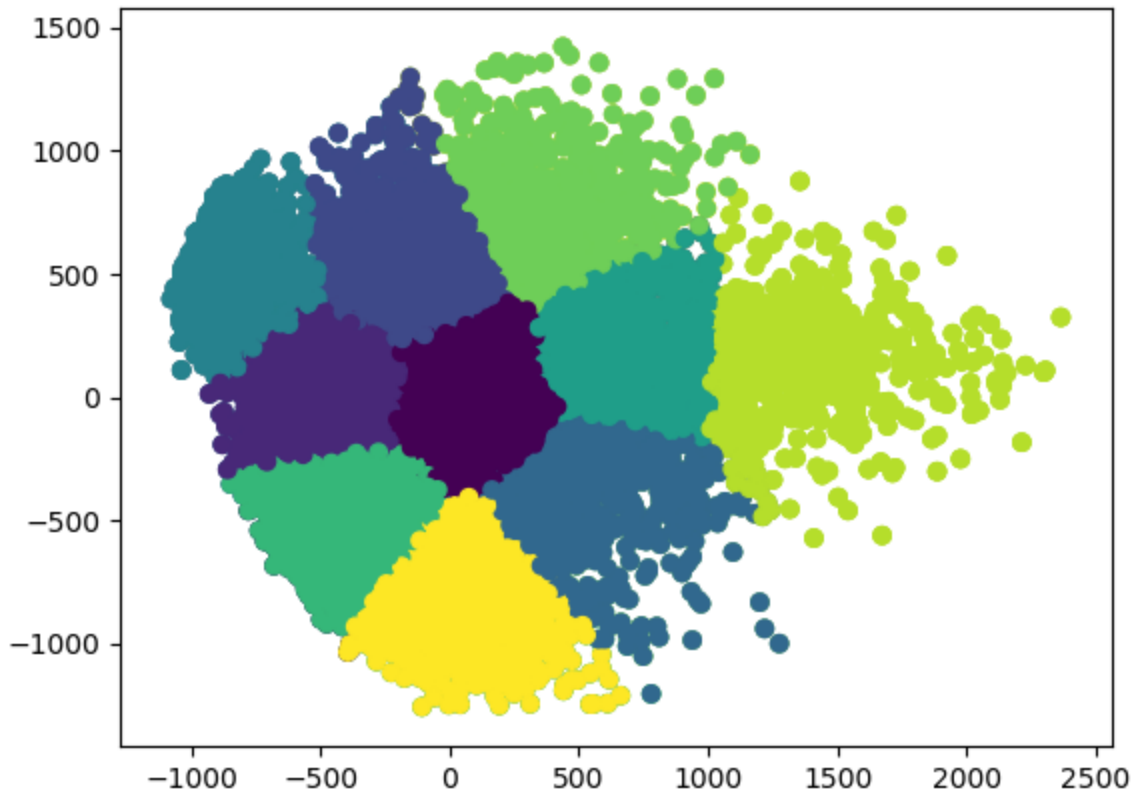
PCA Based Init: 0



PCA Based Init: 1

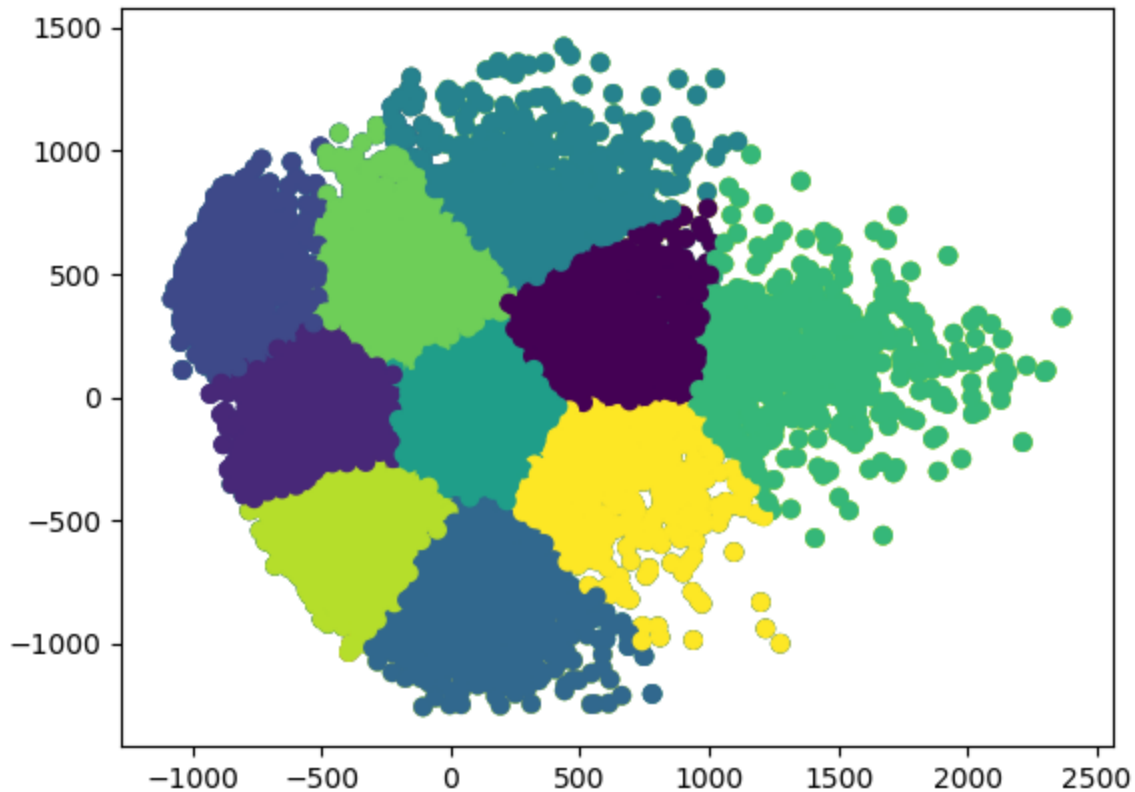


PCA Based Init: 2

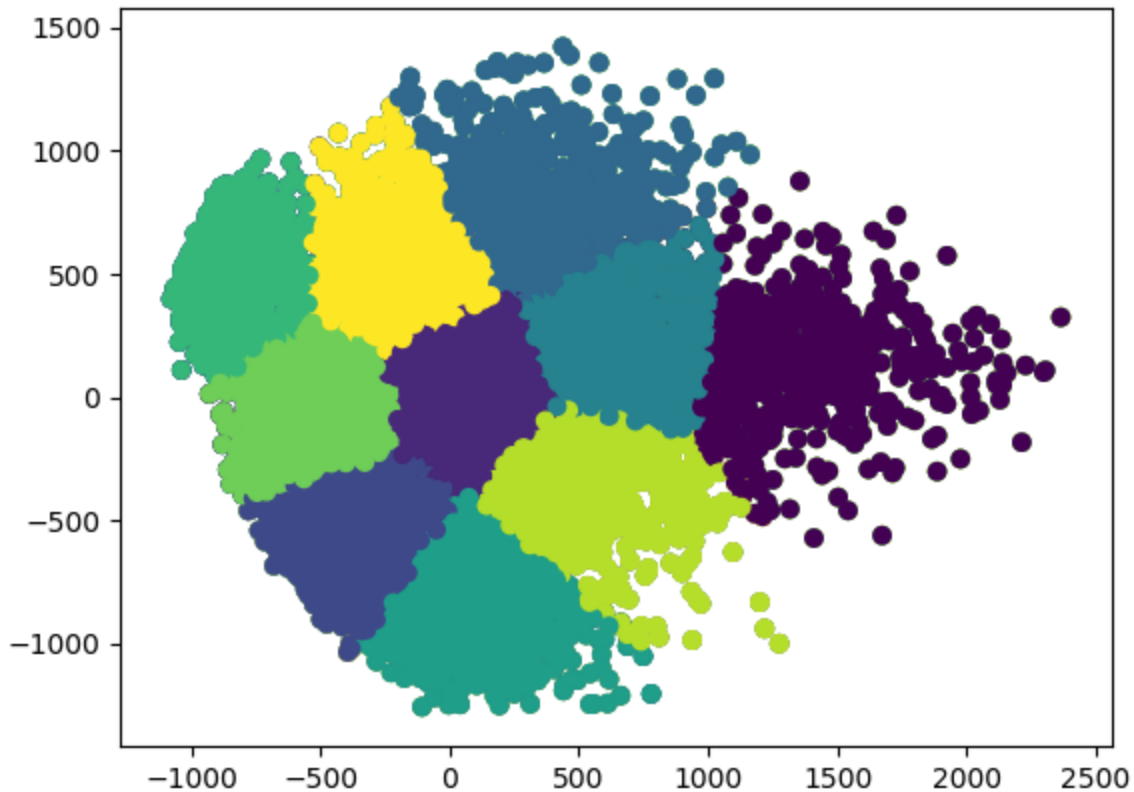




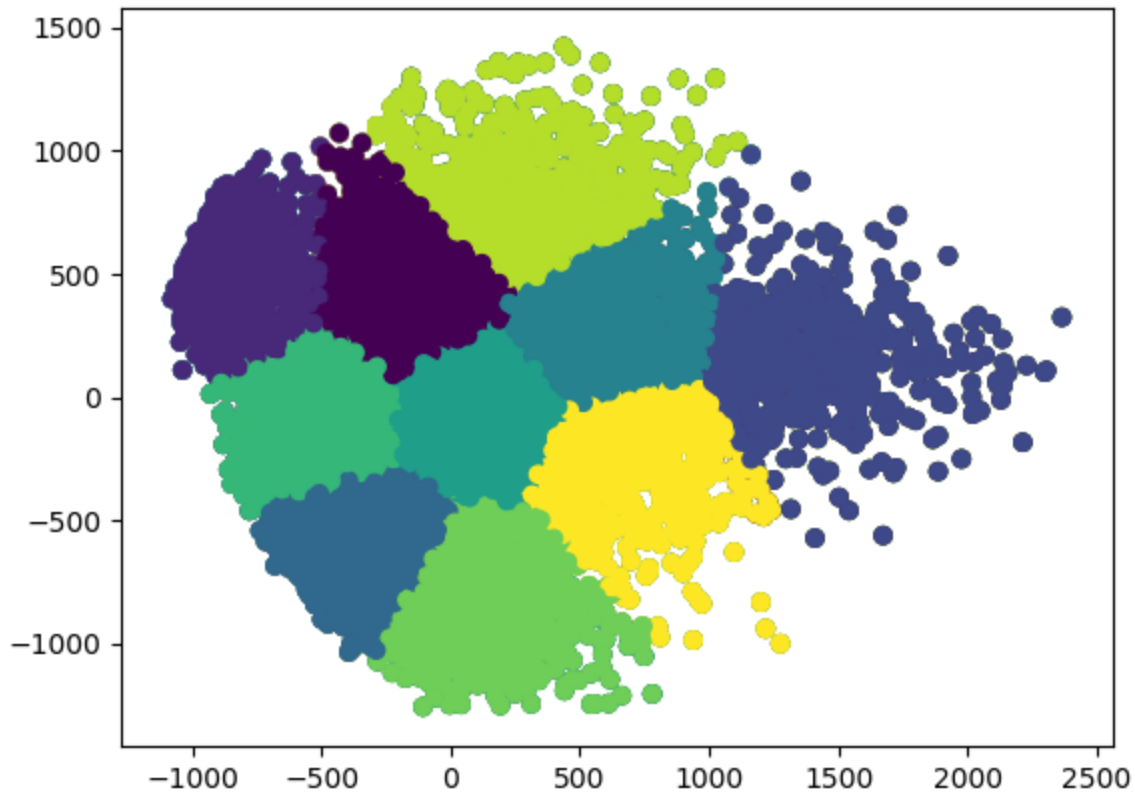
PCA Based Init: 3



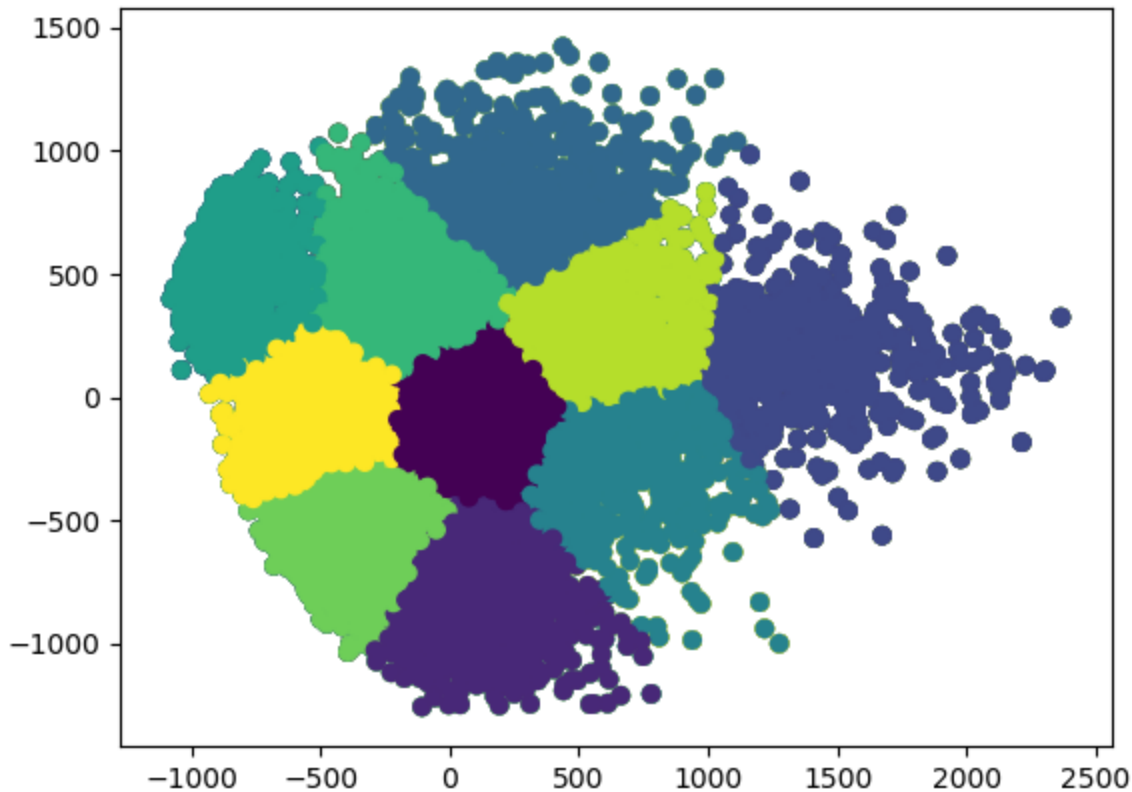
PCA Based Init: 4



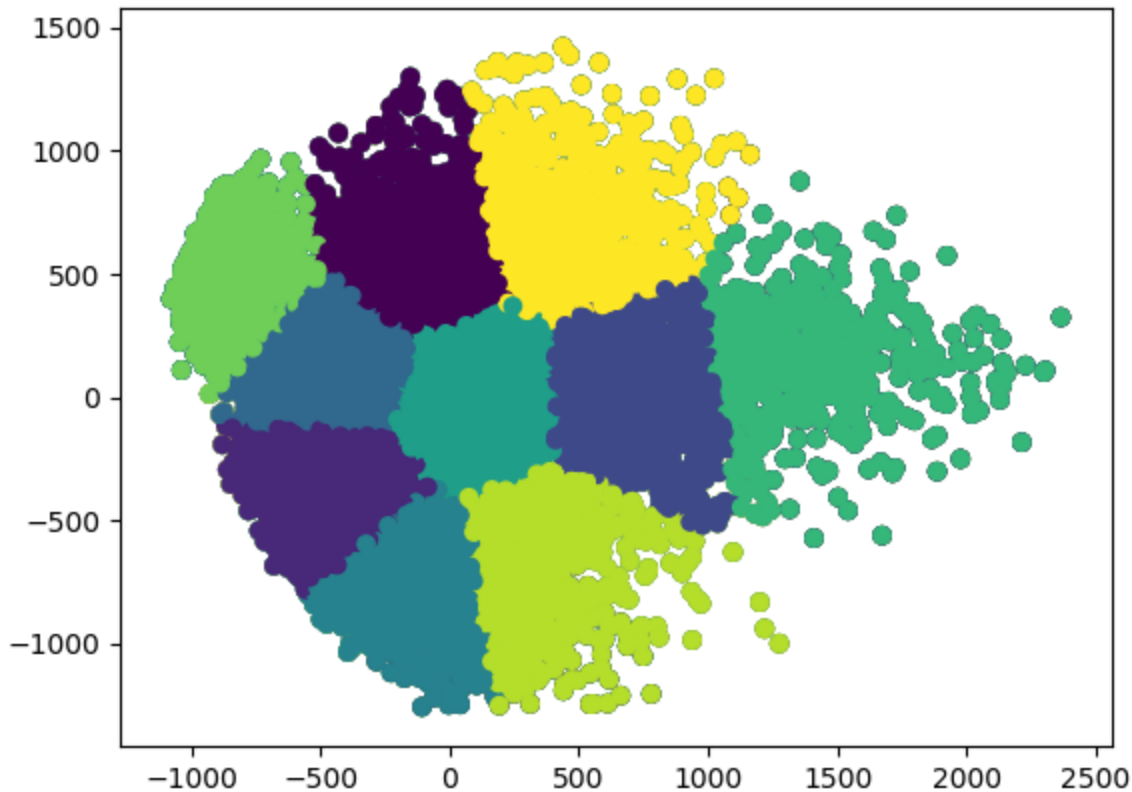
PCA Based Init: 5



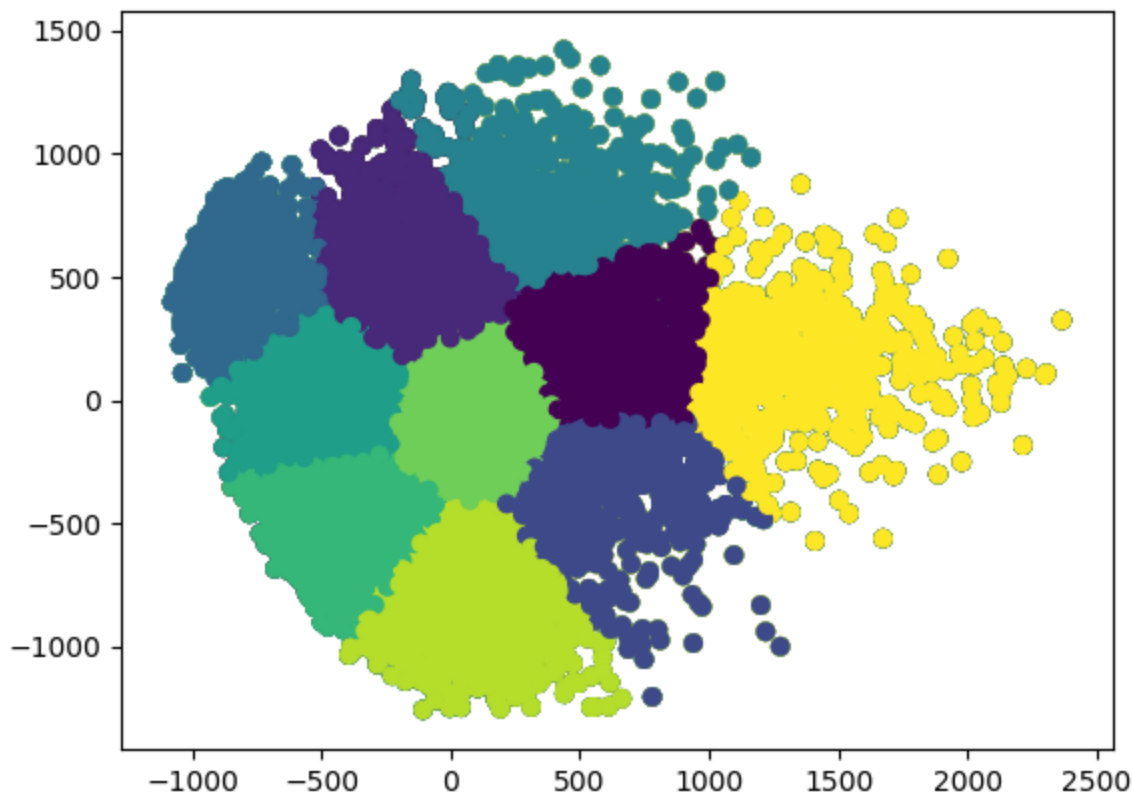
PCA Based Init: 6



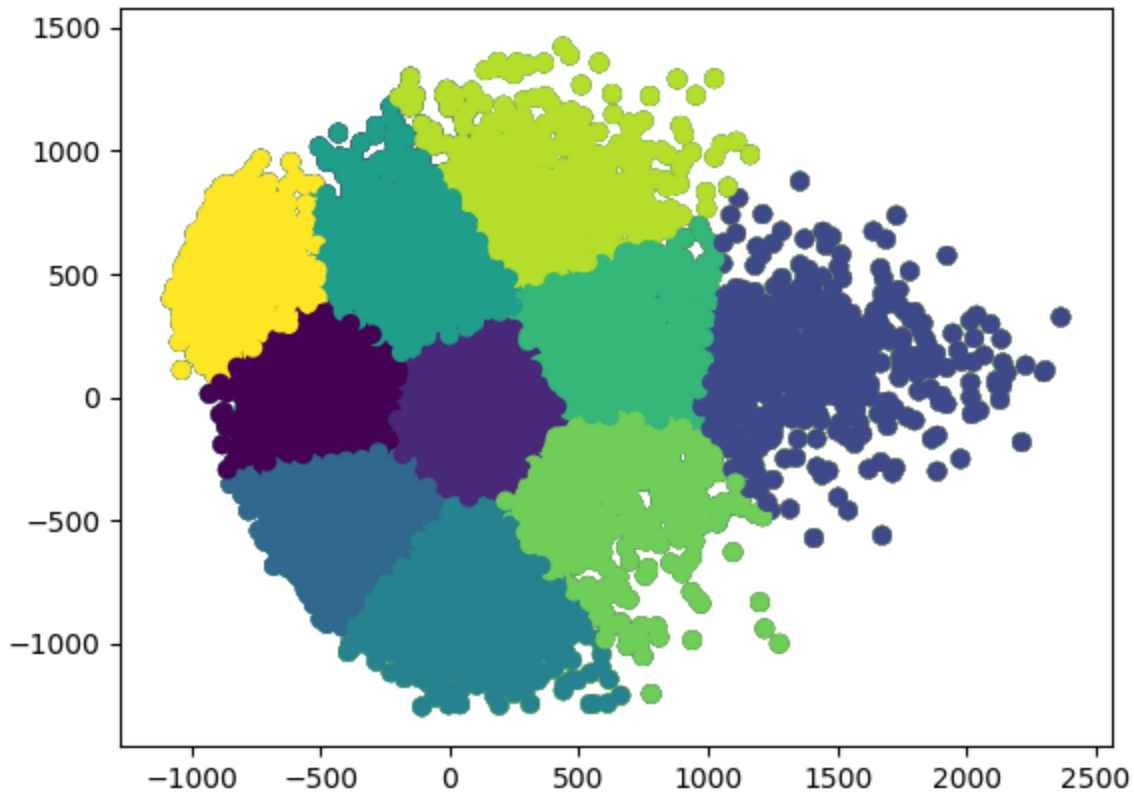
PCA Based Init: 7



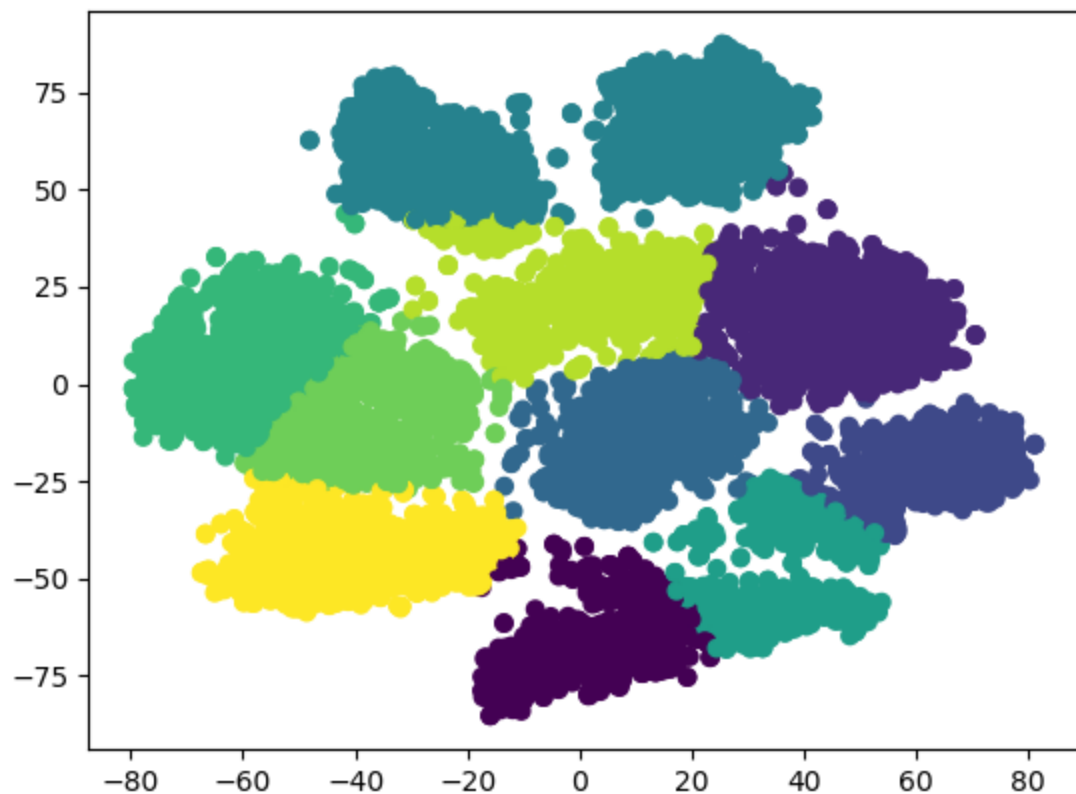
PCA Based Init: 8



PCA Based Init: 9

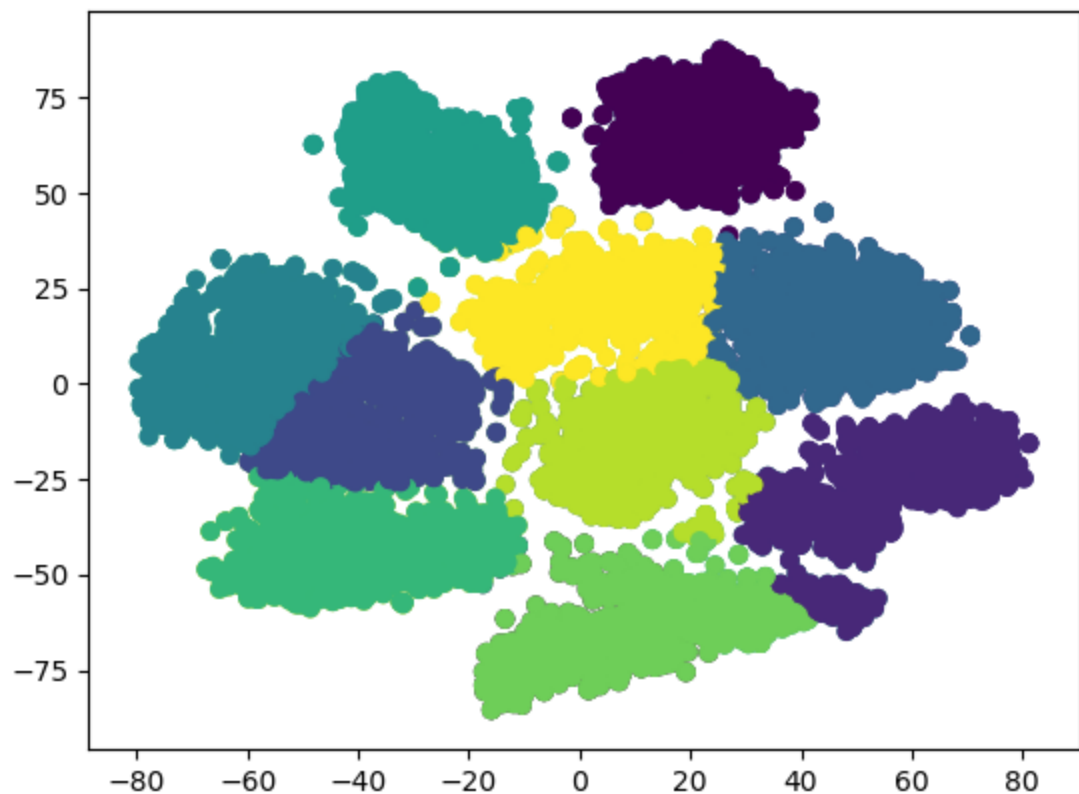


tSNE Based Init: 0

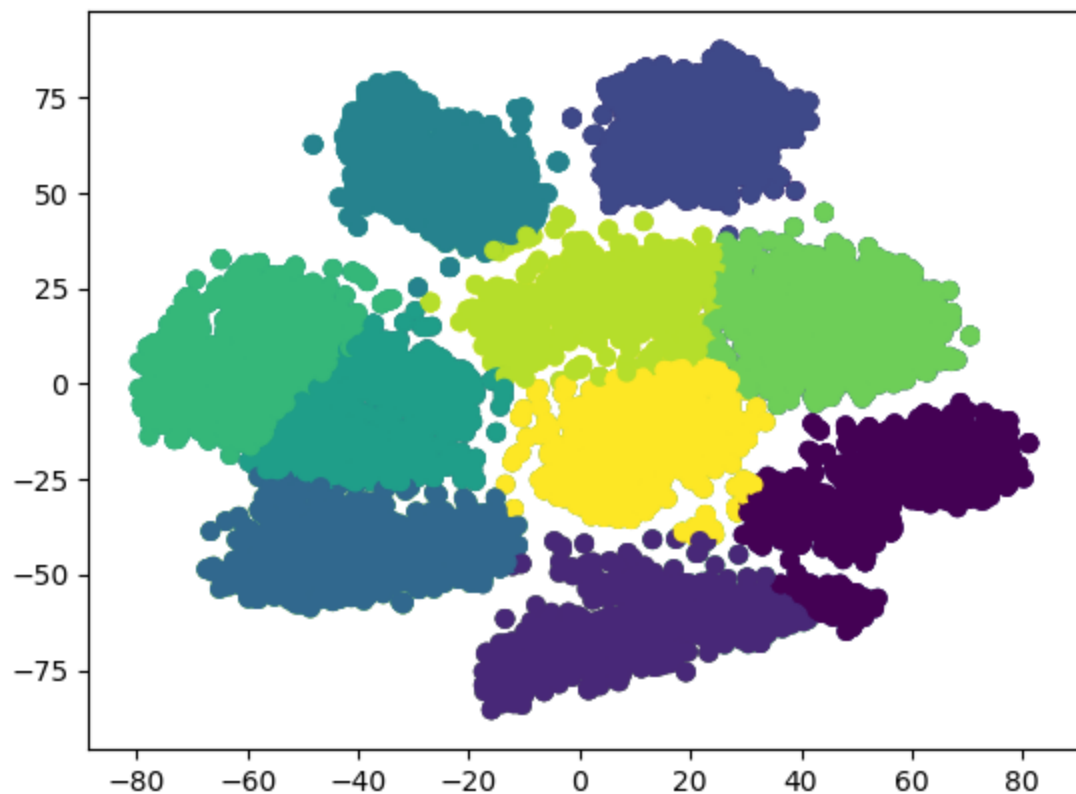




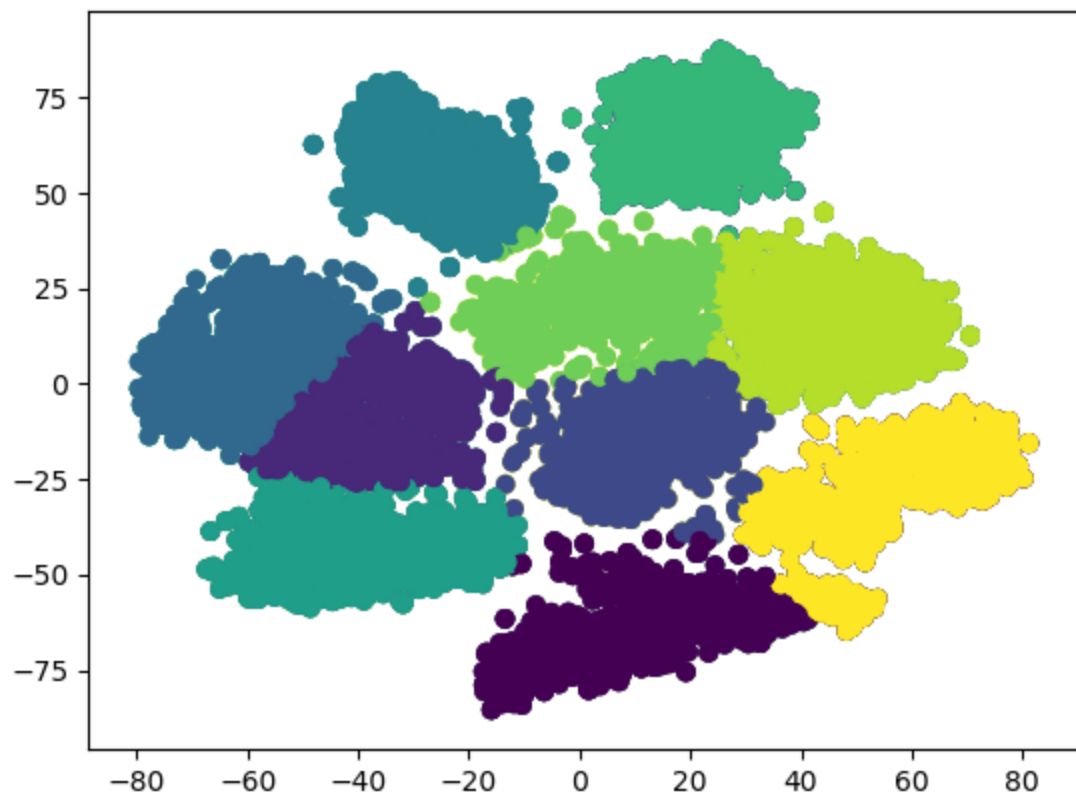
tSNE Based Init: 1



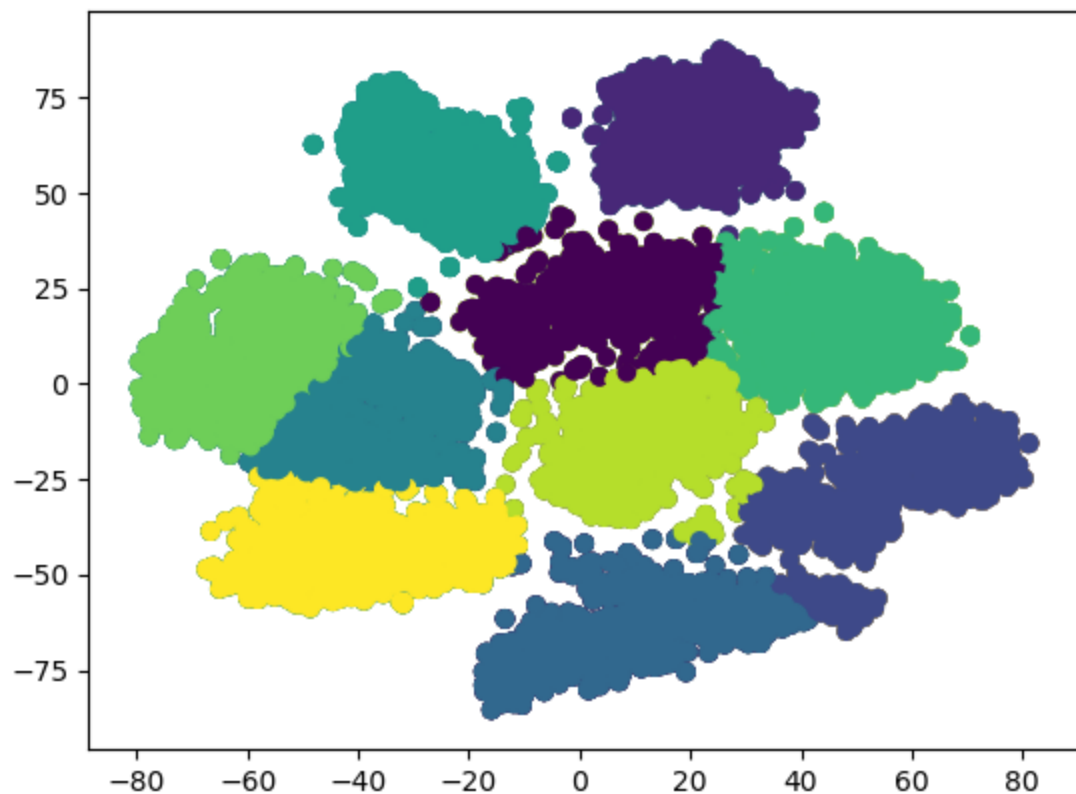
tSNE Based Init: 2



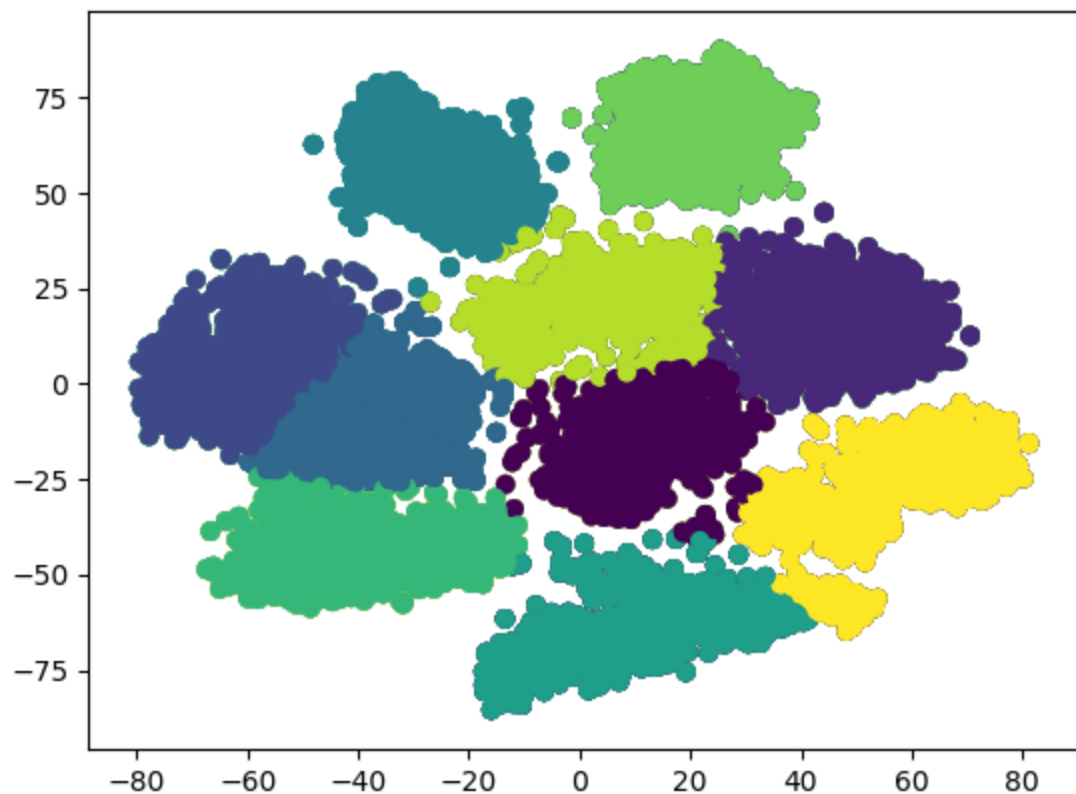
tSNE Based Init: 3



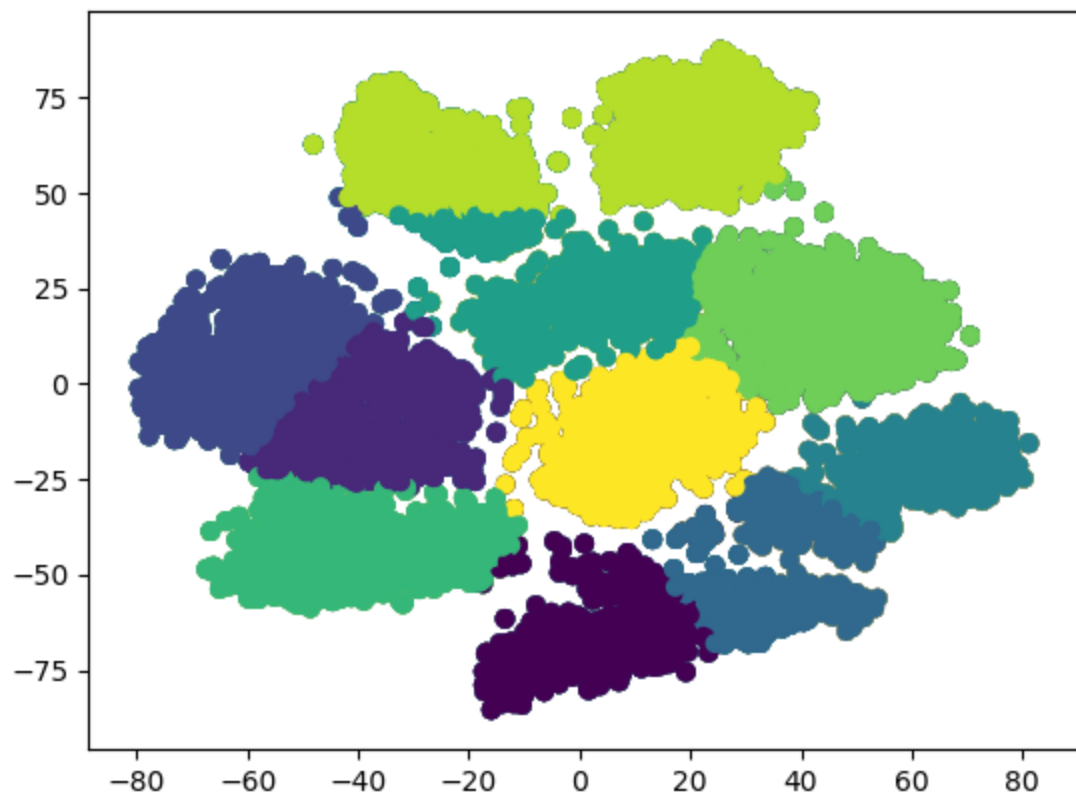
tSNE Based Init: 4



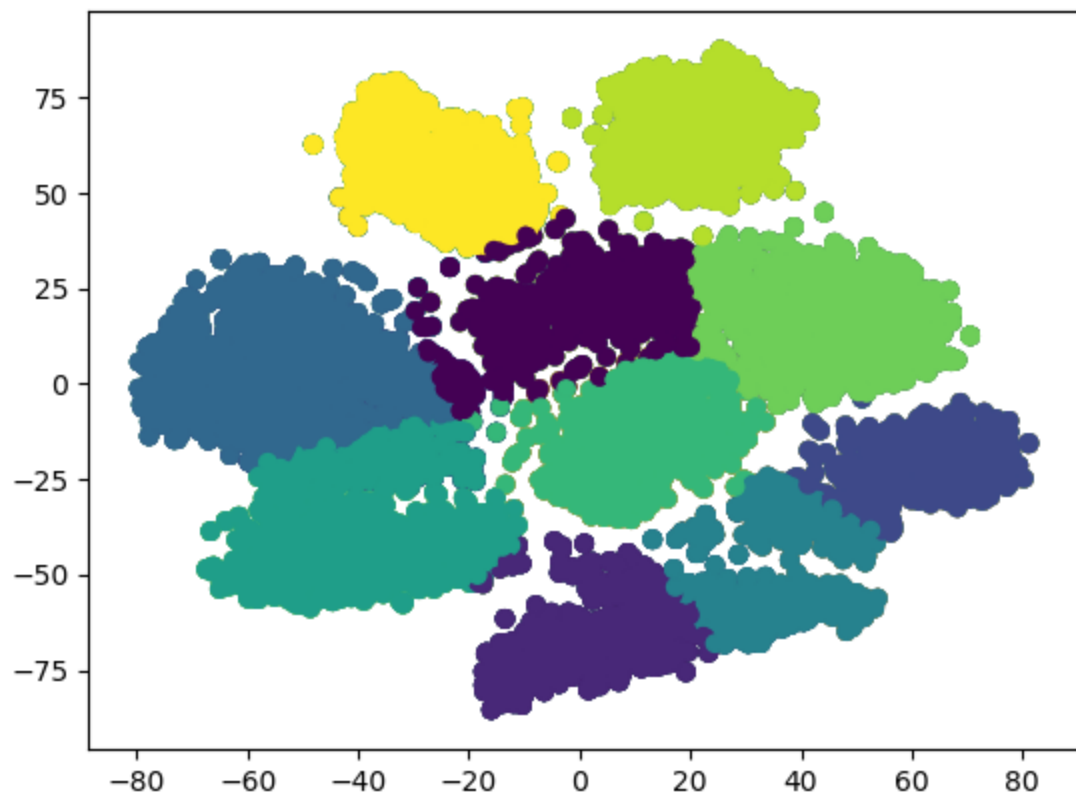
tSNE Based Init: 5



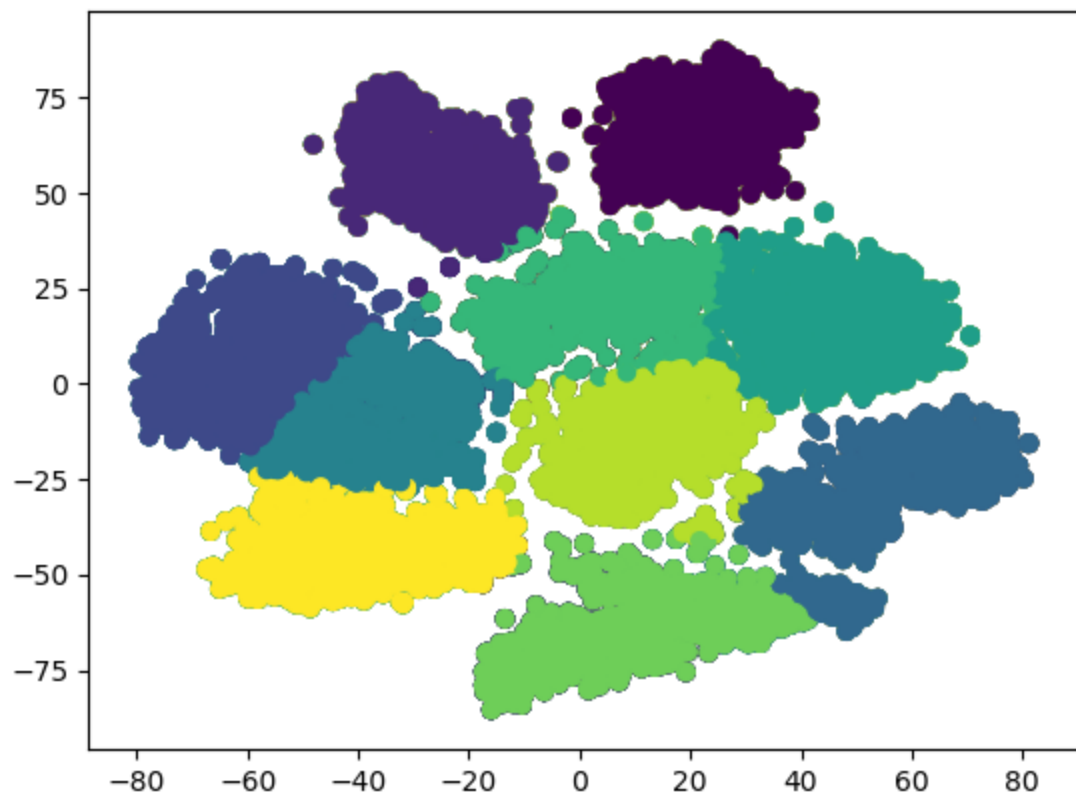
tSNE Based Init: 6



tSNE Based Init: 7



tSNE Based Init: 8





tSNE Based Init: 9

