

Inf2B-CW2

Task 3 Report

→ my_gaussian_classify:

- for each class of the data:
 - find the mean vector $\mu = E[x]$ from training data
 - calculate covariance matrix using $\Sigma = E[(x - \mu)(x - \mu)^T]$ and add epsilon along diagonal
 - calculate for the covariance matrix: natural logarithm determinant; inverse
 - subtract class mean vector from testing data
 - for each vector in testing data:
 - calculate (natural logarithm) posterior probability
- sort posterior probabilities according to `argmax()` to find best-matching class

Time elapsed approx (in seconds) [DICE environment, command line]	covariances	classes	total
	2.38	21.96	24.34

Statistics	N	Nerrs	acc
- initial run gives good accuracy, almost on-par with knn_classify, while only needing ~60% of the total time of knn_classify	7800	1250	83.97%

→ my_improved_gaussian_classify:

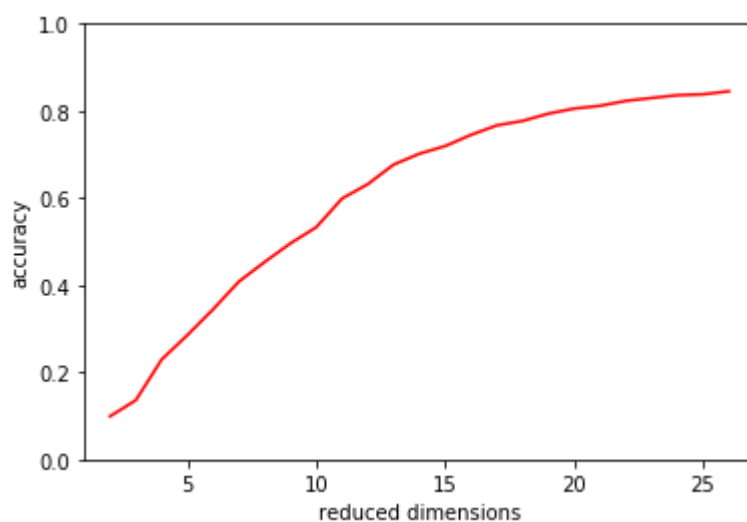
- user-defined dimensionality $d \in [1, 26]$
- for each class of the data:
 - find eigenvalues, eigenvectors from covariance matrix
- select (dims) eigenvectors with highest eigenvalues
- apply eigenvector transformations to training and testing data
- continue with gaussian classifier on modified data as before

→ my_improved_gaussian_system:

- parse arguments to run different preset experiments or set number of dimensions to reduce to
 - usage: `$ my_improved_gaussian_system.py [-e experiment] [-d dims]`
 - choose either to override default behaviour, indicated below

Time elapsed approx (in seconds) [DICE environment, command line]	covs x26 (784x784)	eigen- vectors	PCA transf.	covs x26 (26x26)	class prob.	total
	2.39	7.05	0.41	0.02	0.47	10.34

Statistics / Observations	exp.	dims	ϵ	N	Nerrs	acc
<ul style="list-style-type: none"> - runs ~2x quicker than raw gaussian_classify - accuracy slightly better with similar ϵ - ϵ makes a difference in higher dimensions - ($\epsilon = 0$ is unsuitable, and therefore highlighted in red, as it can result in a matrix with no inverse) - tradeoff here is ~-1% accuracy for 4x higher performance compared to knn_classify - finding matrix eigenvectors results in complex values, that although are probably small enough to be ignored, are still present and may influence results - takes longer real time to implement, more things could go wrong (compared to lower up-and-running time of a knn implementation) 	1	1	0.01	7800	7028	9.90%
	2	1	1e-10	7800	7027	9.91%
	3	2	0.01	7800	6740	13.59%
	4	2	1e-10	7800	6740	13.59%
	5	4	0.01	7800	5581	28.45%
	6	4	1e-10	7800	5573	28.55%
	7	8	1e-10	7800	3930	49.62%
	8	16	1e-10	7800	1818	76.68%
	9	21	1e-10	7800	1381	82.29%
Legend: <div> <div></div> [point of comparison ($\epsilon=0.01$)] <div></div> [ϵ altered] <div></div> [best PCA run ($\epsilon=0.01$)] <div></div> [best PCA result with modded $\epsilon=1e-10$] <div></div> [best theoretical result, $\epsilon=0$, risky] </div>	10	(26)	0.02	7800	1257	83.88%
	11	(26)	0.01	7800	1216	84.40%
	12	(26)	1e-10	7800	1167	85.03%
	13	(26)	0	7800	1167	85.03%



plot of reduced PCA dimensions vs accuracy (with $\epsilon=1e-10$)