

	Hypotheses formulation: Hypotheses formulation: Three machine learning models are used namely Random Forest, Logistic Regression, and Multinomial Naive Bayes. Random Forest is used as baseline model to experiment with various subsets of dataset keeping other two machine learning models to only train with single subset of dataset. All three models are trained on TFIDF comment feature values and Random Forest performed well than others with a testing accuracy of 82.49% while MLP Classifier has 78.88% and Multinomial Naive Bayes has 81.32%. Additionally, few experiments conducted with baseline model for different subsets of datasets. Firstly the model trained with all identities and tfidf feature values which gave testing accuracy of 82.49%. Secondly the baseline model trained with tfidf feature values + religion identities and yielded testing accuracy of 82.37%. Lastly the baseline model trained with tfidf feature values + gender identities and yielded testing accuracy of 82.46%. The Baseline model performed well on all identities + tfidf feature values and also on only gender identity + tfidf values as shown in above accuracy table.
	What are the strength and weaknessses of each machine learning paradigms? Random Forest Strengths: Decision trees are capable of learning non-linear correlations and are somewhat resistant to outliers. In reality, ensembles perform well, winning several conventional (non-deep-learning) machine learning contests. Weaknesses: Individual trees are prone to overfitting when left unconstrained because they might keep branching until they recall the training data. This can be mitigated by employing ensembles. MLP Classifier MLP Classifier Strengths: Multilayer Perceptrons have the advantage of learning non-linear models and the ability to train models in real-time (online learning). MLP include too many parameters because it is fully connected. Each node is connected to another in a very dense web — resulting in redundancy and inefficiency. Multinomial Naive Bayes Multinomial Naive Bayes Strengths: Even though the conditional independence assumption is rarely valid, NB models perform fairly well in practise, especially given their simplicity. They are simple to implement and may expand with your data.
	fairly well in practise, especially given their simplicity. They are simple to implement and may expand with your data. Because of their simplicity, NB models are frequently outperformed by models properly trained and tweaked utilising the preceding approaches. How to adapt the models to close performance gap? For You need to collect relevant data for training, and deploy pipelines that will feed data to the model when it is in production. Which features are most predictive for the task? For When training baseline model on all identities + tfidf feature values meaning complete dataset we find features that are more important in algorithm's decision making which are: Unnamed: 500, Unnamed: 28, Unnamed: 41, Unnamed: 63, Unnamed: 735, White, Comment, Black, Christian, and homosexual gay or lesbian. When trained baseline model with a subset of dataset including tfidf values + gender identity the important features were: Unnamed: 500, Unnamed: 28, Unnamed: 41, Unnamed: 86, Unnamed: 63, Unnamed: 735, other gender. Does method X outperform method Y on the task of comment toxicity because of theoretical property Z. For The baseline model trained on complete dataset outperformed the baseline models trained on subsets of datasets because of all the features (all identities) were present in the complete dataset.
In [496	<pre># Training DBSCAN db = DBSCAN(eps=0.3, min_samples=10).fit(X_train) core_samples_mask = np.zeros_like(db.labels_, dtype=bool) core_samples_mask[db.core_sample_indices_] = True labels = db.labels_ # Number of clusters in labels, ignoring noise if present. n_clusters_ = len(set(labels)) - (1 if -1 in labels else 0) n_noise_ = list(labels).count(-1) print("Estimated number of clusters: %d" % n_clusters_) print("Estimated number of noise points: %d" % n_noise_) print("Homogeneity: %0.3f" % metrics.homogeneity_score(YY, labels)) print("Completeness: %0.3f" % metrics.completeness_score(YY, labels)) print("V-measure: %0.3f" % metrics.v_measure_score(YY, labels)) print("Adjusted Rand Index: %0.3f" % metrics.adjusted_rand_score(YY, labels)) print("Adjusted Rand Index: %0.3f" % metrics.adjusted_rand_score(YY, labels)) print(</pre>
In [49°	
	<pre># y_true = dev_raw_data['Toxicity'] # 1 = len(y_true) 1 = len(y_test) acc = sum([db.labels_[i]==y_test[i] for i in range(1)])/1 print("Test Accuracy: ",acc) Test Accuracy: 0.00173333333333333333333333333333333333</pre>
Out[498]:	0.09135615], [0.03997351, 0.03049004, 0.03508613,, 0.00224874, 0.04441736, 0.14293678]]) Evaluation
In [501 Out[501]:	<pre>3. K-Means Clustering Algorithm # Training K-Means rng = RandomState(42) kmeans = KMeans(2, random_state=rng).fit(X_train) np.round(kmeans.cluster_centers_, decimals=2)</pre>
In [503	# Training Accuracy #y_true = dev_raw_data['Toxicity'] 1 = len(y_test) acc = sum([kmeans.labels_[i]==y_test[i] for i in range(1)])/1 print("Test Accuracy: ",acc) Test Accuracy: 0.7376 Answering Research Question#1: The accuracy obtained from using unlabelled data to train unsupervised learning algorithms is far less than that of labelled data with supervised learning algorithms due to lot of variance in the distributaion of toxicity in different identities. Also a huge amount of sparse data is present in the dataset which makes it harder for unsupervised algorithms to draw a line, a common problem in machine learning is sparse data, which alters the performance of machine learning algorithms and their ability to
In [28]:	calculate accurate predictions. The result is less accurate as we do not have any input data to train from. The model is learning from raw data without any prior knowledge. It is also a time-consuming process.
In []:	Gaussian Mixture Model: