Work Flowchart-DRAFT: Machine Learning Based Prediction To Detect Anti-Cancer Polypeptides Using Deep Learning, Support Vector Machine Sigmoidal Function & Naive Bayes. Start Obtain Tyagi-B dataset & Controls, {2 x *.fasta Format} Convert Polypeptides to % AA & Dipeptide Composition 2D-matrix [~500 polypeptides, ~423 Features] **Exploratory Data Analysis Testing** 1. Investigate: 1st, 2nd 3rd moments of AA & DP, 2. Check for Near Zero Variance of features, 3. Use Boruta RF for Feature Selection, 3. Check Pearson's Correlation Coefficients for multi-collinearity, 4. Conduct Principle Component Analysis for feature selection, 5. K-Means Cluster analysis **Remove** N.Z.V. Features <u>Remove</u> **Boruta RF Features** |R| >= 0.75YES NO Test & Remove Boruta RF **Attribute Columns: Features** NO Change **Combine Columns:** To Obtain Working Dataset. Skewness >= 2.0 YES NO Test: Box-Cox Transformation, **Attribute Columns:** NO Change $\lambda = xi$ & **Transform** Feature i **Combine Columns** To Obtain Working Dataset. **Generate** Skree Plot from PCA <u>Test</u> PC(i) VS Model Accuracy Model Accuracy >= 0.9 NO **ADD** YES PC(i+1) **Combine Columns** PC(i) + Indicator Variable

To Obtain Working Dataset.