

Task #3

1. **Apply a Logistic Regression model** using the functions in the notebook [«Logistic Regression as a Neural Network – BP alg.ipynb»](#) to predict the biological response of a molecule:
data: *bioresponse.csv*,
description from Kaggle: “*The data is in the comma separated values (CSV) format. Each row in this dataset represents a molecule. The first column contains experimental data describing a real biological response; the molecule was seen to elicit this response (1), or not (0). The remaining columns represent molecular descriptors (d1 through d1776), these are calculated properties that can capture some of the characteristics of the molecule - for example size, shape, or elemental constitution. The descriptor matrix has been normalized.*”.
Use **75% of the dataset to train the model**, and the rest of the data to estimate its accuracy.
2. **Modify optimize() function** to implement the **stochastic gradient descent (SGD)** method. Apply it to solve the problem from p.1.
3. **Implement the Adam optimization** method [using the numpy library](#) and compare the accuracy of the model fitted with it with the models trained by the classic GD (p.1) and SGD (p2) algorithms.
4. **For three modifications** of gradient descent (pp. 1-3), **plot learning curves** (dependence of the value of the loss function on the iteration number).
5. **Train the models with different values of the learning rate (at least 5 different values of learning rate)**. Plot the learning curves for all these cases. How does it affect the accuracy of the model?