

Resource Summary Report

Generated by [dkNET](#) on Sep 6, 2026

GDSC data set and random forest model scripts

DOI:10.17504/protocols.io.3j9gkr6

Type: Protocol

Proper Citation

Peter Anderson 2019. GDSC data set and random forest model scripts . protocols.io
dx.doi.org/10.17504/protocols.io.3j9gkr6

Protocol Information

URL: <https://dx.doi.org/10.17504/protocols.io.3j9gkr6>

Authors: Peter Anderson

Summary: This location contains the GDSC data set with 145 oncogene mutation statuses and ~1200 chemical descriptors. For instructions how to run binary classification (to predict compound activity vs inactivity) or regression (to predict values of log IC50), see the README.classification.txt and README.regression.txt files, respectively.

Associated Publications: Lind AP, Anderson PC (2019) Predicting drug activity against cancer cells by random forest models based on minimal genomic information and chemical properties. PLoS ONE 14(7): e0219774. doi: [10.1371/journal.pone.0219774](https://doi.org/10.1371/journal.pone.0219774)

Affiliations: Physical Sciences Division, University of Washington Bothell

External URL: <https://doi.org/10.1371/journal.pone.0219774>

Version: 1

Publication Date: 2019

Proper Citation: Peter Anderson 2019. GDSC data set and random forest model scripts . protocols.io dx.doi.org/10.17504/protocols.io.3j9gkr6

Record Creation Time: 20210329T220529+0000

Record Last Update: 20210329T220938+0000

Data and Source Information

Source: