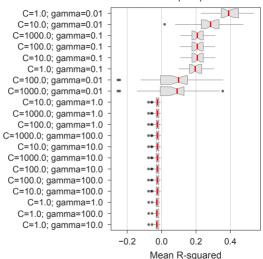SVM - Top 30 proteins

Hyperparameters (top to bottom):
- C=1.0; gamma=0.01
- C=10.0; gamma=0.01
- C=1000.0; gamma=0.1
- C=100.0; gamma=0.1
- C=10.0; gamma=0.1
- C=1.0; gamma=0.1
- C=100.0; gamma=0.01
- C=1000.0; gamma=0.01
- C=10.0; gamma=1.0
- C=1000.0; gamma=1.0
- C=100.0; gamma=1.0
- C=1000.0; gamma=100.0
- C=10.0; gamma=10.0
- C=1000.0; gamma=10.0
- C=100.0; gamma=10.0
- C=100.0; gamma=100.0
- C=10.0; gamma=100.0
- C=1.0; gamma=1.0
- C=1.0; gamma=100.0
- C=1.0; gamma=10.0

Mean R-squared