

**Machine Learning Models to Predict Kidney
Stone Composition and 24-Hour Urine
Abnormalities From Electronic Health Record-
Derived Features
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Introduction:

- Machine learning models allow for analyzing complex data
- Could predict stone type or 24hr urine parameters based on demographic information
- We sought to demonstrate the feasibility of predicting this outcomes using machine learning methods

Methods

Retrospective review of 1245 stone patients w/ 24hr urines

- Demographics
- Past medical history
- Medications
- 24-hour urine values

Target variable: Kidney stone composition

- 0 = Calcium oxide monohydrate
- 1 = Calcium oxide dihydrate
- 2 = Hydroxyapatite
- 3 = Uric acid
- 4 = Other

Input Features

- Demographics:
 - Age
 - BMI
 - Race
 - Sex
- Past medical history:
 - Hypertension
 - Gout
 - Diarrhea/IBD
 - Diabetes
 - CAD/MI
 - Osteoporosis, immobility, hyperparathyroidism
 - Epilepsy, migraine
 - CVA
 - GERD
- Medications:
 - Allopurinol
 - Hydrochlorothiazide
 - Potassium citrate
- 24 hours urine data
 - Volume at 24 hours (Vol24_closest)
 - SSCaOx_closest
 - Ca24_closest
 - Ox24_closest
 - Clt24_closest
 - SSCaP_closest
 - pH_closest
 - SSUA_closest
 - UA24_closest
 - Na24_closest
 - Mg24_closest
 - P24_closest
 - NH424_closest
 - Sul24_closest
 - UUN24_closest
 - Cr24_corrected
 - Cr24kg_corrected
 - Ca24Cr24_corrected
 - Ca24Cr24_closest

80% training and 20% for testing

Model Overview

Pre-processing:

- Scaling continuous variables: mean of 0, SD of 1
- One-hot encoding of categorical variables

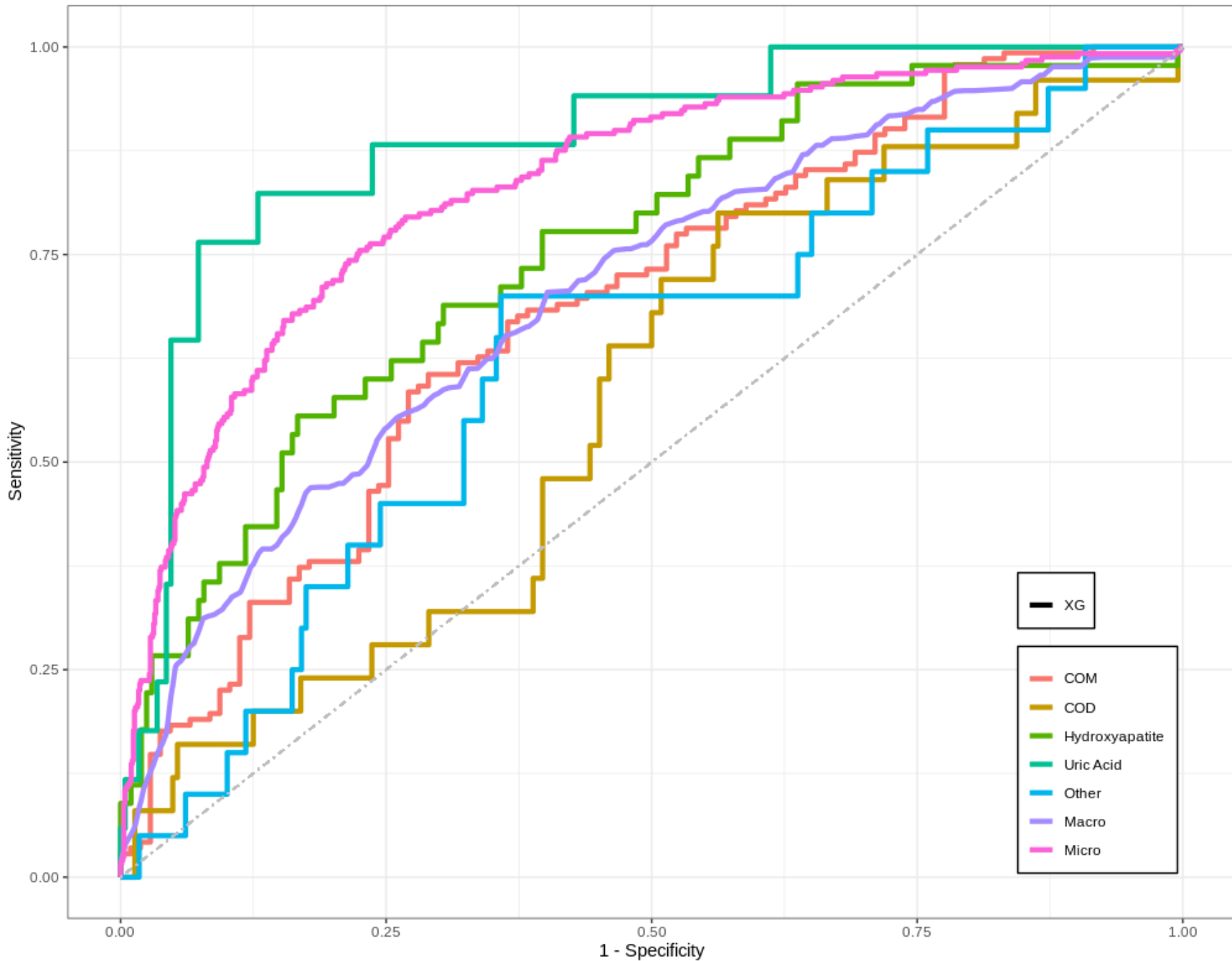
Baseline performance of XGBoost model with

- Booster: gbtree
- Objective: multi:softprob
- Evaluation metrics: mlogloss
- Accuracy: 58.23%

Hyperparameter tuning using random sampling (100 iterations)

- Accuracy: 62.25%

Examining SHAP values of each target class



Accuracy : 0.6225
 95% CI : (0.5591, 0.6829)
 No Information Rate : 0.5703
 P-Value [Acc > NIR] : 0.05416

Kappa : 0.2408

Mcnemar's Test P-Value : NA

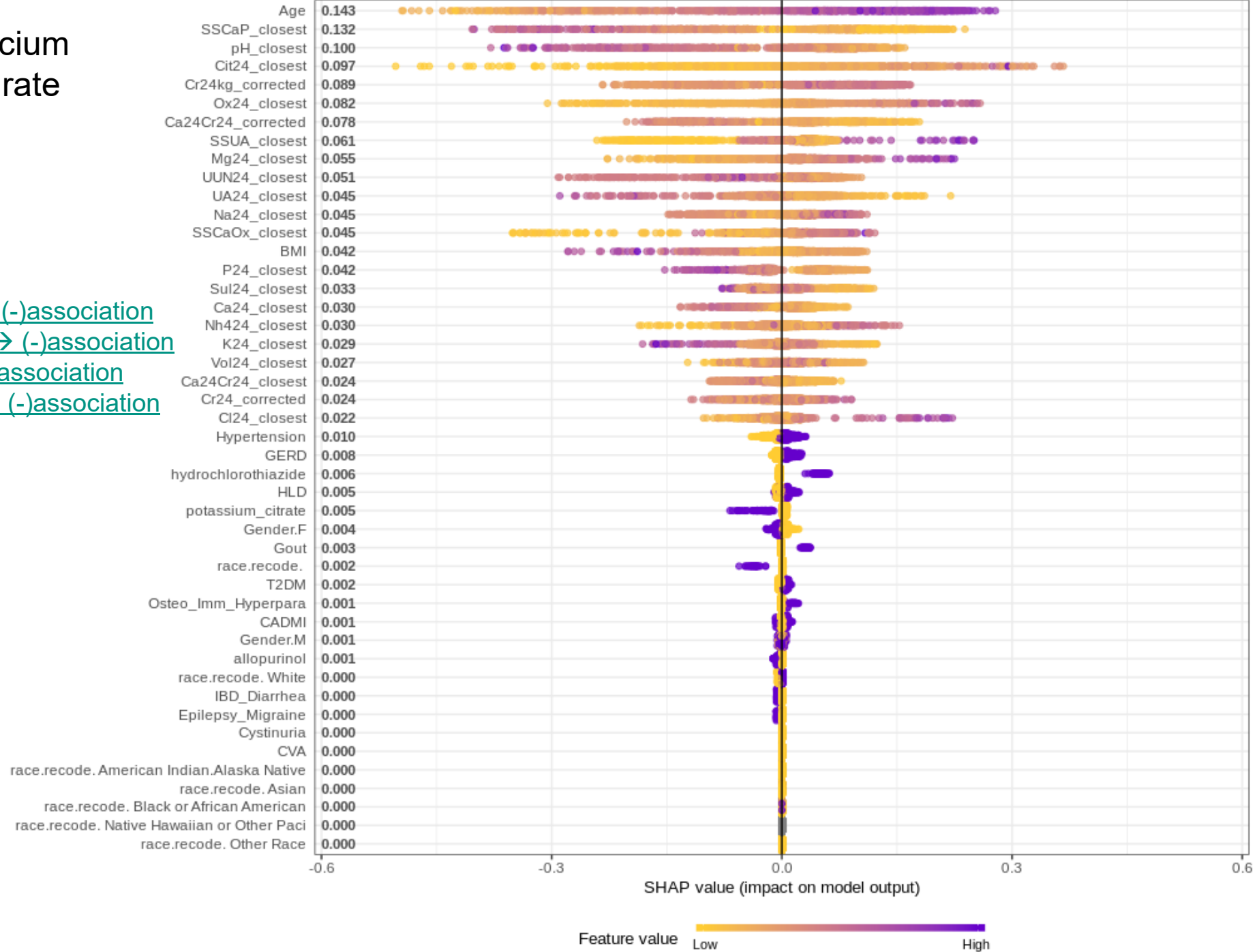
Statistics by Class:

	COM	COD	HA	UA
Sensitivity	0.9366	0.000000	0.40000	0.23529
Specificity	0.2991	0.991071	0.92647	0.99138
Pos Pred Value	0.6394	0.000000	0.54545	0.66667
Neg Pred Value	0.7805	0.898785	0.87500	0.94650
Prevalence	0.5703	0.100402	0.18072	0.06827
Detection Rate	0.5341	0.000000	0.07229	0.01606
Detection Prevalence	0.8353	0.008032	0.13253	0.02410
Balanced Accuracy	0.6178	0.495536	0.66324	0.61334

Target Class: Calcium Oxalate Monohydrate

Associations:

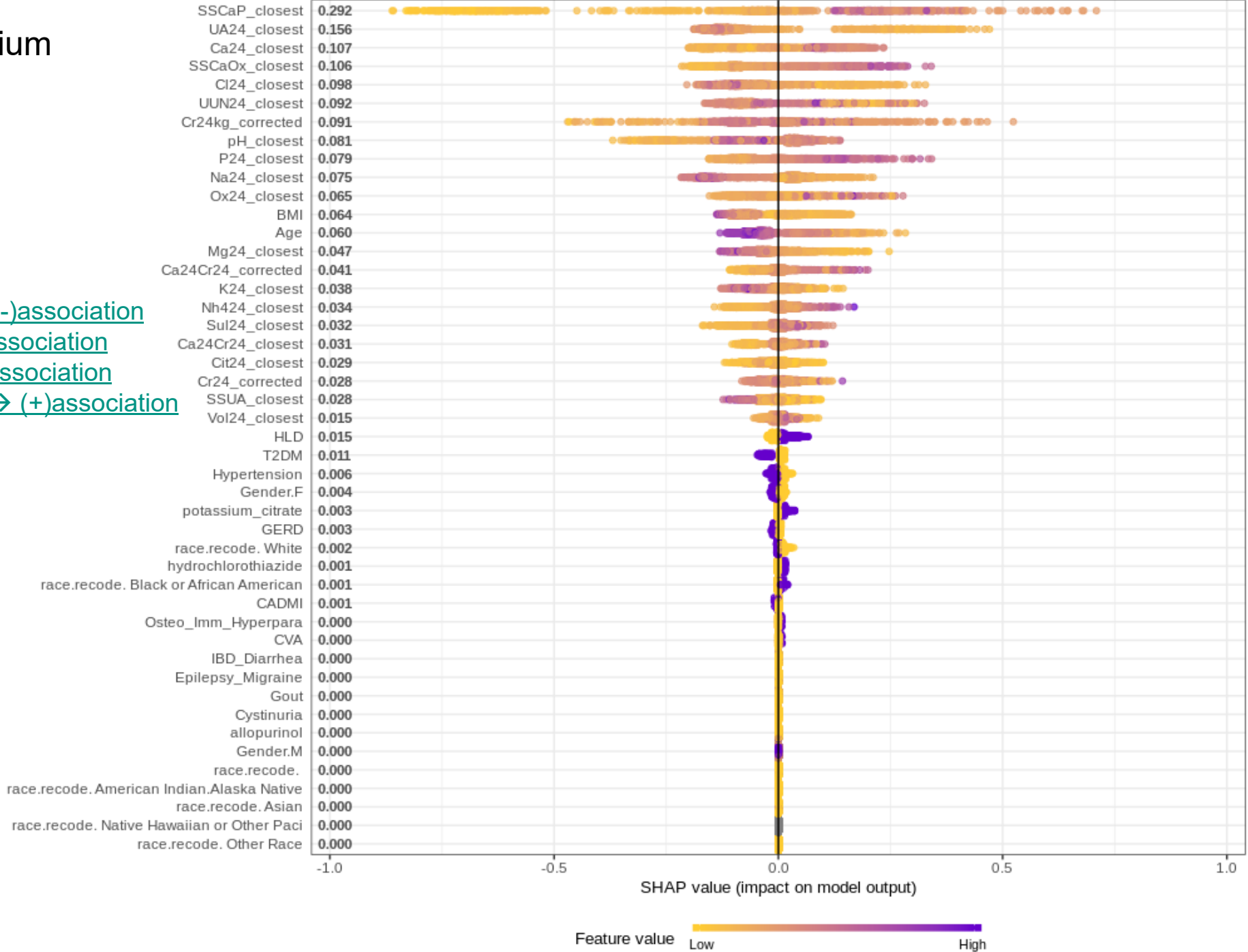
- [Younger Age → \(-\)association](#)
- [Higher SSCaP → \(-\)association](#)
- [Higher pH → \(-\)association](#)
- [Lower Citrate → \(-\)association](#)



Target Class: Calcium Oxalate Dihydrate

Associations:

- [Lower SSCaP → \(-\)association](#)
- [Lower UA → \(+\)association](#)
- [Higher Ca → \(+\)association](#)
- [Higher SSCaOx → \(+\)association](#)

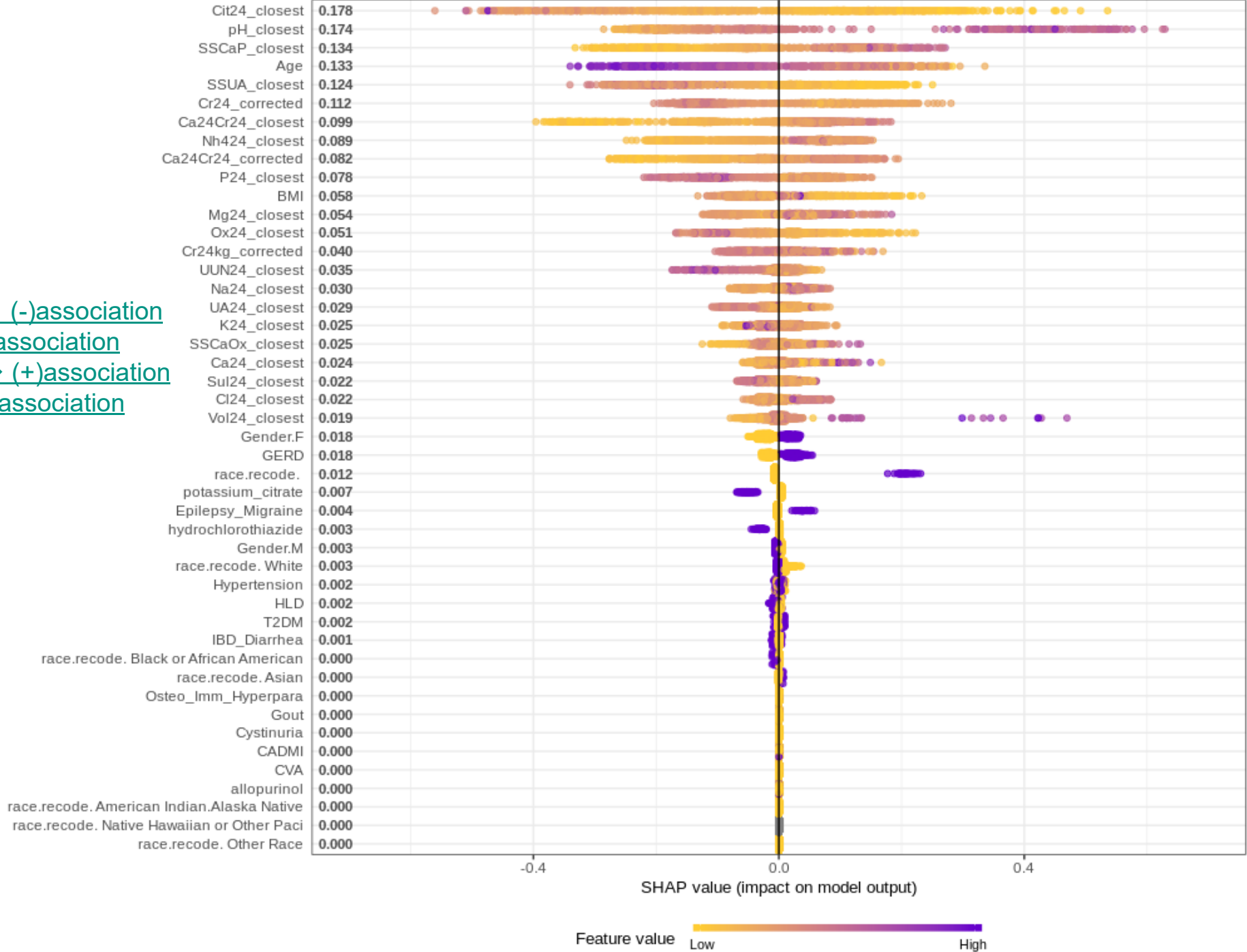


Feature value Low High

Target Class: Hydroxyapatite

Associations:

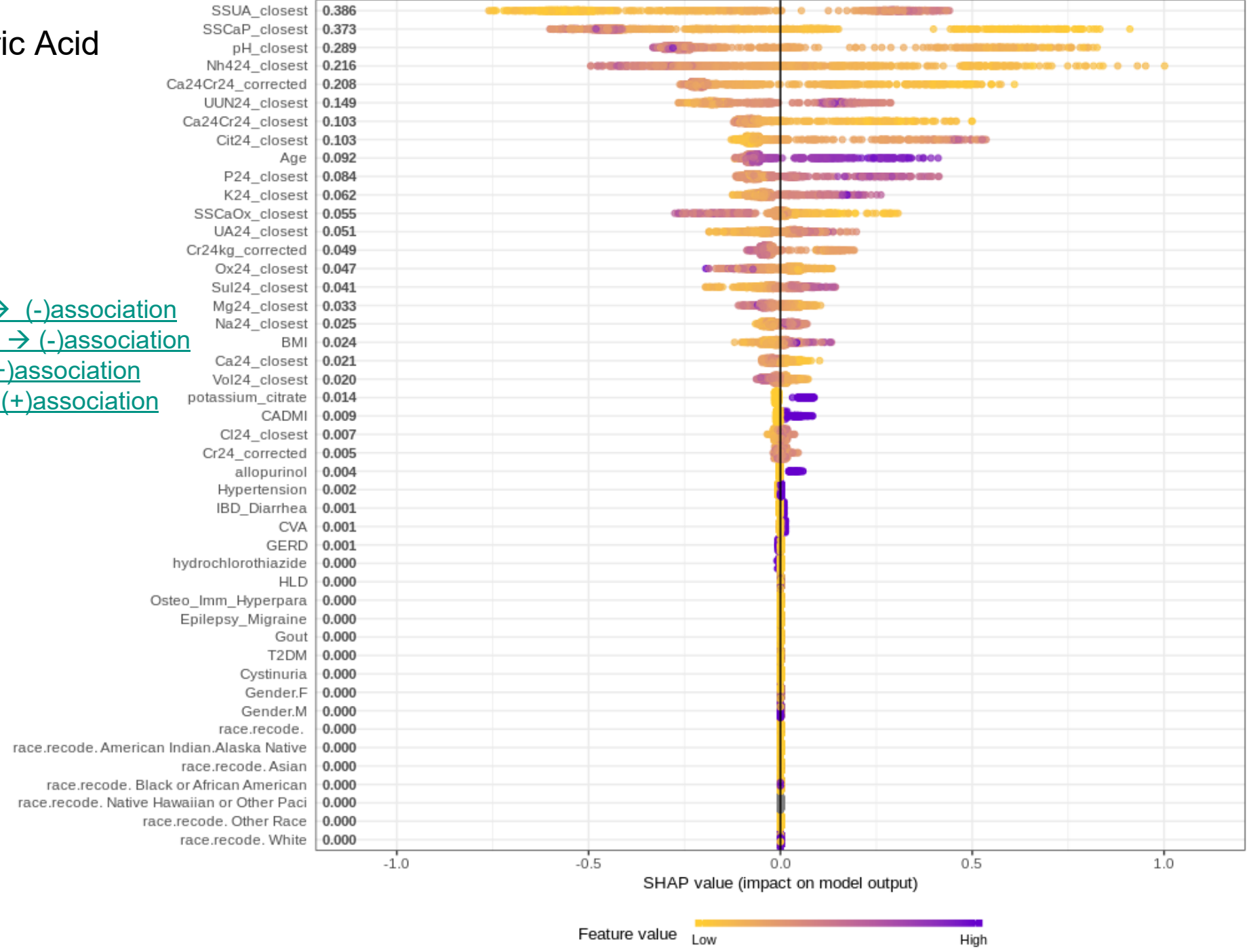
- [Higher Citrate → \(-\)association](#)
- [Higher pH → \(+\)association](#)
- [Higher SSCaP → \(+\)association](#)
- [Higher Age → \(-\)association](#)



Target Class: Uric Acid Stones

Associations:

- [Lower SSUA → \(-\)association](#)
- [Higher SSCaP → \(-\)association](#)
- [Lower pH → \(+\)association](#)
- [Lower NH4 → \(+\)association](#)



Conclusions

- Stone composition prediction appears feasible with machine learning methods
- Features of EHR-derived data were identified and associated with stone composition
- Additional work is necessary to further optimize model performance