

Machine Learning Prediction of High-Current Disruptions with Low-Current Training Data

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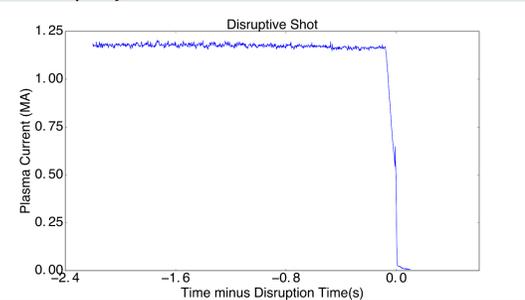
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Objective

- Develop a predictor of high plasma current disruptions using ensembles of regression trees trained with low plasma current data
- Explore methods of stacking the predictions of multiple types of regression ensembles.

Disruptions

- Disruption events occur in tokamaks when the plasma current rapidly decreases to zero.



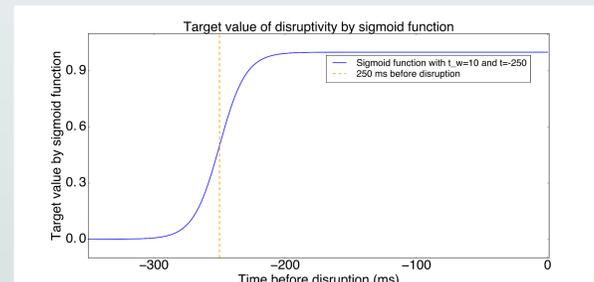
- Disruptions can lead to “runaway” electrons, which cause significant damage to plasma-facing components.

Disruption Prediction

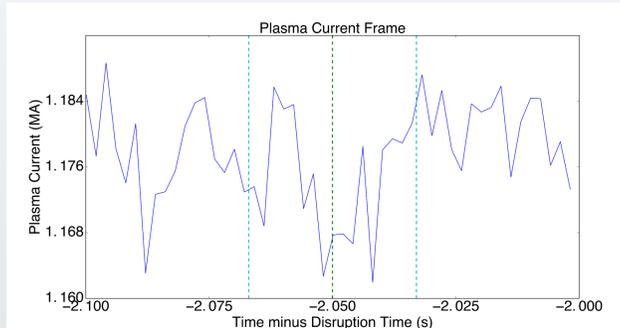
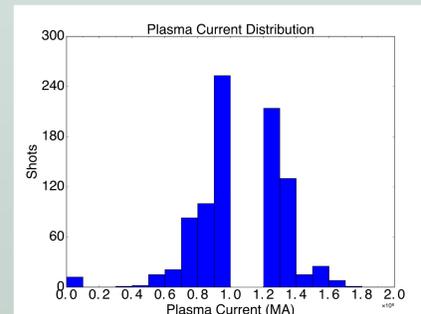
- Complex, nonlinear interactions between parameters preclude pure physics-based disruptors.
- Machine learning techniques for pattern recognition are often employed.
- Artificial Neural Networks (ANNs) used at JET [1], ASDEX [2], ADITYA [3], and DIII-D [4] classify disruptions with signals available in real-time.
- ANNs are “black boxes”; prediction methods are not transparent.
- **Classification and regression trees [5] offer a more transparent approach to prediction.**
- **Challenges to solve before ITER:**
 - Develop an accurate predictor without having existing database of high plasma current shots at ITER
 - Develop a predictor which highlights patterns in parameters of interest for disruption modeling

Data Reduction Process

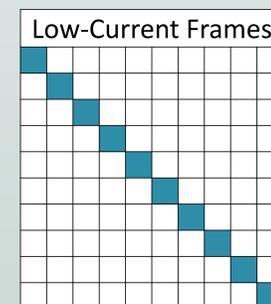
- Time series data for 29 plasma parameters from 630 disrupting shots and 500 non-disrupting shots are collected.
- Time frames are created for each parameter by splitting the time series data into 100 ms windows.
- For disruptive shots, disruptivity target values are created using a sigmoid activation function:



- Ensembles of regression trees are trained with low-current data using k-fold cross validation.
- A disruptivity prediction is made for each high-current frame using the trained ensembles.
- Accuracy is evaluated at a series of disruptivity thresholds between 0 and 1.



- Each frame is split into sub-frames (whole, halves and thirds).
- For each sub-frame, the mean, variance, and trend are calculated.
- The coefficients of a cubic polynomial fit are calculated from the data of the whole frame.
- Thus, one frame has 638 associated values in the dataset (29 parameters, each with 22 values).

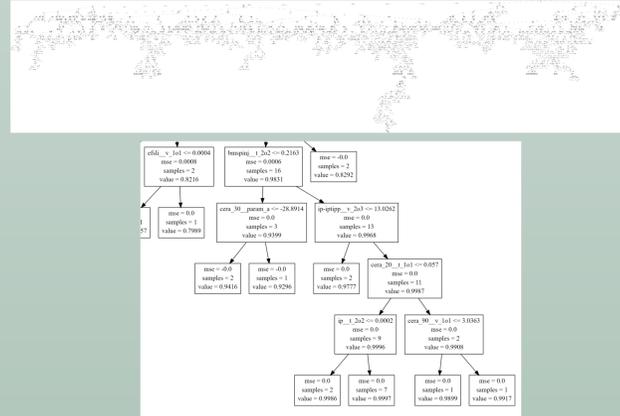


k-fold cross validation:

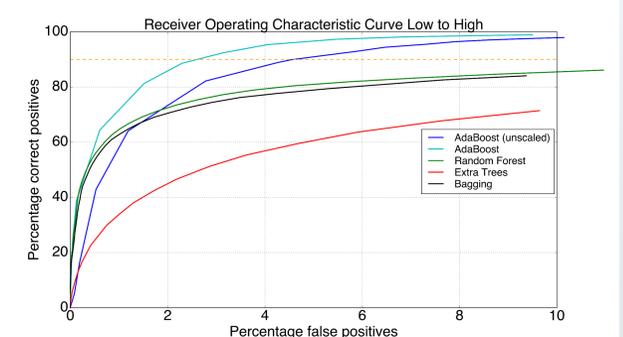
With k=10, low-current data is split into 10 folds. Ten ensembles are trained, each with a different fold excluded from the training set. The final prediction is the mean of the ensembles' predictions.

Regression Tree Ensembles

- Regression trees group frames into terminal nodes, “leaves”, by a series of decisions:
- At each level, the parameter value which reduces the mean squared error of the disruptivity predictions is determined.
- In each final leaf, the disruptivity value predicted is the mean of the values in that leaf.
- Four ensemble methods employed using the scikit-learn [6] machine learning package for the Python programming language:
- Bootstrap aggregating (**Bagging**) [7] trains trees in parallel using subsets of the same size as the full training set, drawn with replacement.
- **Random Forests** [8] extend the bagging method by choosing split candidates from a random subspace of the parameters.
- **Extremely Randomized Trees (Extra Trees)** [9] further extend random forests by choosing the best split from a random set of uniform splits, from a random subspace of the parameters.
- **Adaptive Boosting (AdaBoost)** [10-11] progressively trains an ensemble of weak learners by emphasizing and improving the worst predictions in the previous iteration.



Results



- **93% success rate with 3.2% false positive predictions for AdaBoost with scaled parameters**
- **Using scaled parameters almost halved the false positive predictions at 90% success rate.**
- **Non-boosting methods were less successful than AdaBoost was, but they were more robust.**
- **Low success rates of other algorithms precluded accuracy improvements from stacking regressors.**

Future Work

- Repeat with data spanning a wider range of plasma currents.
- Develop robust weighting algorithm that combines strengths of AdaBoost and Random Forests.
- Optimize parameter list.
- Perform cross-device analysis with normalized parameters.
- Train with 1-d radial profile data.
- Study cases where predictions fail.

References

- [1] B. Cannas et al 2004 Nucl. Fusion 44 68
- [2] B. Cannas et al 2010 Nucl. Fusion 50 075004
- [3] A. Sengupta and P. Ranjan 2000 Nucl. Fusion 40 1993
- [4] D. Wroblewski et al 1997 Nucl. Fusion 37 725
- [5] Breiman, Leo; Friedman, J. H.; Olshen, R. A.; Stone, C. J. (1984). *Classification and regression trees*. Monterey, CA: Wadsworth & Brooks/Cole Advanced Books & Software.
- [6] Pedregosa et al., “Scikit-learn: Machine Learning in Python”, JMLR 12, pp. 2825-2830, 2011
- [7] Breiman, Leo (1996). “Bagging predictors”. *Machine Learning*. 24 (2): 123-140.
- [8] Breiman, Leo (2001). “Random Forests”. *Machine Learning*. 45 (1): 5-32.
- [9] P. Geurts, D. Ernst., and L. Wehenkel, “Extremely randomized trees”, *Machine Learning*, 63(1), 3-42, 2006.
- [10] Y. Freund, R. Schapire, “A Decision-Theoretic Generalization of on-Line Learning and an Application to Boosting”, 1995.
- [11] H. Drucker, “Improving Regressors using Boosting Techniques”, 1997.

Acknowledgements

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