Salty Project



High Performance Computing and Machine Learning for Engineering Solvents

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Motivation and Background

Salty Project is a webapp where researchers can:

- 1. Experiment with predictive computational models.
- 2. Easily access IL thermophysical data.

ILs (ionic liquids) are excellent solvents due to their low volatility, low vapor pressure, and range of polarities. This has lead to exponential growth in interest for developing new ILs. ILs have applications in batteries and biomass transport.

What is the problem?

NIST (The National Institute of Standards and Technology) provides ILThermo, a database with over 1 million IL thermophysical datapoints. However, accessing the data requires navigating a complicated filter then pasting the results into an Excel document. This is a significant barrier to IL researchers who require thousands of datapoints.

What is the solution?

Salty Project is driven by data visualization. The webapp provides an interactive model in the form of a graphical interface. Users can test a predictive computational model and easily read thousands of datapoints. Additionally, all IL data on the webapp is available in .csv file format.

Computational Modeling

Why use high performance computing (HPC)? HPC gives researchers a novel method for materials development. Traditional labs undergo exhaustive synthesis and testing when developing new materials. With HPC, researchers can build computational models that predict thermophysical properties of molecules by analyzing empirically obtained data.

Which model does Salty Project use?

Salty Project utilizes a model which predicts the viscosities of ILs. The model, LASSO (Least Absolute Shrinkage and Selection Operator), is a variation of linear regression. Basic = linear regression finds the best linear function for fitting a dataset. LASSO improves on the method by adding the hyperparameter alpha. Increasing alpha removes certain features (input) by reducing their coefficient to 0.

This penalization of features prevents the model from becoming overly flexible and fitting to irrelevant outlier datapoints, a phenomena known as overfitting. Choosing alpha depends on minimizing mean squared error. Features in Salty Project include experimental conditions, and physiochemical descriptors of the cationic and anionic moieties.

Linear function: $y = m_1x + m_2x + ... + m_n x + b$

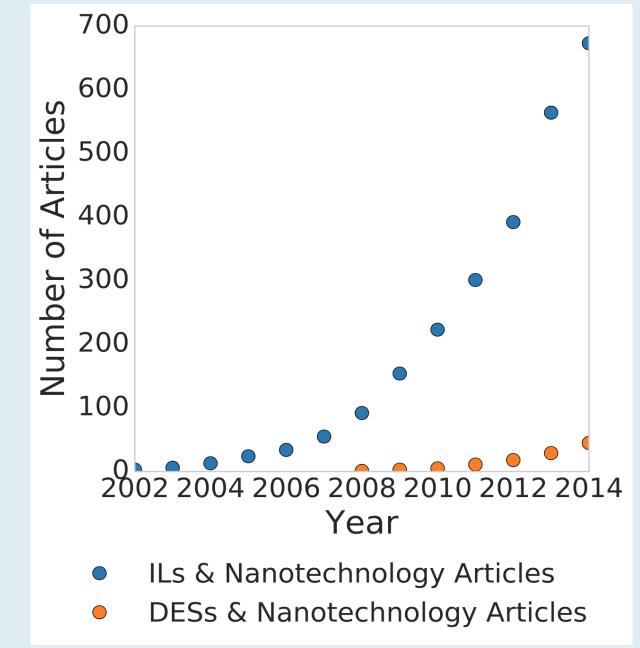


Figure 1: Ionic liquid publications by year. Abo-Hamad, et al. (2015). Potential applications of deep eutectic solvents in nanotechnology. Chemical Enginering Journal, 273, 551–567. http://doi.org/10.1016/j.cej.2015.03.091.

- 1-ethyl-3-methylimidazolium ethyl sulfate 1-ethyl-3-methylimidazolium dicyanamide 1-hexyl-3-methylimidazolium bis[(trifluoromethyl)sulfonyl]imide 1-ethyl-3-methylimidazolium bis[(trifluoromethyl)sulfonyl]imide 1-methyl-3-octylimidazolium tetrafluoroborate
- 1-hexyl-3-methylimidazolium tris(pentafluoroethyl)trifluorophosphate 1-ethyl-3-methylimidazolium tetrafluoroborate
- 1-ethyl-3-methylimidazolium trifluoromethanesulfonate 1-hexyl-3-methylimidazolium hexafluorophosphate
- 1-hexyl-3-methylimidazolium trifluoromethanesulfonate
- trihexyl(tetradecyl)phosphonium chloride 1-ethyl-3-methylimidazolium hexafluorophosphate
- 1-hexyl-3-methylimidazolium tetrafluoroborate trihexyl(tetradecyl)phosphonium tris(pentafluoroethyl)trifluorophospha
- propylammonium acetate
- 1-octyl-3-methylimidazolium hexafluorophosphate
- 1-ethyl-3-methylimidazolium acetate
- trihexyl(tetradecyl)phosphonium bis[(trifluoromethyl)sulfonyl]imide
- 1-ethyl-3-methylimidazolium thiocyanate
- 1-octyl-3-methylimidazolium trifluoromethanesulfonate

Figure 2: D3 produced legend with names of IL salts.

Webapp Development

The framework of the webapp uses HTML. The frontend (user interface) uses D3. The backend (sever, application, and database) uses Python.

What features does Salty Project include?

D3 is a JavaScript library for data visualization. As the frontend includes the interactive LASSO model, D3 provided the best option for interacting with the IL data and viewing the model output. The interactive LASSO model includes:

- 1. 9,000 viscosity datapoints.
- 2. Tooltips for identifying datapoints.
- 3. A color-coded legend for each type of salt.
- 4. Panning and zooming functionality.
- 5. A slider bar for adjusting the alpha value.

With plans to add:

- 1. Filter-sort interactivity to the legend.
- 2. Dynamic updates of model improvements.

The current model interacts with static underlying datasets from the Python backend. In the future, creating a dynamic link between the D3 frontend and Python backend will allow the user to train and test their model interactively.

RDKit Model Data Feature Training Aquirement Generation

31, 326 density datapoints received.

9,000 datapoints for 22 types of salts used in current model.

Figures 3 & 4: Model flexibility and accuracy.

Figures 5 & 6: Effect of alpha on a LASSO model.

www.analyticsvidhya.com/blog/2016/01/complete-tutorial-ridge-lasso-regression-python/.

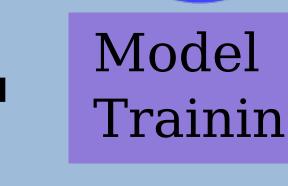
1.0 1.5 2.0 2.5 3.0 3.5 4.0 4.5 5.0 5.5

James et. al. Introduction to Statistical Learning, 2015, 34.

Experimental conditions.

Molecular features.

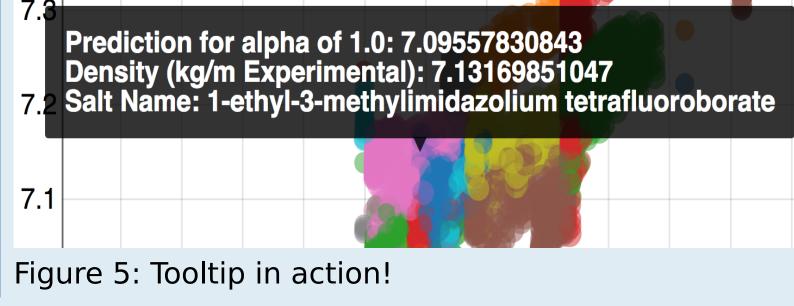
Python-based machine learning package.



6.80 6.85 6.90 6.95 7.00 7.05 7.10 7.15 7.20 7.25 7.30 7.35

Used in development of LASSO model.

Open source tool for developing statistical models.



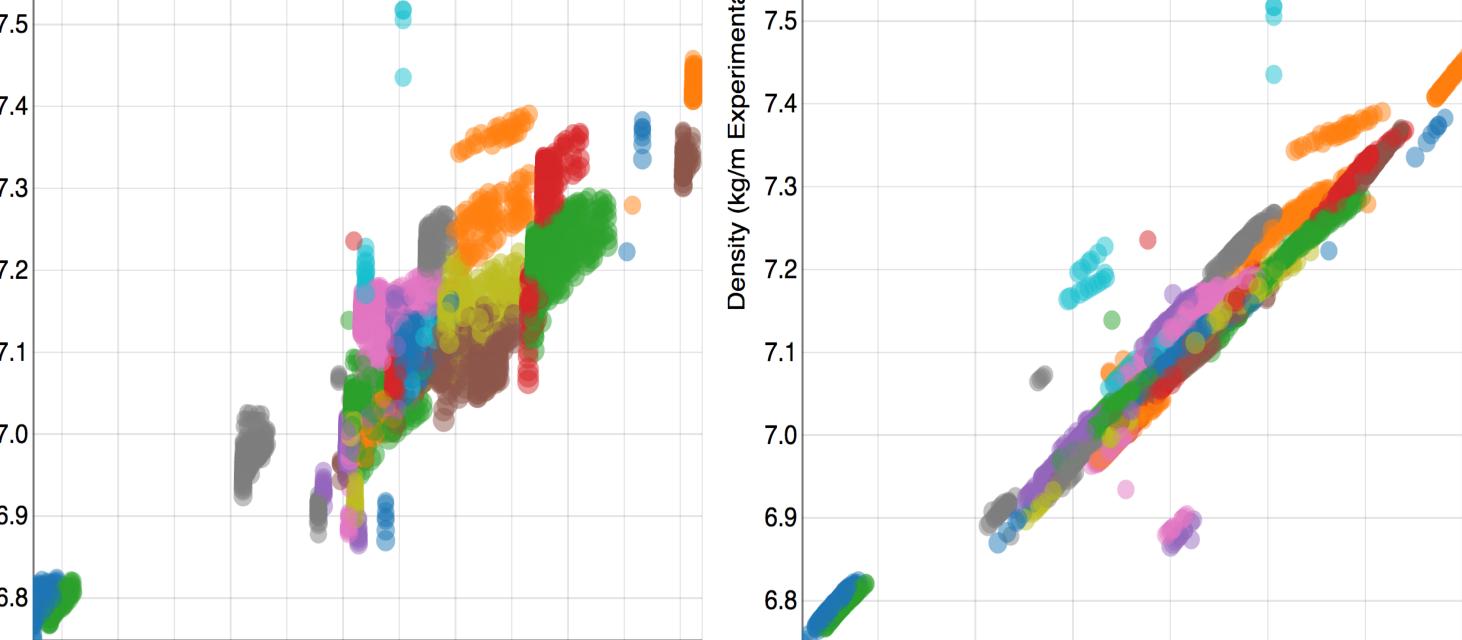


Figure 6: Webapp LASSO model with alpha of 0.8. Figure 7: Webapp LASSO model with alpha of 0.1.

Prediction for alpha of 0.8

Experim 7.4

Tooltips update as the value of alpha changes.

Results

Extra Details

Project webapp

is available for

The Salty

download

An included

allows users to

and watch the

LASSO model

slider bar

adjust the

change.

Prediction for alpha of 0.1

alpha value

through

GitHub.