

Evaluating Statistical Methods for Nuclear Forensics Analysis

Preliminary Examination

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Outline



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2 Literature Review

Nuclear Forensics

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Research Overview



How does the ability to determine forensic-relevant spent nuclear fuel attributes using machine learning techniques degrade as less information is available?

Determine

The inverse problem: given end measurements, calculate the model parameters that created them

Information

Nuclide vectors, measurements of isotope ratios

Forensic-relevant Attributes

Reactor type, enrichment, cooling time, burnup

Machine Learning Techniques

Creating statistical models (not physical)

Degrade

Model prediction performance

Less Information

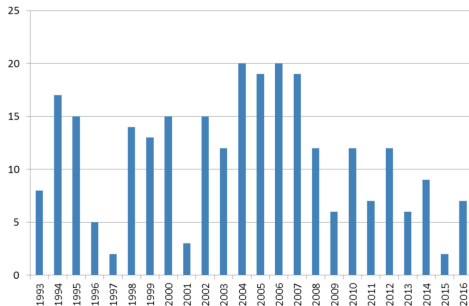
Error in nuclide vectors, fewer measurements, etc

Figure 1: Definitions of terms within the main research question



Nuclear Security and Forensics

Incidents related to trafficking or malicious use, 1993–2016



- FY2016 DHS DNDO budget : 0.3 bill
- FY2016 DOE NNSA nonpro budget : 1.6 bill

Figure 2: 24 years of incidents: HEU (12), Pu (2), Pu-Be neutron sources (4) [Obtained from: <https://www.iaea.org/sites/default/files/17/12/itdb-factsheet-2017.pdf>]



Needs in Nuclear Forensics

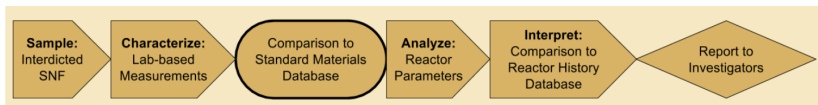


Figure 3: Typical technical nuclear forensics workflow

Material-specific:

- Measurement needs
- Measurement techniques
- Forensic signatures

Challenges:

- Rapid characterization
- Forensics databases
 - Multidimensional
 - Inconsistent uncertainties
 - International cooperation



Computational Methods

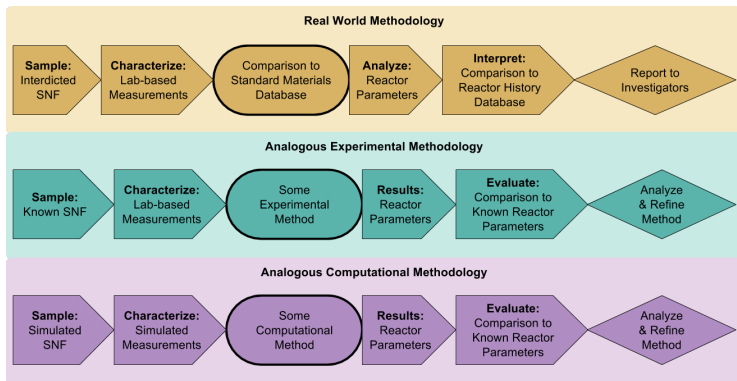


Figure 4: Nuclear forensics research: physical, experimental, and computational

Computational Methods

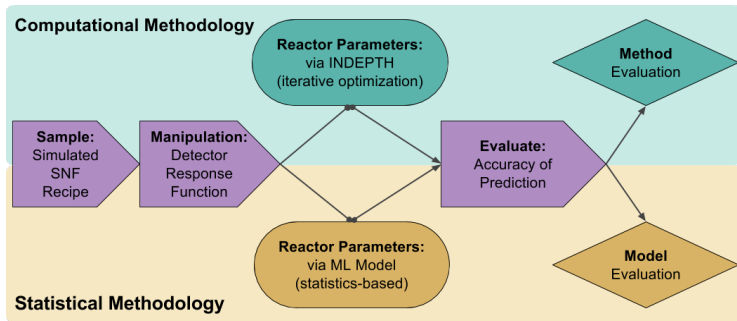


Figure 5: Comparison of two different computational approaches



Statistical Methods

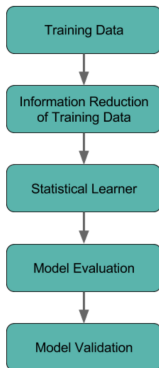


Figure 6: Workflow of a methodology using statistical models

- Training data: large set of SNF measurements
 - Labels (e.g., burnup)
 - Features (e.g., nuclide concs)
 - Instances (individual SNF recipe)
- Statistical learner
 - Machine learning algorithms
 - Algorithm parameters
 - Predict label of new instance
- Model evaluation
 - Diagnostic curves
 - Learning curves
 - Validation curves
 - Prediction error
 - Bias versus variance
 - Generalizability

Statistical Methods



	TRAINING DATA	TESTING DATA
Physical Motivation		
Ideal World	<i>Lab-Measured Mass Spectra</i>	<i>Lab-Measured Mass Spectra</i>
Real World	<i>Lab-Measured Gamma Spectra</i>	<i>Field-Measured Gamma Spectra</i>
Computational Representation		
Ideal World	<i>Simulation-Created Isotopics</i>	<i>Simulation-Created Isotopics</i>
Real World	<i>DRF-Derived Gamma Spectra</i>	<i>DRF-Derived Gamma Spectra</i>

Figure 7: Illustration of data set modularity



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Technical Nuclear Forensics



something showing illicit trafficking probs. SNF is really just for RDDs, but least guarded

Types of Investigations



Post-detonation

- Debris collection (fallout pred/dose rate)
- Rapid (field) analysis (isotope ratios, debris characterization)
- Data evaluation (uncertainty quantification)

Pre-detonation

- Material characterization (separations, etc)
- Material provenance (inverse prob)



Nuclear Forensics as an Inverse Problem

Necessary to determine the quality of prediction
Use Bayes' Framework:

$$P(A|B) = \frac{P(B|A)P(A)}{P(B)}$$

$$P(M|D) = \frac{P(D|M)P(M)}{P(D)}$$

Machine Learning



machine vs. statistical (domain knowledge-¿none)
supervised and unsupervised
clustering, dimensionality reduction
classification, regression – discrete and continuous variables

Supervised Regression

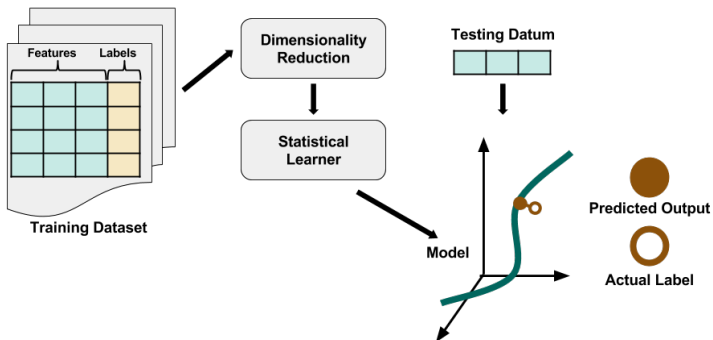


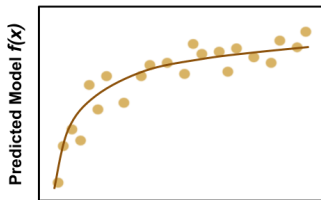
Figure 8: Schematic of a representative prediction workflow



Linear Models

Objective: minimize error over all training data wrt their labels

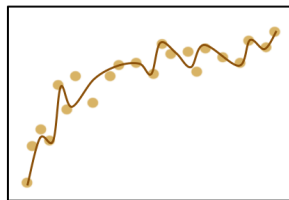
$$F(\mathbf{X}) = \beta_0 + \sum_{j=1}^p x_j \beta_j$$



Input Feature (x), $\lambda > 0$

Regularization using λ

$$F(\mathbf{X}) = \beta_0 + \sum_{j=1}^p x_j \beta_j + \lambda \sum_{j=1}^p \beta_j^2$$



Input Feature (x), $\lambda = 0$

Figure 9: How regularization might affect the generalizability of an ML model



Nearest Neighbor Methods

Objective: minimum distance between test sample and training instance(s)

$$Y(\mathbf{x}) = \frac{1}{k} \sum_{x_i \in N_k(\mathbf{x})} y_i$$

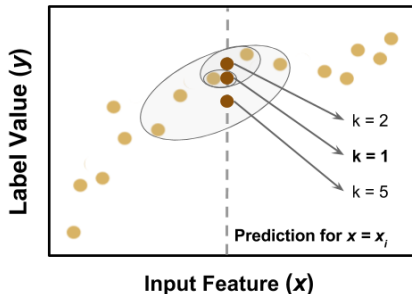


Figure 10: Illustration of the regularization effects by choosing k

Support Vector Machines

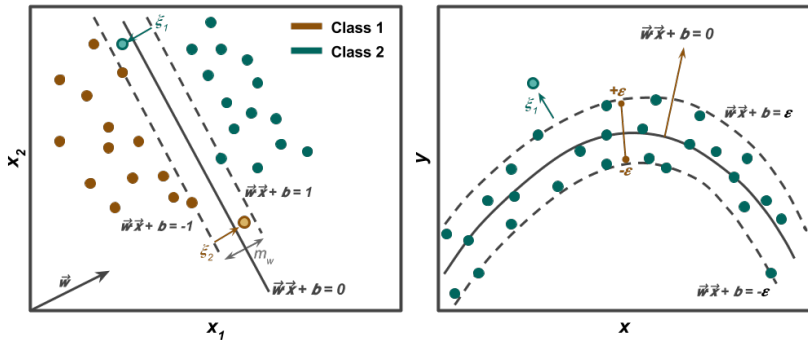


Figure 11: Classification with SVM and regression with SVR

Support Vector Regression with Many Dimensions



Objective: minimize margin width and outliers

$$\min \frac{1}{2} \|w\|^2 + C \sum_i \xi_i$$

$$\text{subject to : } |y_i - (w\phi(x_i) + b)| \leq \varepsilon + \xi_i$$

$$\text{where : } w = \sum_i \alpha_i y_i \phi(x_i)$$

$$\text{and : } K(x_i, x_j) = \phi(x_i)\phi(x_j) = e^{\gamma \|x_i - x_j\|^2}$$

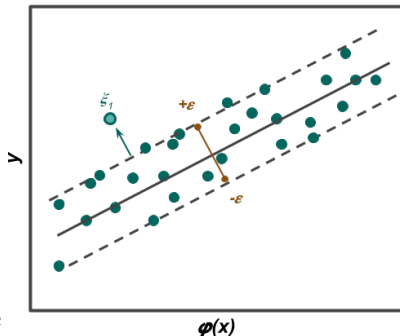


Figure 12: Diagram showing the use of the kernel trick with SVR

Dimensionality Reduction



Manual via domain knowledge or some measure

PCA

Factor Analysis

ICA

Sources of Error

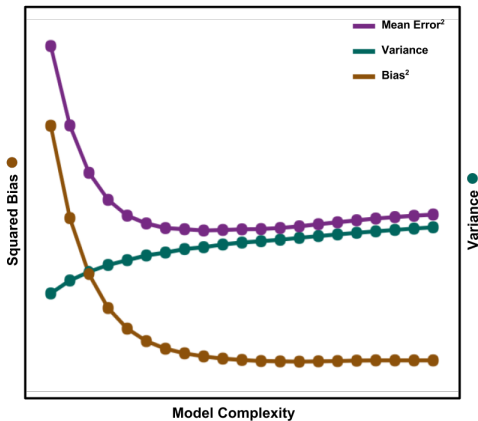


Figure 13: Bias and variance comprise the prediction error



Types of Error

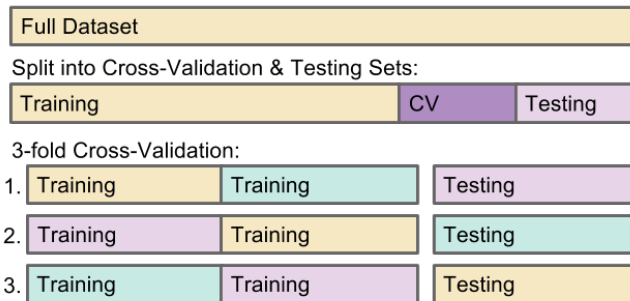


Figure 14: Diagram explaining the concept of k -fold cross-validation

Error Metrics



L1, L2: absolute error and squared error

Others: r^2 score, percent error

Used for model prediction error and optimization of algs in obj funcs



Training Set Size: Learning Curves

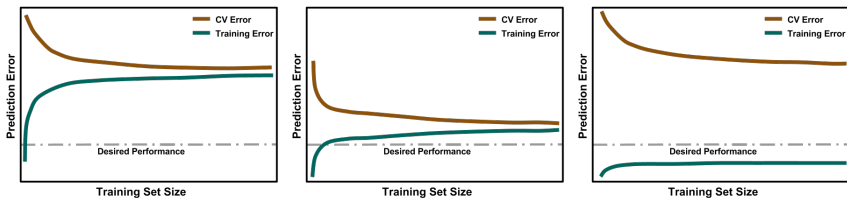


Figure 15: Learning curves for three training scenarios: high bias, balanced bias and variance, and high variance



Model Complexity: Validation Curves

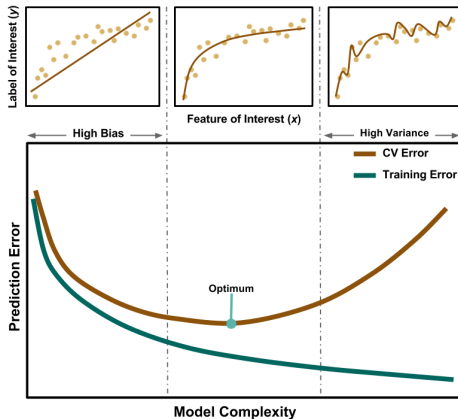


Figure 16: Validation curve showing different fitness of models



Model Comparison

Probabilities	Calculation Method and Example
P(D M) Prior	Given: M = BWR U-oxide with burnup = x GWd/MTU This is true: D = nuclide vector with Pu-239 = $y\%$ With: $z\%$ probability Calc'd from: ORIGEN simulations in training set
P(M) Likelihood	Given: No direct information on D This is true: M = BWR U-oxide with burnup = x GWd/MTU With: $z\%$ probability Calc'd from: Machine-learned model prediction
P(D) Marginal L.	Given: No direct information on M This is true: D = nuclide vector with Pu-239 = $y\%$ With: $z\%$ probability Calc'd from: Summation of training set instances
P(M D) Posterior	Given: D = nuclide vector with Pu-239 = $y\%$ This is true: M = BWR U-Ox with burnup = x GWd/MTU With: $z\%$ probability Calc'd from: All quantities above

$$\text{Posterior} = \frac{\text{Likelihood} * \text{Prior}}{\text{Marginal Likelihood}}$$

Table 1: Bayes

Computational Tools



cite stuff

- Training Data : SCALE/ORIGEN-ARP
- Statistics Toolkit : scikit-learn (python)
- Information Reduction
 - Gamma energies: ORIGEN
 - Computational gamma spectra: GADRAS

Pre-detonation Materials of Interest



UOC

UOX powder

SNF

Reprocessed SNF

Statistical Methods Employed



get images

Factor Analysis

SFCOMPO extension

Dayman paper on prediction ability wrt info reduction



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Proposed Experiment Methodology

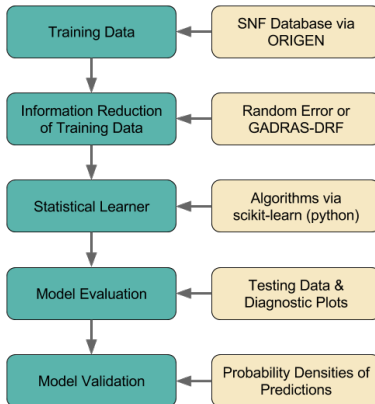


Figure 17: Workflow of the experiments with tools used for each step



Training Set

ORIGEN Rxtr	Rxtr Type	Enrichment
CE14x14	PWR	2.8
CE16x16	PWR	2.8
W14x14	PWR	2.8
W15x15	PWR	2.8
W17x17	PWR	2.8
S14x14	PWR	2.8
VVER440	PWR	3.60
VVER440_3.82	PWR	3.82
VVER440_4.25	PWR	4.25
VVER440_4.38	PWR	4.38
VVER1000	PWR	2.8
GE7x7-0	BWR	2.9
GE8x8-1	BWR	2.9
GE9x9-2	BWR	2.9
GE10x10-8	BWR	2.9
Abb8x8-1	BWR	2.9
Atrium9x9-9	BWR	2.9
SVEA64-1	BWR	2.9
SVEA100	BWR	2.9
CANDU28	PHWR	0.711
CANDU37	PHWR	0.711



Independent Testing Set

Reactor	Type	Enrichment	Cooling Time
CANDU28	PHWR	0.711	{1m, 7d, 30d, 1y}
CANDU28	PHWR	0.711	{3m, 9d, 2y}
CE16x16	PWR	2.8	{1m, 7d, 30d, 1y}
CE16x16	PWR	2.8	{3m, 9d, 2y}
CE16x16	PWR	3.1	{7d, 9d}
GE7x7-0	BWR	2.9	{1m, 7d, 30d, 1y}
GE7x7-0	BWR	2.9	{3m, 9d, 2y}
GE7x7-0	BWR	3.2	{7d, 9d}

Information Reduction



Random error here
gamma not implemented here



Algorithm Parameters

Algorithm	Parameter	Value
Nearest Neighbor Regression	n -neighbors	1
	Weights	uniform
	Distance Metric	L2: Euclidian Distance
Ridge Regression	Regularization, α	1.0
	Normalization	False
	Stopping Tolerance	0.001
Support Vector Regression	Kernel	Radial Basis Function
	Gamma, γ	0.001
	C	1000
	Epsilon, ϵ	0.1
	Stopping Tolerance	0.001

Table 5: caption



Initial Results

Algorithm	Error Origin	MAPE	RMSE [MWd/MTU]
Nearest Neighbor Regression	Testing Set	9.82	812.43
	5-fold Cross-Validation	2.24	421.41
Ridge Regression	Testing Set	15.68	1049.66
	5-fold Cross-Validation	0.08	13.08
Support Vector Regression	Testing Set	12.28	769.97
	5-fold Cross-Validation	2.08	188.07

Table 6: caption

ML Model Prediction with Reduced Information

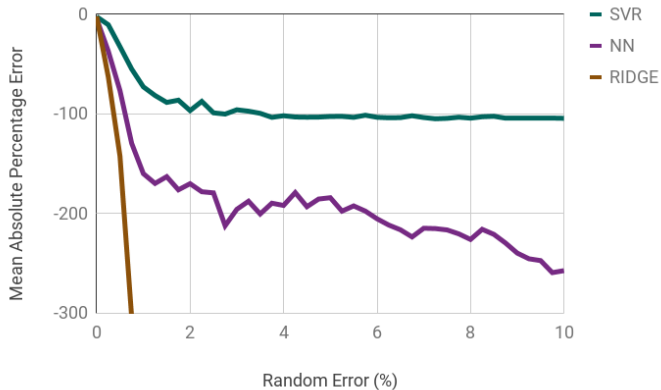
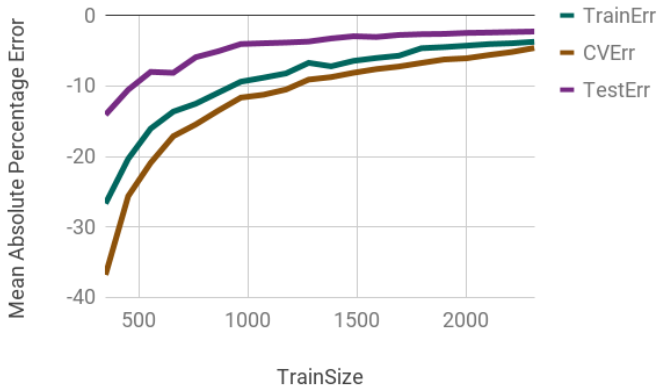


Figure 18: caption



SVR Learning Curve

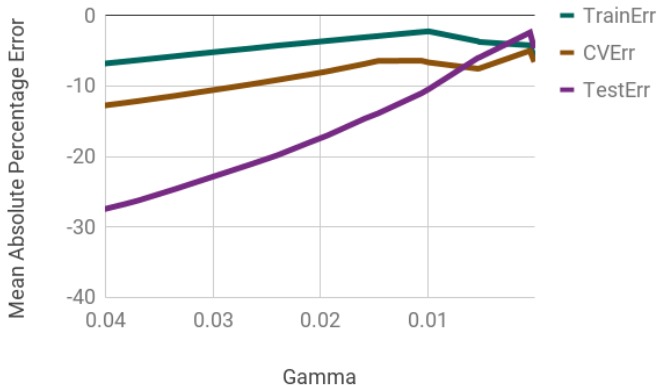
add in example LCs for comparison add in NN or Ridge LC (They look the same)





SVR Validation Curve

add in example VCs for comparison add in NN or Ridge VC (They look the same)



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Research Proposal Preparations



Previous Work - λ SFCOMPO-based
Finalizing set of algorithms
computational resources

Statistical Learning with Direct Isotopics



Goals : Understand limits of simplest scenario

- ① Usefulness of statistical methods for reactor parameter prediction
- ② Best performing methods

Variables

- ① the complexity of the ML algorithm used,
- ② feature reduction, and
- ③ different subsets of the decision space.



Statistical Learning with Direct Isotopics

Qualitative Hypotheses

- Complex algorithm will provide best behavior
- Manual preprocessing (feature reduction): speed, accuracy
- Reduction of decision space should help: PWR vs. BWR?

Risk Mitigation

- New algorithms: tree-based, neural nets, Bayesian MLE
- Statistical preprocessing: PCA, ICA
- New materials: Pu, UOC, Post-detonation (urban canyon [1])

Statistical Learning with Gamma Spectra



Goals : Understand limits of real-world scenario

- ① Level of reduction in reactor parameter prediction
- ② Best performing methods

Variables

- ① the complexity of the ML algorithm used,
- ② feature reduction (implicit), and
- ③ quality of training and/or testing data set.

Statistical Learning with Gamma Spectra



Qualitative Hypotheses

- Complex algorithm will provide best behavior
- Indirect isotopics = implicit feature reduction: less accurate
- Higher quality gamma spectra will yield better results

Risk Mitigation

- New algorithms: tree-based, neural nets, Bayesian MLE
- Further manual or statistical preprocessing
- Add isotope identification step

Statistical Learning with Reprocessed Fuel



Goals : Probe prediction performance in reprocessing scenario

- ① Experiment with both direct and indirect isotopics
- ② Fresh evaluation of preprocessing
- ③ Best performing methods for materials with multiple sources

Variables

- ① the complexity of the ML algorithm used,
- ② quality of training data set, and
- ③ type of preprocessing for feature reduction.

Statistical Learning with Reprocessed Fuel



Qualitative Hypotheses

- Complex algorithm will provide best behavior
- Reduced information will provide less accurate results
- ICA may outperform PCA, but factor analysis may outperform components analysis [4, 5, 6, 8, 7, 2, 3]

Risk Mitigation

- New algorithms: tree-based, neural nets, Bayesian MLE
- Manual preprocessing
- Results may be interesting even if prediction fails
- Ensemble methods or other creative solutions [8, 7]



Probability Distributions

Include uncertainty for measures of confidence, posterior probs become prob distrib

C : constant given by marginal likelihood

\mathbf{d} : training data set

\mathbf{m} : model parameters

$P(\mathbf{d}|\mathbf{m})$: likelihood distribution function

$P(\mathbf{m})$: prior probability distribution

$P(\mathbf{m}|\mathbf{d})$: posterior probability distribution

$$P(\mathbf{m}|\mathbf{d}) = C * P(\mathbf{d}|\mathbf{m}) * P(\mathbf{m})$$

Integrate over prob densities to get prob distrib

\mathbf{m} : range of predicted model parameters

\mathbf{d} is a set of nuclide vectors

$$\rho(\mathbf{x}) = \prod_i \rho(x_i)$$

$$P(\mathbf{m}) = \int_{\mathbf{m}} \rho(\mathbf{d}) d\mathbf{d}$$

Likelihood distribution function:

$$P(\mathbf{d}|\mathbf{m}) = \int_{\mathbf{d}, \mathbf{m}} \rho(\mathbf{d}|\mathbf{m}) d\mathbf{m}$$

But, we infer them...



Estimating Density Functions

estimate ρ , have a 'sense' or try different
prior probability distributions are given by the model space, e.g., reactor
parameters as predicted from the ML models. [?] Note: This implies the
posterior is now only dependent on the likelihood.

likelihood function: the training phase provides the maximum likelihood
distribution through the use of CV, since the results are reported as a
mean error with a standard deviation (which can be converted to
accuracy for likelihood) [?]

MLE is not this simple for other methods that do not employ CV [?, ?]



Posterior Odds

citations plz

calc a non-normalized posterior probability distribution, $P(m_i|d)$ then do it for a model obtained from a different algorithm, $P(m_j|d)$

relative posterior probability distribution : *posterior odds* $B_{ij} = \frac{\rho(d|m_i)}{\rho(d|m_j)}$: *Bayes factor*.

$$\frac{P(m_i|d)}{P(m_j|d)} = B_{ij} \frac{P(m_i)}{P(m_j)}$$

$ \ln B_{ij} $	Probability	Likelihood Strength
< 1.0	< 0.750	Inconclusive
1.0	0.750	Weak
2.5	0.923	Moderate
5.0	0.993	Strong

Table 9: Model comparison using likelihood strength

posterior probabilities calculated from $|\ln B_{ij}|$

Summarize:

Given a mean-squared error and its standard deviation from using CV with any alg, get MLE

compare two models : MLE to MLE



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Summarize



References I

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