

# Section 2.

# Stan Components

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**Part I**

**Stan Top Level**

# Stan's Namesake

- Stanislaw Ulam (1909–1984)
- Co-inventor of Monte Carlo method (and hydrogen bomb)



- Ulam holding the Fermiac, Enrico Fermi's physical Monte Carlo simulator for random neutron diffusion

# What is Stan?

- Stan is an **imperative** probabilistic programming language
  - cf., BUGS: declarative; Church: functional; Figaro: object-oriented
- Stan **program**
  - declares data and (constrained) parameter variables
  - defines log posterior (or penalized likelihood)
- Stan **inference**
  - MCMC for full Bayesian inference
  - VB for approximate Bayesian inference
  - MLE for penalized maximum likelihood estimation

# Platforms and Interfaces

- **Platforms:** Linux, Mac OS X, Windows
- **C++ API:** portable, standards compliant (C++03)
- **Interfaces**
  - **CmdStan:** Command-line or shell interface (direct executable)
  - **RStan:** R interface (Rcpp in memory)
  - **PyStan:** Python interface (Cython in memory)
  - **MatlabStan:** MATLAB interface (external process)
  - **Stan.jl:** Julia interface (external process)
  - **StataStan:** Stata interface (external process) [under testing]
- **Posterior Visualization & Exploration**
  - **ShinyStan:** Shiny (R) web-based

# Who's Using Stan?

- 1200 **users group** registrations; 10,000 manual **downloads** (2.5.0); 300 Google scholar citations (100+ fitting)
- **Biological sciences**: clinical drug trials, entomology, ophthalmology, neurology, genomics, agriculture, botany, fisheries, cancer biology, epidemiology, population ecology, neurology
- **Physical sciences**: astrophysics, molecular biology, oceanography, climatology
- **Social sciences**: population dynamics, psycholinguistics, social networks, political science
- **Other**: materials engineering, finance, actuarial, sports, public health, recommender systems, educational testing

# Documentation

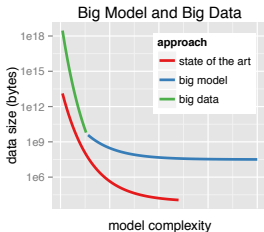
- *Stan User's Guide and Reference Manual*
  - 500+ pages
  - Example models, modeling and programming advice
  - Introduction to Bayesian and frequentist statistics
  - Complete language specification and execution guide
  - Descriptions of algorithms (NUTS, R-hat,  $n_{\text{eff}}$ )
  - Guide to built-in distributions and functions
- Installation and getting started manuals by interface
  - RStan, PyStan, CmdStan, MatlabStan, Stan.jl
  - RStan vignette

# Books and Model Sets

- **Model Sets** Translated to Stan
  - BUGS and JAGS examples (most of all 3 volumes)
  - Gelman and Hill (2009) *Data Analysis Using Regression and Multilevel/Hierarchical Models*
  - Wagenmakers and Lee (2014) *Bayesian Cognitive Modeling*
- **Books** with Sections on Stan
  - Gelman et al. (2013) *Bayesian Data Analysis*, 3rd Edition.
  - Kruschke (2014) *Doing Bayesian Data Analysis, Second Edition: A Tutorial with R, JAGS, and Stan*
  - Korner-Nievergelt et al. (2015) *Bayesian Data Analysis in Ecology Using Linear Models with R, BUGS, and Stan*

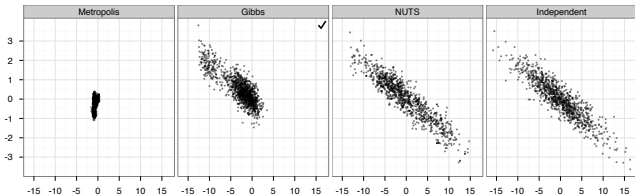


# Scaling and Evaluation



- Types of Scaling: data, parameters, **models**
- Time to converge and per effective sample size:  
0.5- $\infty$  times faster than BUGS & JAGS
- Memory usage: 1-10% of BUGS & JAGS

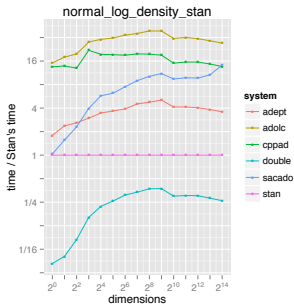
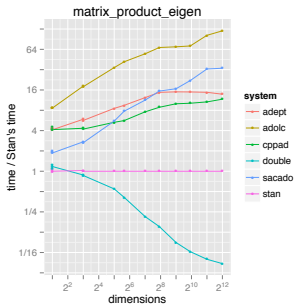
# NUTS vs. Gibbs and Metropolis



- Two dimensions of highly correlated 250-dim normal
- **1,000,000 draws** from Metropolis and Gibbs (thin to 1000)
- **1000 draws** from NUTS; 1000 independent draws

# Stan's Autodiff vs. Alternatives

- Among C++ open-source offerings: Stan is **fastest** (for gradients), **most general** (functions supported), and **most easily extensible** (simple OO)



**Part II**

# **Stan Language**

# Stan is a Programming Language

- **Not** a graphical specification language like BUGS or JAGS
- Stan is a Turing-complete imperative programming language for specifying differentiable log densities
  - reassignable local variables and scoping
  - full conditionals and loops
  - functions (including recursion)
- With automatic “black-box” inference on top (though even that is tunable)
- Programs computing same thing may have different efficiency

# Basic Program Blocks

- **data** (once)
  - *content*: declare data types, sizes, and constraints
  - *execute*: read from data source, validate constraints
- **parameters** (every log prob eval)
  - *content*: declare parameter types, sizes, and constraints
  - *execute*: transform to constrained, Jacobian
- **model** (every log prob eval)
  - *content*: statements defining posterior density
  - *execute*: execute statements

# Derived Variable Blocks

- **transformed data** (once after data)
  - *content*: declare and define transformed data variables
  - *execute*: execute definition statements, validate constraints
- **transformed parameters** (every log prob eval)
  - *content*: declare and define transformed parameter vars
  - *execute*: execute definition statements, validate constraints
- **generated quantities** (once per draw, double type)
  - *content*: declare and define generated quantity variables;  
includes pseudo-random number generators  
(for posterior predictions, event probabilities, decision making)
  - *execute*: execute definition statements, validate constraints

# Variable and Expression Types

Variables and expressions are **strongly, statically typed**.

- **Primitive:** `int`, `real`
- **Matrix:** `matrix[M,N]`, `vector[M]`, `row_vector[N]`
- **Bounded:** primitive or matrix, with  
`<lower=L>`, `<upper=U>`, `<lower=L,upper=U>`
- **Constrained Vectors:** `simplex[K]`, `ordered[N]`,  
`positive_ordered[N]`, `unit_length[N]`
- **Constrained Matrices:** `cov_matrix[K]`, `corr_matrix[K]`,  
`cholesky_factor_cov[M,N]`, `cholesky_factor_corr[K]`
- **Arrays:** of any type (and dimensionality)



# Integers vs. Reals

- Different types (conflated in BUGS, JAGS, and R)
- Distributions and assignments care
- Integers may be assigned to reals but not vice-versa
- Reals have not-a-number, and positive and negative infinity
- Integers single-precision up to +/- 2 billion
- Integer division rounds (Stan provides warning)
- Real arithmetic is inexact and reals should not be (usually) compared with `==`

# Arrays vs. Matrices

- Stan separates arrays, matrices, vectors, row vectors
- Which to use?
- Arrays allow most efficient access (no copying)
- Arrays stored first-index major (i.e., 2D are row major)
- Vectors and matrices required for matrix and linear algebra functions
- Matrices stored column-major
- Are not assignable to each other, but there are conversion functions

# Logical Operators

<i>Op.</i>	<i>Prec.</i>	<i>Assoc.</i>	<i>Placement</i>	<i>Description</i>
	9	left	binary infix	logical or
&&	8	left	binary infix	logical and
==	7	left	binary infix	equality
!=	7	left	binary infix	inequality
<	6	left	binary infix	less than
<=	6	left	binary infix	less than or equal
>	6	left	binary infix	greater than
>=	6	left	binary infix	greater than or equal

# Arithmetic and Matrix Operators

<i>Op.</i>	<i>Prec.</i>	<i>Assoc.</i>	<i>Placement</i>	<i>Description</i>
+	5	left	binary infix	addition
-	5	left	binary infix	subtraction
*	4	left	binary infix	multiplication
/	4	left	binary infix	(right) division
\	3	left	binary infix	left division
.*	2	left	binary infix	elementwise multiplication
./	2	left	binary infix	elementwise division
!	1	n/a	unary prefix	logical negation
-	1	n/a	unary prefix	negation
+	1	n/a	unary prefix	promotion (no-op in Stan)
^	2	right	binary infix	exponentiation
'	0	n/a	unary postfix	transposition
()	0	n/a	prefix, wrap	function application
[]	0	left	prefix, wrap	array, matrix indexing

# Built-in Math Functions

- All built-in **C++ functions and operators**  
C math, TR1, C++11, including all trig, pow, and special log1 m, erf, erfc, fma, atan2, etc.
- Extensive library of **statistical functions**  
e.g., softmax, log gamma and digamma functions, beta functions, Bessel functions of first and second kind, etc.
- Efficient, arithmetically stable **compound functions**  
e.g., multiply log, log sum of exponentials, log inverse logit

# Built-in Matrix Functions

- **Basic arithmetic:** all arithmetic operators
- **Elementwise arithmetic:** vectorized operations
- **Solvers:** matrix division, (log) determinant, inverse
- **Decompositions:** QR, Eigenvalues and Eigenvectors, Cholesky factorization, singular value decomposition
- **Compound Operations:** quadratic forms, variance scaling, etc.
- **Ordering, Slicing, Broadcasting:** sort, rank, block, rep
- **Reductions:** sum, product, norms
- **Specializations:** triangular, positive-definite,

# User-Defined Functions

- **functions** (compiled with model)
  - *content*: declare and define general (recursive) functions (use them elsewhere in program)
  - *execute*: compile with model
  
- Example

```
functions {  
  
    real relative_difference(real u, real v) {  
        return 2 * fabs(u - v) / (fabs(u) + fabs(v));  
    }  
  
}
```

# Differential Equation Solver

- System expressed as function
  - given state ( $y$ ) time ( $t$ ), parameters ( $\theta$ ), and data ( $x$ )
  - return derivatives ( $\partial y / \partial t$ ) of state w.r.t. time
- Simple harmonic oscillator diff eq

```
real[] sho(real t,          // time
           real[] y,       // system state
           real[] theta,   // params
           real[] x_r,     // real data
           int[] x_i) {    // int data
  real dydt[2];
  dydt[1] <- y[2];
  dydt[2] <- -y[1] - theta[1] * y[2];
  return dydt;
}
```



# Differential Equation Solver

- Solution via functional, given initial state ( $y_0$ ), initial time ( $t_0$ ), desired solution times ( $t_s$ )

```
mu_y <- integrate_ode(sho, y0, t0, ts, theta, x_r, x_i);
```

- Use noisy measurements of  $y$  to estimate  $\theta$

```
y ~ normal(mu_y, sigma);
```

- Pharmacokinetics/pharmacodynamics (PK/PD),
- soil carbon respiration with biomass input and breakdown

# Diff Eq Derivatives

- Need derivatives of solution w.r.t. parameters
- Couple derivatives of system w.r.t. parameters

$$\left( \frac{\partial}{\partial t} y, \frac{\partial}{\partial t} \frac{\partial y}{\partial \theta} \right)$$

- Calculate coupled system via nested autodiff of second term

$$\frac{\partial}{\partial \theta} \frac{\partial y}{\partial t}$$

# Distribution Library

- Each distribution has
  - log density or mass function
  - cumulative distribution functions, plus complementary versions, plus log scale
  - Pseudo-random number generators
- Alternative parameterizations  
(e.g., Cholesky-based multi-normal, log-scale Poisson, logit-scale Bernoulli)
- New multivariate correlation matrix density: LKJ  
degrees of freedom controls shrinkage to (expansion from) unit matrix

# Statements

- **Sampling:** `y ~ normal(mu, sigma)` (increments log probability)
- **Log probability:** `increment_log_prob(lp);`
- **Assignment:** `y_hat <- x * beta;`
- **For loop:** `for (n in 1:N) ...`
- **While loop:** `while (cond) ...`
- **Conditional:** `if (cond) ...; else if (cond) ...; else ...;`
- **Block:** `{ ... }` (allows local variables)
- **Print:** `print("theta=", theta);`
- **Reject:** `reject("arg to foo must be positive, found y=", y);`

# “Sampling” Increments Log Prob

- A Stan program defines a log posterior
  - typically through log joint and Bayes’s rule
- Sampling statements are just “syntactic sugar”
- A shorthand for incrementing the log posterior
- The following define the same\* posterior
  - `y ~ poisson(lambda);`
  - `increment_log_prob(poisson_log(y, lambda));`
- \* up to a constant
- Sampling statement drops constant terms

# Local Variable Scope Blocks

- `a ~ bernoulli(theta);`

is more efficient with sufficient statistics

```
{  
  real a; // local variable  
  a <- 0;  
  for (n in 1:N)  
    a <- a + x[n]; // reassignment  
  a ~ binomial(N, theta);  
}
```

- Simpler, but roughly same efficiency:

```
sum(a) ~ binomial(N, theta);
```

# Print and Reject

- Print statements are for **debugging**
  - printed every log prob evaluation
  - print values in the middle of programs
  - check when log density becomes undefined
  - can embed in conditionals
- Reject statements are for **error checking**
  - typically function argument checks
  - cause a rejection of current state (0 density)

# Prob Function Vectorization

- Stan's probability functions are vectorized for speed
  - removes repeated computations (e.g.,  $-\log \sigma$  in normal)
  - reduces size of expression graph for differentiation
- Consider:  $y \sim \text{normal}(\mu, \sigma)$ ;
- Each of  $y$ ,  $\mu$ , and  $\sigma$  may be any of
  - scalars (integer or real)
  - vectors (row or column)
  - 1D arrays
- All dimensions must be scalars or having matching sizes
- Scalars are broadcast (repeated)



**Part III**

# **Transformed Parameters**

# Transforms: Precision

```
parameters {  
  real<lower=0> tau; // precision  
  ...  
}  
transformed parameters {  
  real<lower=0> sigma; // sd  
  sigma <- 1 / sqrt(tau);  
}
```

# Transforms: “Matt Trick”

```
parameters {  
  vector[K] beta_raw; // non-centered  
  real mu;  
  real<lower=0> sigma;  
}  
transformed parameters {  
  vector[K] beta; // centered  
  beta <- mu + sigma * beta_raw;  
}  
model {  
  mu ~ cauchy(0, 2.5);  
  sigma ~ cauchy(0, 2.5);  
  beta_raw ~ normal(0, 1);  
}
```

**Part IV**

# **Generated Quantities for Prediction**

# Linear Regression (Normal Noise)

- **Likelihood:**

- $y_n = \alpha + \beta x_n + \epsilon_n$

- $\epsilon_n \sim \text{Normal}(0, \sigma)$

for  $n \in 1:N$

- Equivalently,

- $y_n \sim \text{Normal}(\alpha + \beta x_n, \sigma)$

- **Priors** (improper)

- $\sigma \sim \text{Uniform}(0, \infty)$

- $\alpha, \beta \sim \text{Uniform}(-\infty, \infty)$

- Stan allows **improper prior**; requires **proper posterior**.

# Linear Regression in Stan

```
data {  
  int<lower=0> N;  
  vector[N] x;  
  vector[N] y;  
}  
parameters {  
  real alpha;  
  real beta;  
  real<lower=0> sigma;  
}  
model {  
  y ~ normal(alpha + beta * x, sigma);  
}  
  
// for (n in 1:N)  
//   y[n] ~ normal(alpha + beta * x[n], sigma);
```

# Posterior Predictive Inference

- Parameters  $\theta$ , observed data  $y$  and data to predict  $\tilde{y}$

$$p(\tilde{y}|y) = \int_{\Theta} p(\tilde{y}|\theta) p(\theta|y) d\theta$$

- data {  
 int<lower=0> N\_tilde;  
 matrix[N\_tilde,K] x\_tilde;  
 ...  
parameters {  
 vector[N\_tilde] y\_tilde;  
 ...  
model {  
 y\_tilde ~ normal(x\_tilde \* beta, sigma);

# Predict w. Generated Quantities

- Replace sampling with pseudo-random number generation

```
generated quantities {  
  vector[N_tilde] y_tilde;  
  
  for (n in 1:N_tilde)  
    y_tilde[n] <- normal_rng(x_tilde[n] * beta, sigma);  
}
```

- Must include noise for predictive uncertainty
- PRNGs only allowed in generated quantities block
  - more computationally efficient per iteration
  - more statistically efficient with i.i.d. samples (i.e., MC, not MCMC)



**End (Section 2)**