

Fazang: A Reverse-mode Automatic differentiation tool in Fortran

**User's Guide
(Version pre-release)**

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February 16, 2022

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CHAPTER 1

Introduction

Fazang is a reverse-mode automatic differentiation (AD) tool. The project is heavily influenced by **Stan/Math** [1], a project the author is also involved in. **Fazang** is intended to support general scientific computing in Fortran beyond Bayesian inference and Markov Chain Monte Carlo that **Stan/Math** is designed for.

User should be aware that the project is at early stage and still under development. For any questions, suggestions, and contributions, please visit the project at <https://github.com/yizhang-yiz/fazang>.

CHAPTER 2

Quick Start

Currently **Fazang** has been tested on Linux and MacOS platform, with Fortran compiler Intel Fortran 19.0.1+ and GNU Fortran 11.2.0+.

After downloading **Fazang**, user can use **meson** to build the library.

```
git clone git@github.com:yizhang-yiz/fazang.git
cd fazang && mkdir build && cd build
meson compile
```

This generates a shared library at **build/src/**. User needs to link this library when building an application. This can be done in **meson** by setting

```
executable('app_name', files('path/to/app_file.F90'), dependencies : fazang_dep)
```

Fazang provides a user-facing derived type **var**. This is the type for the dependent and independent variables of which the adjoint (derivative) will be calculated.

For example, consider the log of the Gaussian distribution density with mean μ and standard deviation σ

$$f(\mu, \sigma) = \log \left(\frac{1}{\sigma \sqrt{2\pi}} \exp \left(-\frac{1}{2} \left(\frac{y - \mu}{\sigma} \right)^2 \right) \right) \quad (2.1)$$

The following programe calculates $\frac{df}{d\mu}$ and $\frac{df}{d\sigma}$ at $y = 1.3$, $\mu = 0.5$, and $\sigma = 1.2$.

```
program log_demo
  use fazang ! load Fazang library

  implicit none

  real(rk) :: y
  type(var) :: f, sigma, mu

  ! data
  y = 1.3d0

  ! independent variables
  mu = var(0.5d0)
  sigma = var(1.2d0)

  ! dependent
  f = var(-0.5d0 * log(2 * pi))
```

```
f = f - log(sigma)
f = f - 0.5d0 * ((y - mu) / sigma) ** 2.d0;

! use grad() to calculate df/d(mu) and df/d(sigma). Each var's
! derivative (also called adjoint) can be access through var%adj().

call f%grad()
write(*, *) "df/d(mu): ", mu%adj()
write(*, *) "df/d(sigma): ", sigma%adj()
end program log_demo
```

CHAPTER 3

Use Fazang

Fazang uses `var` type to record numerical and gradient operations. The type supports three functions

- `var%val()` : returns value
- `var%adj()` : returns derivative, henceforth referred as *adjoint*.
- `var%grad()` : takes gradient operation with respect to the current `var` variable.

1. Constructors

`var` can be constructed using overloaded `var` interface.

```
real(real64) :: a, b(3), c(2, 3)
real(real64) :: new_a, new_b(3), new_c(2, 3)
type(var) :: x, y(3), z(2, 3)
! ...
x = var()           ! x%val() == 0.d0
x = var(a)          ! x%val() == a
y = var(b)          ! y%val() == b
z = var(c)          ! z%val() == c
```

2. Assignment

`var` can be assigned from consistent `var` and `real(real64)`.

```
! ....
x = new_a           ! x%val() == new_a
y = new_b           ! y%val() == new_b
z = new_c           ! z%val() == new_c
```

3. Gradient

Consider a variable z calculated by the composition of a series of operations

$$z = f_1(z_1), \quad z_1 = f_2(z_2), \quad \dots, \quad z_{n-1} = f_n(z_n).$$

For $z_i, i = 1, \dots, n$ we refer dz/dz_i as the *adjoint* of z_i , denoted by z_i^{adj} . The chain rule says the adjoints can be calculated recursively [2],

$$z^{\text{adj}} = 1, \quad z_1^{\text{adj}} = z^{\text{adj}} \frac{df_1}{dz_1}, \quad \dots, \quad z_i^{\text{adj}} = z_{i-1}^{\text{adj}} \frac{df_i}{dz_i}.$$

We often refer each (f_i, z_i) pair as a *node*, and z_i the *operand* of operation f_i . The above recursion through the nodes requires a way to store and visit the *callstack* of nodes. It is embodied in Fazang by the `var%grad()` function. When `z%grad()` is called, `z`'s adjoint is

set to 1, and every other `var` variable is transversed with its adjoint updated. In order to calculate the adjoint with respect to another variable, user must call `set_zero_all_adj()` first to reset all adjoints to zero.

An alternative to invoke gradient calculation is to define the dependent as a function and feed it to Fazang's `gradient` function. Take Eq.(2.1) for example, we can first define the function for $f(\mu, \sigma)$.

```
module func
  use fazang ! load Fazang library
  implicit none

  real(rk), parameter :: y = 1.3d0

contains
  type(var) function f(x)
    type(var), intent(in) :: x(:)
    type(var) :: mu, sigma
    mu = x(1)
    sigma = x(2)
    f = -0.5d0 * log(2 * pi) - log(sigma) - 0.5d0 * ((y - mu) / sigma) ** 2.d0;
  end function f
end module func
```

Then we can supply function `f` as a procedure argument.

```
program log_demo2
  use iso_c_binding
  use fazang
  use func

  implicit none

  real(real64) :: fx(3), x(2)
  x = [0.5d0, 1.2d0]

  fx = gradient(f, x)
  write(*, *) "f(x): ", fx(1)
  write(*, *) "df/d(x(1)):", fx(2)
  write(*, *) "df/d(x(2)):", fx(3)
end program log_demo2
```

The output of `gradient(f, x)` is an array of size `1 + size(x)`, with first component being the function value, and the rest the partial derivatives.

Note that the above approach of using `gradient` function does not involve explicitly setting up `var` variables. Fazang achieves this by using a *nested* AD environment.

4. Nested AD environment

Let us take a look of the internals of Fazang's `gradient` function. The `dependent_function` interface requires f to follow the above example's signature, and `x` is the `real64` array of

independent variables. We then create the `var` version of `x` and introduce it to `f`. The evaluation result is saved in `f_var` variable. The adjoints are obtained by calling `f_var%grad()`. Unlike what we have seen, the above process happens within a pairing `begin_nested()` and `end_nested()` calls.

```
function gradient(f, x) result (f_df)
  procedure(dependent_function) :: f
  real(real64), intent(in) :: x(:)
  real(real64) :: f_df(1 + size(x))
  type(var) :: x_var(size(x)), f_var

  call begin_nested()

  x_var = var(x)
  f_var = f(x_var)
  f_df(1) = f_var%val()
  call f_var%grad()
  f_df(2:(1+size(x))) = x_var%adj()

  call end_nested()
end function gradient
```

When we use these two functions, all the `var` variables created in between are "temporary", in the sense that the values and adjoints of these variables are no longer available after `call end_nested()`. User can use this function pair to construct a local gradient evaluation procedure.

5. Jacobian

Similar to `gradient`, using the same nested technique `Fazang` provides a `jacobian` function that calculates the Jacobian matrix of `f`, a multivariate function $f : \mathbb{R}^m \rightarrow \mathbb{R}^n$ for an input array `x` of dimension `m`.

```
function jacobian(f, n, x) result (f_df)
```

The input function must follow the interface

```
abstract interface
  function jac_dependent_function (x, n) result (fx)
    import :: var
    integer, intent(in) :: n
    type(var), intent(in) :: x(:)
    type(var) :: fx(n)
  end function jac_dependent_function
end interface
```

where `n` is the output dimension. Like `gradient`, the output `f_df` has dimension $n \times (m+1)$, with the first column being the function results and the rest columns the adjoints.

6. Functions

Numeric functions supported by **Fazang** are listed in Appendix A. All unary and binary functions are **elemental**. The binary functions allow mixed argument types, namely, either argument can be **real64** type while the other the **var** type.

Probability distributions supported by **Fazang** are list in Appendix B.

7. Ordinary differential equations

Fazang supports ODE solutions through CVODES from SUNDIALS library [3]. One can solve ODE like this.

```
! user defined ODE
module ode_mod
  use fazang
  use, intrinsic :: iso_c_binding
  implicit none

  real(rk), parameter :: params(2) = [0.2d0, 0.1d0]

contains
  ! user defined right-hand-side
  subroutine eval_rhs(t, y, fy)
    real(c_double), intent(in) :: t, y(:)
    real(c_double), intent(inout) :: fy(size(y))
    fy(1) = y(2)
    fy(2) = t * y(1) * sum(params%val())
  end subroutine eval_rhs
end module ode_mod

program cvodes_solve_data
  use ode_mod
  use fazang
  implicit none

  real(rk) :: yt(2, 3)
  real(rk) :: y0(2)
  type(cvodes_tol) :: tol

  ! import ODE
  ! import Fazang

  ! output array
  ! initial condition
  ! basic solver control

  ! use BDF method with given relative tolerance, absolute tolerance,
  ! and max number of steps between outputs
  tol = cvodes_tol(CV_BDF, 1.d-10, 1.d-10, 1000_8)

  ! initial condition
  y0 = [1.2d0, 1.8d0]

  ! solve the ODE with initial time 0.d0 and
  ! output time 1.d0, 2.d0, 3.d0
  yt = cvodes_sol(0.d0, y0, [1.d0, 2.d0, 3.d0], eval_rhs, tol)

end program cvodes_solve_data
```

In the above example, we first define an ODE following `Fazang`'s interface on the RHS. The defined RHS function `eval_rhs` will be later used as an argument to `cvodes_sol`. In addition to initial condition, one must also define an object for solver control. Such an object must be of a type that `extend` the `cvodes_options` abstract type. Here we use `Fazang`'s basic type `cvodes_tol`, which gives: the integration scheme (`CV_BDF` for BDF method or `CV_ADAMS` for Adams-Moulton method), relative tolerance, absolute tolerance, and the maximum number of steps allowed between output.

Then to the solver interface `cvodes_sol` we give the initial time, initial condition, array for output time, the RHS subroutine, and the solver control. It returns a 2D array with each column at a requested output time.

7.1. Forward sensitivity. Combining the sensitivity capability of `CVODES` and `AD` from `Fazang`, we can solve for ODE sensitivity with respect to given parameters without explicitly supplying Jacobian. For that the user-defined ODE must include an additional RHS definition with `var` parameters, following `Fazang`'s RHS interface.

```

module ode_mod
  use fazang
  use, intrinsic :: iso_c_binding
  implicit none

  real(rk), parameter :: omega = 0.5d0
  real(rk), parameter :: d1 = 1.0d0
  real(rk), parameter :: d2 = 1.0d0

contains
  ! right-hand-side for data input
  subroutine eval_rhs(t, y, fy)
    implicit none
    real(c_double), intent(in) :: t, y(:)
    real(c_double), intent(inout) :: fy(size(y))
    fy(1) = y(2)
    fy(2) = sin(omega * d1 * d2 * t)
  end subroutine eval_rhs

  ! right-hand-side for var input with parameters
  ! y, p, and output fy must all be of var type
  subroutine eval_rhs_pvar(t, y, fy, p)
    implicit none
    real(c_double), intent(in) :: t
    type(var), intent(in) :: y(:), p(:)
    type(var), intent(inout) :: fy(size(y))
    fy(1) = y(2)
    fy(2) = sin(p(1) * p(2) * p(3) * t)
  end subroutine eval_rhs_pvar
end module ode_mod

```

Now we can solve the defined ODE in a similar way.

```

program cvodes_demo
  use ode_mod
  use fazang
  implicit none

  type(var) :: yt(2, 3)
  type(cvodes_tol) :: tol
  real(rk), parameter :: ts(3) = [1.2d0, 2.4d0, 4.8d0]
  real(rk), parameter :: y00(2) = [0.2d0, 0.8d0]
  type(var) :: param(3)
  real(rk) :: y0(2), ga(2)
  integer :: i, j

  y0 = y00                                ! init condition
  param = var([omega, d1, d2])             ! parameters
  tol = cvodes_tol(CV_BDF, 1.d-10, 1.d-10, 1000_8)

  yt = cvodes_sol(0.d0, y0, ts, param, eval_rhs,&
    & eval_rhs_pvar, tol)
  ! ...
end program cvodes_demo

```

Note that now the call to `cvodes_sol` includes additional argument `param` as the sensitivity parameters, as well as the RHS function for `var` inputs. When the sensitivities are desired, one can take the gradient of the result in the same way.

```

call yt(1, 1) % grad()
write(*, *) "dy_1/ d_omega at time ts(1):", param(1)%adj()

```

7.2. Functions.

(1) Data solution

```

function cvodes_sol(t, y, ts, rhs, cvs_options) result(yt)
  real(real64), intent(in) :: t           ! initial time
  real(real64), intent(inout) :: y(:)     ! initial condition
  real(real64), intent(in) :: ts(:)       ! output time
  procedure(cvs_rhs_func) :: rhs          ! RHS definition (see below)
  class(cvodes_options), intent(in) :: cvs_options ! solver control
  real(real64) :: yt(size(y), size(ts)) ! solution

```

(2) Sensitivity solution with respect to the initial condition

```

function cvodes_sol(t, y, ts, rhs, rhs_yvar, cvs_options) result(yt)
  real(real64), intent(in) :: t           ! initial time
  type(var), intent(inout) :: y(:)       ! initial condition
  real(real64), intent(in) :: ts(:)       ! output time
  procedure(cvs_rhs_func) :: rhs          ! data-only RHS (see below)
  procedure(cvs_rhs_func_yvar) :: rhs_yvar ! var-type RHS (see below)

```

```

class(cvodes_options), intent(in) :: cvs_options ! solver control
type(var) :: yt(size(y), size(ts)) ! solution

```

- (3) Sensitivity solution with respect to the parameters

```

function cvodes_sol(t, y, ts, param, rhs, rhs_pvar, cvs_options) result(yt)
  real(real64), intent(in) :: t ! initial time
  real(real64), intent(inout) :: y(:) ! initial condition
  real(real64), intent(in) :: ts(:) ! output time
  type(var), target, intent(in) :: param(:) ! parameters
  procedure(cvodes_rhs_func) :: rhs ! data-only RHS (see below)
  procedure(cvodes_rhs_func_pvar) :: rhs_pvar ! var-type RHS (see below)
  class(cvodes_options), intent(in) :: cvs_options ! solver control
  type(var) :: yt(size(y), size(ts)) ! solution

```

- (4) Interfaces for different solvers

```

abstract interface
  subroutine cvs_rhs_func(t, y, fy)
    import c_double
    real(c_double), intent(in) :: t, y(:)
    real(c_double), intent(inout) :: fy(size(y))
  end subroutine cvs_rhs_func

  subroutine cvs_rhs_func_yvar(t, y, f)
    import c_double, var
    real(c_double), intent(in) :: t
    type(var), intent(in) :: y(:)
    type(var), intent(inout) :: f(size(y))
  end subroutine cvs_rhs_func_yvar

  subroutine cvs_rhs_func_pvar(t, y, f, p)
    import c_double, var
    real(c_double), intent(in) :: t
    type(var), intent(in) :: y(:), p(:)
    type(var), intent(inout) :: f(size(y))
  end subroutine cvs_rhs_func_pvar
end interface

```

- (5) Solver controls The last argument of the solver call is a solver control object. User-defined type must be able to follow CVODES user guide to modify CVODES memory object, by extending the abstract type `cvodes_options`.

```

type, abstract :: cvodes_options
  integer :: cv_method = -1
  contains
    procedure(set_cvodes), deferred :: set
end type cvodes_options

abstract interface

```

```

subroutine set_cvodes(this, mem)
  import c_ptr, cvodes_options
  class(cvodes_options), intent(in) :: this
  type(c_ptr), intent(inout) :: mem ! CVODES memory
end subroutine set_cvodes
end interface

```

One can follow Fazang 's tolerance control type as an example.

```

type, extends(cvodes_options) :: cvodes_tol
  real(c_double) :: rtol, atol
  integer(c_long) :: max_nstep
contains
  procedure :: set
end type cvodes_tol

contains

subroutine set(this, mem)
  class(cvodes_tol), intent(in) :: this
  type(c_ptr), intent(inout) :: mem ! CVODES memory
  integer :: ierr
! call cvodes functions
  ierr = FCVodeStolerances(mem, this % rtol, this % atol)
  ierr = FCVodeSetMaxNumSteps(mem, this % max_nstep)
end subroutine set

```

CHAPTER 4

Design

The core of any reverse-mode automatic differentiation is the data structure to store and visit the callstack. **Fazang** achieves this through two derived types, **tape** and **vari**.

1. tape data structure

A **tape** is an **int32** array emulating a stack, with an integer marker **head** pointing to the head to the current stack top.

```
type :: tape
  integer(ik) :: head = 1
  integer(ik), allocatable :: storage(:)
  !...
```

Each time a new AD node is created, space in **storage** is allotted to store the node's

- value $f_i(z_i)$,
- adjoint z_{i-1}^{adj} ,
- number of **var** operands of f_i ,
- The **var** operands' index in the same **tape** array,
- number of **real64** operands of f_i ,
- The **real64** operands' value.

Since a node's value, adjoint, and data operands are **real64**, they are first converted to **int32** using **transfer** function before stored in the **tape** array, so that each such a value occupies two **storage** entries. After each allotation, the **head** is moved to point to the next empty slot in the array after saving its current value to a **vari** type variable for future retrieval.

2. vari type

The **vari** type is simply a proxy of a node's storage location in the tape

```
type :: vari
  integer(ik) :: i = 0
  procedure(chain_op), pass, pointer :: chain
contains
  !....
```

where **i** is the index to the beginning of a node's storage, and the **chain** procedure encodes the node's operation f_i . **chain** follows an interface that describes the chain rule operation

```

abstract interface
  subroutine chain_op(this)
    import :: vari
    class(vari), intent(in) :: this
  end subroutine chain_op
end interface

```

An alternative to integer index is to a **pointer** to the according entry in the **tape** array. However, we will need to expand the **storage** when it is filled up, and **Fazang** does this by doubling the **storage** size and use **move_alloc** to restore the original values. Since there is no guarantee that **move_alloc** will keep the original memory, a pointer to the original address would be corrupted.

As a **Fazang** program steps forward, a series of **vari** variables are generated, with their *values* calculated and stored. This is called a *forward pass*. The generated **vari** variables in the forward pass are stored in array **varis**. Each entry in **varis** is a dependent (operation output) of one or more previous entries.

3. var type

The user-facing **var** type serves as proxy to **vari**. Each **var** stores the index of a **vari** in the **varis** array.

```

type :: var
  integer(int32) :: vi
  contains
    procedure :: val
    procedure :: adj
    procedure :: grad
    procedure :: set_chain
end type var

```

After the forward pass, when adjoints are desired, we call **grad** or **gradient** procedure. This initiates a *backward pass*, in which the **varis** array is traversed backward so that each **vari**'s **chain** procedure is called to update the operand adjoints.

```

subroutine grad(this)
  class(var), intent(in) :: this
  integer i
  call callstack % varis (this%vi) % init_dependent()
  do i = callstack % head - 1, 1, -1
    call callstack % varis(i) % chain()
  end do
end subroutine grad

```

Here **callstack** is the module variable that encapsulate **tape** and **varis** arrays.

4. Nested tape

Fazang use `begin_nested()` and `end_nested()` to record and terminate a nested tape. With call `begin_nested()` Fazang records the current `tape` and `varis` array head. When `end_nested()` is called, the storage between the recorded head and current head are wiped, and the head is moved back to the recorded location. Multiple levels of nested environment are supported this way.

CHAPTER 5

Add operation functions

Adding an operation f_i involves creating functions for forward pass and backward pass. Let us first use `log` function as a simple example.

First, we create a `log_v` function for the forward pass.

```
impure elemental function log_v(v) result(s)
  type(var), intent(in) :: v
  type(var) :: s
  s = var(log(v%val()), [v])
  call s%set_chain(chain_log)
end function log_v
```

The function generates a new `var` variable `s` using a special constructor `var(value, array of operands)` which stores the value as well as the single operand `v`'s index (in the `tape` storage array). It also points `s`'s chain to a dedicated procedure `chain_log`.

```
subroutine chain_log(this)
  class(vari), intent(in) :: this
  real(rk) :: adj(1), val(1)
  val = this%operand_val()
  adj(1) = this%adj() / val(1)
  call this%set_operand_adj(adj)
end subroutine chain_log
```

To understand this function, recall the recursion in Section 3, assume the `log` operation is node i , then $f_i = \log()$ and z_i is the operand `v`, and the new `var s` would be z_{i-1} . During the backward pass when the node is visited, `chain_log` first retrieves current (z_i, z_i^{adj}) using `operand_val()` and `operand_adj()`, then updates z_i^{adj} with an additional

$$z_{i-1}^{\text{adj}} \frac{df_i}{dz_i} = z_{i-1}^{\text{adj}} \frac{d \log(z_i)}{dz_i} = \frac{z_{i-1}^{\text{adj}}}{z_i}.$$

Adding a binary operation $f_i(z_i^{(1)}, z_i^{(2)})$ is slightly more complex, as we will need to address possibly different scenarios when $z_i^{(1)}$ and $z_i^{(2)}$ are either `var` or `real64`. Let us use overloaded division operator `/` as an example.

With

$$f_i(z_i^{(1)}, z_i^{(2)}) = z_i^{(1)} / z_i^{(2)}$$

we need to account for

- both $z_i^{(1)}$ and $z_i^{(2)}$ are `var`'s
- $z_i^{(1)}$ is `var`, $z_i^{(2)}$ is `real64`,

- $z_i^{(1)}$ is real64, $z_i^{(2)}$ is var,

For the first scenario, we create

```

impure elemental function div_vv(v1, v2) result(s)
  type(var), intent(in) :: v1, v2
  type(var) :: s
  s = var(v1%val() / v2%val(), [v1, v2])
  call s%set_chain(chain_div_vv)
end function div_vv

```

Similar to the log example, we create a new **s** with both operands stored. In the corresponding chain procedure, we need update the adjoints of both **v1** and **v2**.

```

subroutine chain_div_vv(this)
  class(vari), intent(in) :: this
  real(rk) :: adj(2), val(2)
  val = this%operand_val()
  adj(1) = this%adj() / val(2)
  adj(2) = - this%val() * this%adj() / val(2)
  call this%set_operand_adj(adj)
end subroutine chain_div_vv

```

For the second scenario, we create

```

impure elemental function div_vd(v, d) result(s)
  type(var), intent(in) :: v
  real(rk), intent(in) :: d
  type(var) :: s
  s = var(v%val() / d, [v], [d])
  call s%set_chain(chain_div_vd)
end function div_vd

```

Again we create a new var **s**. But this time we use another constructor **var(value, var operands, data operands)** to store value, var operand **v**, and real64 operand **d**. In the corresponding backward pass chain procedure, not only we need retrieve var operand **v** but also data operand **d**, as the new adjoint of $z_i^{(1)}$ is

$$z_i^{(1)\text{new adj}} = z_i^{(1)\text{old adj}} + z_{i-1}^{\text{adj}} \frac{df_i}{dz_i^{(1)}} = z_i^{(1)\text{old adj}} + z_{i-1}^{\text{adj}} \frac{1}{dz_i^{(2)}}$$

So with **v** as $z_i^{(1)}$ and **d** as $z_i^{(2)}$ we have

```

subroutine chain_div_vd(this)
  class(vari), intent(in) :: this
  real(rk) d(1), adj(1)
  d = this%data_operand()
  adj(1) = this%adj() / d(1)
  call this%set_operand_adj(adj)
end subroutine chain_div_vd

```

The third scenario is treated similarly.

APPENDIX A

Fazang Functions

Function	Argument(s)	Operation
<code>sin</code>	scalar or array	same as intrinsic
<code>cos</code>	scalar or array	same as intrinsic
<code>tan</code>	scalar or array	same as intrinsic
<code>asin</code>	scalar or array	same as intrinsic
<code>acos</code>	scalar or array	same as intrinsic
<code>atan</code>	scalar or array	same as intrinsic
<code>log</code>	scalar or array	same as intrinsic
<code>exp</code>	scalar or array	same as intrinsic
<code>sqrt</code>	scalar or array	same as intrinsic
<code>erf</code>	scalar or array	same as intrinsic
<code>erfc</code>	scalar or array	same as intrinsic
<code>abs</code>	scalar or array	same as intrinsic
<code>norm2</code>	1D array	same as intrinsic
<code>hypot</code>	scalars or arrays	same as intrinsic
<code>sinh</code>	scalar of array	same as intrinsic
<code>cosh</code>	scalar of array	same as intrinsic
<code>tanh</code>	scalar of array	same as intrinsic
<code>asinh</code>	scalar of array	same as intrinsic
<code>acosh</code>	scalar of array	same as intrinsic
<code>atanh</code>	scalar of array	same as intrinsic
<code>log_gamma</code>	scalar or array	same as intrinsic
<code>square</code>	scalar or array	For input <code>x</code> , calculate <code>x**2</code>
<code>inv</code>	scalar or array	For input <code>x</code> , calculate <code>1/x</code>
<code>inv_square</code>	scalar or array	For input <code>x</code> , calculate <code>1/x**2</code>
<code>inv_sqrt</code>	scalar or array	For input <code>x</code> , calculate <code>1/sqrt(x)</code>
<code>logit</code>	scalar or array	For input <code>x</code> , calculate <code>log(x/(1-x))</code>
<code>inv_logit</code>	scalar or array	For input <code>x</code> , calculate <code>1/(1+exp(-x))</code>
<code>operator (+)</code>	scalars or arrays	same as intrinsic
<code>operator (-)</code>	scalars or arrays	same as intrinsic
<code>operator (*)</code>	scalars or arrays	same as intrinsic
<code>operator (/)</code>	scalars or arrays	same as intrinsic
<code>operator (**)</code>	scalars	same as intrinsic
<code>sum</code>	1D array	same as intrinsic
<code>dot_product</code>	1D arrays	same as intrinsic
<code>log_sum_exp</code>	1D array	For input <code>x</code> , calculate <code>log(sum(exp((x))))</code>
<code>matmul</code>	2D arrays	same as intrinsic

APPENDIX B

Fazang Probability distributions

1. Normal distribution

$$\text{Normal}(y, \mu, \sigma) = \prod_{i=1}^n \frac{1}{\sqrt{2\pi}\sigma} \exp\left(-\frac{1}{2} \left(\frac{y_i - \mu}{\sigma}\right)^2\right), \quad \forall y \in \mathbb{R}^n, \mu \in \mathbb{R}, \sigma \in \mathbb{R}^+. \quad (\text{B.1})$$

- `normal_lpdf(y, mu, sigma)`
 - `y`: real64 array.
 - `mu`: real64 or var.
 - `sigma`: real64 or var.
 - Return: the log of $\text{Normal}(y, \mu, \sigma)$.

2. LogNormal distribution

$$\text{LogNormal}(y, \mu, \sigma) = \prod_{i=1}^n \frac{1}{\sqrt{2\pi}\sigma y_i} \exp\left(-\frac{1}{2} \left(\frac{\log y_i - \mu}{\sigma}\right)^2\right), \quad \forall y \in (\mathbb{R}^+)^n, \mu \in \mathbb{R}, \sigma \in \mathbb{R}^+. \quad (\text{B.2})$$

- `lognormal_lpdf(y, mu, sigma)`
 - `y`: real64 array.
 - `mu`: real64 or var.
 - `sigma`: real64 or var.
 - Return: the log of $\text{LogNormal}(y, \mu, \sigma)$.

3. **TODO** additional distributions

Bibliography

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