stdin

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```

7 directories, 30 files

./reverse.h

```
/*
* Simple Tape-Based Reverse Mode Autodiff
 * This header-only C implementation provides reverse mode automatic
* differentiation using a dynamic computation tape and operator overloading
 * on a custom `var t` type.
* Each `var_t` variable corresponds to a node on a global tape. The tape
* records the computation graph by tracking the operation and parent variables
 * for each intermediate result.
* After constructing a function from input `var t`s, a reverse pass propagates
 * gradients from the output node backward through the tape using the chain
* rule.
* Usage Example:
* To compute \partial f/\partial x and \partial f/\partial y for f(x, y) = \sin(x) + y^2:
* tape_t tape = tape_create(64);
* tape load(tape);
* var_t x = var_create(1.0f);
* var t y = var create(2.0f);
* var t f = var sin(x) + var pow(y, var create(2.0f));
* tape_reverse_pass(tape, f);
   // var_adjoint(x) returns \partial f/\partial x, var_adjoint(y) returns \partial f/\partial y
* - Always call `tape_load()` before creating variables.
#ifndef H AUTODIFF
#define H AUTODIFF
#include <stddef.h>
#include <stdio.h>
#include <assert.h>
#include <math.h>
#include <string.h>
const uint32 t MAX TAPE LENGTH = 1 << 24; /* correspond to a ~330mb tap */
typedef enum {
 NIL = 0,
 NEG,
 ADD,
 SUB,
 MUL,
 DIV,
 POW,
 EXP,
 COS.
 SIN,
 SQRT,
} operator t;
```

./reverse.h

```
typedef struct {
  float value;
  float adjoint;
  uint32 t left parent;
  uint32_t right_parent;
  operator_t op;
} tape_entry_t;
typedef struct {
  uint32_t length;
  uint32_t capacity;
  tape entry t *entries;
} tape_t;
typedef struct {
  uint32_t index;
} var_t;
/* should not be set directly, use `tape_load` instead */
static tape_t global_tape = {
  .length = 0,
  .capacitv = 0.
  .entries = NULL,
};
/*
 * setting the initial capacity of the tape to a number like 64 will prevent too
 * much calls to realloc
static tape_t tape_create(uint32_t capacity) {
  assert(capacity <= MAX_TAPE_LENGTH);</pre>
  tape entry t *entries = (tape entry t *) calloc(capacity, sizeof(tape entry t));
  if (entries == NULL) {
    perror("tape malloc");
    exit(1);
    return {};
  return {
    .length = 0.
    .capacity = capacity,
    .entries = entries,
  };
static void tape_destroy(tape_t tape) {
  free(tape.entries);
static void tape_extend(tape_t tape) {
  assert(tape.length < MAX_TAPE_LENGTH);</pre>
  if (tape.length == tape.capacity) {
    tape.entries = (tape entry t *) realloc(tape.entries, 2 * tape.capacity * sizeof(*tape.entries));
    if (tape.entries == NULL) {
      perror("tape realloc");
      exit(1);
```

./reverse.h

```
return;
    memset(tape.entries + tape.capacity, 0, tape.capacity * sizeof(*tape.entries));
    tape.capacity = 2 * tape.capacity:
 ++tape.length;
static void tape clear(tape t tape) {
 memset(tape.entries, 0, tape.length * sizeof(*tape.entries));
 tape.length = 0;
static void tape load(tape t tape) {
 global tape = tape;
static tape_t tape_loaded() {
  return global_tape;
static void tape_reverse_pass(tape_t tape, var_t start) {
 for (size t i = 0: i < tape.length: ++i)
    tape.entries[i].adjoint = 0;
 tape.entries[start.index].adjoint = 1;
 for (size_t i = start.index+1; i-- > 0;) { /* avoid size_t wraps */
    tape entry t *entry = &tape.entries[i];
    tape entry t *left parent entry = &tape.entries[entry->left parent];
    tape_entry_t *right_parent_entry = &tape.entries[entry->right_parent];
    switch (entry->op) {
      case NIL:
        break:
      case NEG:
        left parent entry->adjoint += entry->adjoint * -1;
        break;
      case ADD:
        left_parent_entry->adjoint += entry->adjoint * 1;
        right parent entry->adjoint += entry->adjoint * 1;
        break;
      case SUB:
        left parent entry->adjoint += entry->adjoint * 1:
        right parent entry->adjoint += entry->adjoint * -1;
        break;
      case MUL:
        left_parent_entry->adjoint += entry->adjoint * right_parent entry->value;
        right_parent_entry->adjoint += entry->adjoint * left_parent_entry->value;
        break;
      case DIV:
        left parent entry->adjoint += entry->adjoint / right parent entry->value;
        right_parent_entry->adjoint += entry->adjoint *-1* (entry->value / right_parent_entry->value);
        break:
      case POW:
        left parent entry->adjoint += entry->adjoint * right parent entry->value * (entry->value / left parent entry->value);
        right parent entry->adjoint += entry->adjoint * entry->value * logf(left parent entry->value);
        break;
```

./reverse.h

```
case EXP:
        left parent entry->adjoint += entry->adjoint * entry->value;
        break;
      case COS:
        left parent entry->adjoint += entry->adjoint * -1 * sqrtf(1 - entry->value*entry->value);
        break:
        left_parent_entry->adjoint += entry->adjoint * sqrtf(1 - entry->value*entry->value);
        break;
      case SQRT:
        left_parent_entry->adjoint += entry->adjoint / (2 * entry->value);
        break;
    }
  }
}
/* append new variable to global tape */
static var t var create(float value) {
  var_t a = {global_tape.length};
  tape_extend(global_tape);
  global tape.entries[a.index].value = value;
  return a;
static float var adjoint(var t a) {
  return global_tape.entries[a.index].adjoint;
static float var value(var t a) {
  return global tape.entries[a.index].value;
/* variable operations */
static var t operator-(var t a) {
  tape entry t *a entry = &global tape.entries[a.index];
  var t b = var create(-a entry->value);
  tape_entry_t *b_entry = &global_tape.entries[b.index];
  b_entry->op = NEG;
  b entry->left_parent = a.index;
  return b;
/* variable variable operations */
static var t operator+(var t a, var t b) {
  tape_entry_t *a_entry = &global_tape.entries[a.index];
  tape_entry_t *b_entry = &global_tape.entries[b.index];
  var_t c = var_create(a_entry->value + b_entry->value);
  tape entry t *c entry = &global tape.entries[c.index];
  c entry->op = ADD;
  c_entry->left_parent = a.index;
  c entry->right parent = b.index;
  return c:
static var t operator-(var t a, var t b) {
```

```
tape_entry_t *a_entry = &global_tape.entries[a.index];
  tape_entry_t *b_entry = &global_tape.entries[b.index];
  var_t c = var_create(a_entry->value - b_entry->value);
  tape entry t *c entry = &global tape.entries[c.index];
  c entry->op = SUB;
  c_entry->left_parent = a.index;
  c entry->right parent = b.index;
  return c;
static var_t operator*(var_t a, var_t b) {
  tape_entry_t *a_entry = &global_tape.entries[a.index];
  tape_entry_t *b_entry = &global_tape.entries[b.index];
  var t c = var create(a entry->value * b entry->value);
  tape entry t *c entry = &global tape.entries[c.index];
  c_entry->op = MUL;
  c entry->left parent = a.index;
  c_entry->right_parent = b.index;
  return c;
static var_t operator/(var_t a, var_t b) {
  tape entry t *a entry = &qlobal tape.entries[a.index]:
  tape entry t *b entry = &global tape.entries[b.index];
  var t c = var create(a entry->value / b entry->value);
  assert(b entry->value != 0);
  tape_entry_t *c_entry = &global_tape.entries[c.index];
  c entry->op = DIV;
  c_entry->left_parent = a.index;
  c entry->right parent = b.index;
  return c;
static void operator+=(var_t &a, var_t b) {
  a = a + b;
static void operator==(var_t &a, var_t b) {
  a = a - b;
static void operator*=(var_t &a, var_t b) {
  a = a * b;
static void operator/=(var_t &a, var_t b) {
  a = a / b;
/* variable functions */
static var_t var_pow(var_t a, var_t b) {
  tape entry t *a entry = &global tape.entries[a.index];
  assert(a entry->value > 0):
  tape entry t *b entry = &global tape.entries[b.index];
  var t c = var create(powf(a entry->value, b entry->value));
  tape entry t *c entry = &global tape.entries[c.index];
```

```
c_entry->op = POW;
  c entry->left parent = a.index;
  c_entry->right_parent = b.index;
  return c;
static var_t var_exp(var_t a) {
  tape_entry_t *a_entry = &global_tape.entries[a.index];
  var_t b = var_create(expf(a_entry->value));
  tape entry t *b entry = &global tape.entries[b.index];
  b_entry->op = EXP;
  b_entry->left_parent = a.index;
  return b;
static var t var cos(var t a) {
  tape entry t *a entry = &global tape.entries[a.index];
  var_t b = var_create(cosf(a_entry->value));
  tape_entry_t *b_entry = &global_tape.entries[b.index];
  b_entry->op = COS;
  b entry->left parent = a.index;
  return b;
static var t var sin(var t a) {
  tape_entry_t *a_entry = &global_tape.entries[a.index];
  var_t b = var_create(sinf(a_entry->value));
  tape_entry_t *b_entry = &global_tape.entries[b.index];
  b entry->op = SIN;
  b entry->left parent = a.index;
  return b;
static var t var sqrt(var t a) {
  tape entry t *a entry = &global tape.entries[a.index];
  /* assert(a entry->value > 0); */
  var_t b = var_create(sqrtf(a_entry->value));
  tape_entry_t *b_entry = &global_tape.entries[b.index];
  b entry->op = SQRT;
  b_entry->left_parent = a.index;
  return b;
#endif
```

./forward.h

```
/*
 * Simple Vectorized Forward Mode Autodiff
 * This header-only C implementation provides forward mode automatic
 * differentiation using operator overloading on a custom `var t` type.
 * Each `var_t` variable holds:
 * - `value`: the scalar value of the variable.
   - `grad[GRADLEN]`: a gradient vector representing the derivative of the
     variable with respect to each input in a vector of size `GRADLEN`.
 * Usage Example:
 * To compute \partial f/\partial x and \partial f/\partial y for f(x, y) = \sin(x) + y^2:
 * var t x = {.value = 1.0}; x.grad[0] = 1; // \partial x/\partial x = 1
 * var t y = {.value = 2.0}; y.grad[1] = 1; // \partial y/\partial y = 1
 * var_t f = var_sin(x) + var_pow(y, 2);
   // f.value holds the result, f.grad[0] is \partial f/\partial x, f.grad[1] is \partial f/\partial y
 * - The macro `GRADLEN` must be defined before including this header.
#ifndef H AUTODIFF
#define H_AUTODIFF
#include <stddef.h>
#include <stdio.h>
#include <assert.h>
#include <math.h>
#include <string.h>
/* gradient length */
#ifndef GRADLEN
#error "The GRADLEN macro must set before including fowrard.h"
#define GRADLEN 0
#endif
typedef struct {
  float grad[GRADLEN]:
  float value;
} var_t;
 * initialize an new variable that does not derive from the input vector (see
 * above description)
static void var zero(var t *a) {
  memset(a, 0, \overline{\text{sizeof}}(*a));
/* variable operations */
static var_t operator-(var_t a) {
  for (size t i = 0; i < GRADLEN; i++)
```

./forward.h

```
a.grad[i] = -a.grad[i];
  a.value = -a.value;
  return a;
}
/* variable variable operations */
static var t operator+(var_t a, const var_t &b) {
  for (size_t i = 0; i < GRADLEN; i++)
    a.grad[i] = a.grad[i] + b.grad[i];
  a.value = a.value + b.value;
  return a;
static var t operator-(var t a, const var t &b) {
  for (size t i = 0; i < GRADLEN; i++)
    a.grad[\overline{i}] = a.grad[i] - b.grad[i];
  a.value = a.value - b.value;
  return a;
static var t operator*(var t a, const var t &b) {
  for (size_t i = 0; i < GRADLEN; i++)
    a.grad[i] = b.value * a.grad[i] + a.value * b.grad[i];
  a.value = a.value * b.value;
  return a;
static var t operator/(var_t a, const var_t &b) {
  assert(b.value != 0);
  for (size t i = 0; i < GRADLEN; i++)
    a.grad[i] = (b.value * a.grad[i] - a.value * b.grad[i]) / (b.value * b.value);
  a.value = a.value / b.value;
  return a;
static void operator+=(var_t &a, const var_t &b) {
  for (size_t i = 0; i < GRADLEN; i++)
    a.grad[i] = a.grad[i] + b.grad[i];
  a.value = a.value + b.value;
static void operator-=(var t &a. const var t &b) {
  for (size t i = 0; i < GRADLEN; i++)
    a.grad[i] = a.grad[i] - b.grad[i];
  a.value = a.value - b.value;
static void operator*=(var_t &a, const var_t &b) {
  for (size t i = 0; i < GRADLEN; i++)
    a.grad[i] = b.value * a.grad[i] + a.value * b.grad[i];
  a.value = a.value * b.value;
static void operator/=(var t &a, const var t &b) {
  assert(b.value != 0);
  for (size t i = 0; i < GRADLEN; i++)
```

./forward.h

```
a.grad[i] = (b.value * a.grad[i] - a.value * b.grad[i]) / (b.value * b.value);
  a.value = a.value / b.value;
/* variable float operations */
static var_t operator+(var_t a, float b) {
  a.value += b;
  return a;
static var_t operator-(var_t a, float b) {
  a.value -= b;
  return a;
static var t operator*(var t a, float b) {
  for (size t i = 0; i < GRADLEN; i++)
    a.grad[\overline{i}] *= b;
  a.value *= b;
  return a;
static var t operator/(float a, var t b) {
  for (size t i = 0; i < GRADLEN; i++)
    b.grad[i] = -a * b.grad[i] / (b.value * b.value);
  b.value = a / b.value;
  return b;
static void operator+=(var t &a, float b) {
  a.value += b;
static void operator-=(var t &a, float b) {
  a.value -= b;
static void operator*=(var_t &a, float b) {
  for (size t i = 0; i < GRADLEN; i++)
    a.grad[\overline{i}] *= b;
  a.value *= b;
static void operator/=(float a, var t &b) {
  for (size t i = 0; i < GRADLEN; i++)
    b.grad[i] = -a * b.grad[i] / (b.value * b.value);
  b.value = a / b.value;
/* variable functions */
static var_t var_pow(var_t a, float b) {
  assert(a.value > 0);
  float pow = powf(a.value, b-1);
  for (size t i = 0; i < GRADLEN; i++)
    a.grad[i] = b * a.grad[i] * pow;
  a.value = powf(a.value, b);
```

./forward.h

```
return a;
static var_t var_exp(var_t a) {
  float expa = expf(a.value);
  for (size_t i = 0; i < GRADLEN; i++)
    a.grad[i] = a.grad[i] * expa;
  a.value = expa;
  return a;
static var_t var_cos(var_t a) {
  float sina = -sinf(a.value);
  for (size t i = 0; i < GRADLEN; i++)
    a.grad[i] = a.grad[i] * sina;
  a.value = cosf(a.value);
  return a;
static var_t var_sin(var_t a) {
  float cosa = cosf(a.value);
  for (size_t i = 0; i < GRADLEN; i++)</pre>
    a.grad[i] = a.grad[i] * cosa;
  a.value = sinf(a.value);
  return a;
static var_t var_sqrt(var_t a) {
  /* assert(a.value > 0); */
  for (size_t i = 0; i < GRADLEN; i++)</pre>
    a.grad[\overline{i}] = 0.5 * a.grad[i] / sqrtf(a.value);
  a.value = sqrtf(a.value);
  return a;
#endif
```

```
#include <stdlib.h>
#include <stdio.h>
#include <strings.h>
#include <math.h>
#include <time.h>
const int N = 1000; /* number of terms in the reimann sum */
const int DEG = 10; /* degree of the polynomial proximation */
const float START = 0; /* the start of the integration interval */
const float END = 2; /* the end of the integration interval */
const float ITERATIONS = 5000; /* number of gradient descent iterations */
const float ALPHA = 0.001; /* gradient descent speed */
#include "../../reverse.h"
/* the function to approximate */
float f(float x) {
  if (x == 0) return 0;
  return exp(-1 / (x*x));
var_t poly_eval(var_t P[DEG+1], float x) {
  var t val = P[0]:
  float X = x;
  for (size_t i = 1; i < DEG+1; i++) {
    val += P[i] * var create(X);
    X *= x;
  return val;
void poly_init(var_t P[DEG+1]) {
  for (size t i = 0; i < DEG+1; ++i) {
    P[i] = var create(i+1);
}
void poly_print(float P[DEG+1]) {
  printf("polynomial: ");
  for (size t i = 0; i < DEG; ++i) {
    printf("%f, ", P[i]);
  printf("%f\n", P[DEG]);
var_t reimann_integral(var_t P[DEG+1]) {
  var t loss = \{0\};
  float step size = (END-START) / N;
  for (size_t j = 0; j < N; ++j) {
    float x = START + j*step_size;
    var t delta = poly eval(P, x) - var create(f(x));
    loss = loss + (delta*delta) * var create(step size);
  return loss;
```

./example/polynomial approximation/reverse.cpp

```
}
void polynomial_approximation(float P_coef[DEG+1]) {
  tape_t tape = tape_create(64);
  tape load(tape);
  var_t P[DEG+1];
  poly_init(P);
  for (size_t i = 0; i < ITERATIONS; ++i) {</pre>
    /* reimann integral */
    var_t loss = reimann_integral(P);
    /* printf("loss: %f\n", var value(loss)); */
    /* gradient descent */
    tape_reverse_pass(tape, loss);
    for (size_t j = 0; j < DEG+1; ++j) {
      /* normalize the influance of X^j */
      float one_over_norm_of_xj = j / (powf(END, j+1) - powf(START, j+1));
      P_coef[j] = var_value(P[j]) - ALPHA * var_adjoint(P[j]) * one_over_norm_of_xj;
    /* update polynomial */
    tape clear(tape);
    for (size_t j = 0; j < DEG+1; ++j) {
      P[j] = var_create(P_coef[j]);
  tape_destroy(tape);
int main() {
  float P[DEG+1]:
  polynomial approximation(P);
  poly_print(P);
  return 0;
```

```
#include <stdlib.h>
#include <stdio.h>
#include <strings.h>
#include <math.h>
#include <time.h>
const int N = 1000; /* number of terms in the reimann sum */
const int DEG = 10; /* degree of the polynomial proximation */
const float START = 0; /* the start of the integration interval */
const float END = 2; /* the end of the integration interval */
const float ITERATIONS = 5000; /* number of gradient descent iterations */
const float ALPHA = 0.001; /* gradient descent speed */
#define GRADLEN (DEG+1)
#include "../../forward.h"
/* the function to approximate */
float f(float x) {
  if (x == 0) return 0;
  return exp(-1 / (x*x));
var t poly eval(var t P[DEG+1], float x) {
  var t val = P[0];
  float X = x;
  for (size_t i = 1; i < DEG+1; i++) {
    val += P[i] * X;
    X *= x;
  return val;
void poly init(var t P[DEG+1], size t grad start, size t grad end) {
  for (size_t i = 0; i < DEG+1; ++i) {
    P[i] = \{.grad = \{0\}, .value = (float) i+1\};
    if (i >= grad start && i < grad end) {
      P[i].grad[i - grad_start] = 1;
 }
void polv print(float P[DEG+1]) {
  printf("polynomial: ");
  for (size t i = 0; i < DEG; ++i) {
    printf("%f, ", P[i]);
  printf("%f\n", P[DEG]);
var_t reimann_integral(var_t P[DEG+1]) {
  var_t loss = {0};
  float step size = (END-START) / N;
  for (size t j = 0; j < N; ++j) {
    float x = START + j*step_size;
    var t delta = poly eval(\overline{P}, x) - f(x);
```

./example/polynomial_approximation/forward.cpp

```
loss = loss + (delta*delta) * step_size;
  return loss;
void polynomial_approximation(float P_coef[DEG+1]) {
  var_t P[DEG+1];
  poly_init(P, 0, GRADLEN);
  for (size_t i = 0; i < ITERATIONS; ++i) {</pre>
    /* reimann integral */
    var t loss = reimann integral(P);
    /* printf("loss: %f\n", loss.value); */
    /* gradient descent */
    for (size_t j = 0; j < DEG+1; ++j) {
      float one_over_norm_of_xj = j / (powf(END, j+1) - powf(START, j+1));
      /* normalize the influance of X^i */
      P[j].value -= ALPHA * loss.grad[j] * one_over_norm_of_xj;
  }
  for (size t i = 0; i < DEG+1; ++i) {
    P_coef[i] = P[i].value;
int main() {
  float P[DEG+1];
  polynomial_approximation(P);
  poly_print(P);
  return 0;
```

```
#include <stdlib.h>
#include <stdio.h>
#include <strings.h>
#include <math.h>
#include <time.h>
#include <pthread.h>
#include <assert.h>
const int N = 1000; /* number of terms in the reimann sum */
const int DEG = 10; /* degree of the polynomial proximation */
const float START = 0; /* the start of the integration interval */
const float END = 2; /* the end of the integration interval */
const float ITERATIONS = 5000; /* number of gradient descent iterations */
const float ALPHA = 0.001; /* gradient descent speed */
#define GRADLEN 32
#include "../../forward.h"
/* the function to approximate */
float f(float x) {
  if (x == 0) return 0;
  return exp(-1 / (x*x));
var_t poly_eval(var_t P[DEG+1], float x) {
  var t val = P[0];
  float X = x;
  for (size_t i = 1; i < DEG+1; i++) {
    val += P[i] * X;
    X *= x;
  return val;
void poly print(float P[DEG+1]) {
  printf("polynomial: ");
  for (size_t i = 0; i < DEG; ++i) {
    printf("%f, ", P[i]);
  printf("%f\n", P[DEG]);
#define RI WORKERS 2
const size t RI CHUNKS = ((DEG+1 + GRADLEN-1) / GRADLEN);
typedef struct {
  size_t start_chunk;
  size_t end_chunk;
  float *P;
  float *grad;
  float value;
} ri worker param t;
void *ri worker(void *param ptr) {
  ri_worker_param_t *param = (ri_worker_param_t *) param_ptr;
```

./example/polynomial_approximation/forward_parallel.cpp

```
for (size_t chunk_id = param->start_chunk; chunk_id < param->end_chunk; ++chunk_id) {
    var t loss = \{0\};
    var_t P[DEG+1] = \{0\};
    for (size t i = 0; i < DEG+1; ++i) {
      P[i].value = param->P[i];
      if (i >= chunk_id * GRADLEN && i < chunk_id * GRADLEN + GRADLEN) {</pre>
        P[i].grad[i - chunk_id * GRADLEN] = 1;
    float step_size = (END-START) / N;
    for (size_t j = 0; j < N; ++j) {
      float x = START + j*step_size;
      var t delta = poly eval(P, x) - f(x);
      loss = loss + (delta*delta) * step size;
      if (loss.value != loss.value) {
    if (chunk id == param->start chunk) {
      param->value = loss.value;
    } else {
      /* assert(param->value == loss.value); */
    for (size t i = 0; i < GRADLEN && chunk id * GRADLEN + <math>i < DEG+1; ++i) {
      param->grad[chunk_id * GRADLEN + i] = loss.grad[i];
  return NULL;
float reimann integral(float P[DEG+1], float grad[DEG+1]) {
  pthread t worker threads[RI WORKERS];
  ri_worker_param_t worker_params[RI_WORKERS];
  assert(RI_CHUNKS > 0);
  int handled chunks = 0;
  for (size_t worker_id = 0; worker_id < RI_WORKERS; ++worker_id) {</pre>
    size_t start_chunk = (size_t) ((float) RI_CHUNKS / RI_WORKERS * worker_id);
    size_t end_chunk = (size_t) ((float) RI_CHUNKS / RI WORKERS * (worker id+1));
    worker params[worker id] = {
      .start chunk = start chunk,
      .end chunk = end chunk,
      P = P
      .grad = grad,
    int err = pthread_create(&worker_threads[worker_id], NULL, &ri_worker, &worker_params[worker_id]);
      printf("pthread_create error %d", err);
      exit(1):
      return 0:
    handled_chunks += end_chunk - start_chunk;
```

./example/polynomial approximation/forward parallel.cpp

```
assert(handled_chunks == RI_CHUNKS);
  for (size_t worker_id = 0; worker_id < RI_WORKERS; ++worker_id) {</pre>
    int err = pthread join(worker threads[worker id], NULL);
    if (err) {
      printf("pthread_join error %d", err);
      exit(1);
      return err;
  float value;
  for (size t worker id = 1; worker id < RI WORKERS; ++worker id) {</pre>
    if (worker params[worker id].end chunk - worker params[worker id].start chunk > 0) {
      value = worker_params[worker_id].value;
  return value;
void polynomial_approximation(float P[DEG+1]) {
  for (size_t i = 0; i < DEG+1; ++i) {
    P[i] = i+1:
  for (size_t i = 0; i < ITERATIONS; ++i) {</pre>
    /* reimann integral */
    float loss_grad[DEG+1];
    float loss = reimann_integral(P, loss_grad);
    /* printf("loss: %f\n", loss); */
    /* gradient descent */
    for (size_t j = 0; j < DEG+1; ++j) {
      float one_over_norm_of_xj = j / (powf(END, j+1) - powf(START, j+1));
      /* normalize the influance of X^i */
      P[j] -= ALPHA * loss_grad[j] * one_over_norm_of_xj;
}
int main() {
  float P[DEG+1]:
  polynomial approximation(P);
  poly_print(P);
  return 0;
```

```
#include <stdio.h>
#include "../../reverse.h"

int main() {
    tape_t tape = tape_create(64);
    tape_load(tape);

    var_t a = var_create(4);
    var_t b = var_create(9);
    var_t c = var_create(7);
    var_t d = var_create(-2);

    var_t e = var_pow(var_sqrt(a / (b + c * a) + var_exp(var_create(1) / d)), -var_create(3));
    tape_reverse_pass(tape, e);
    printf("value: %f\n", var_value(e));
    printf("grad: {%f, %f, %f, %f}\n", var_adjoint(a), var_adjoint(b), var_adjoint(c), var_adjoint(d));
    tape_destroy(tape);
    return 0;
}
```

```
1
```

./example/hello_world/forward.cpp

```
#include <stdio.h>
#define GRADLEN 4
#include "../../forward.h"

int main() {
    var_t a = {.grad = {1, 0, 0, 0}, .value = 4};
    var_t b = {.grad = {0, 1, 0, 0}, .value = 9};
    var_t c = {.grad = {0, 0, 1, 0}, .value = 7};
    var_t d = {.grad = {0, 0, 0, 1}, .value = -2};

    var_t e = var_pow(var_sqrt(a / (b + c * a) + var_exp(1 / d)), -3);
    printf("value: %f\n", e.value);
    printf("grad: {%f, %f, %f, %f}\n", e.grad[0], e.grad[1], e.grad[2], e.grad[3]);
    return 0;
}
```

```
#include <stdlib.h>
#include <stdio.h>
#include <strings.h>
#include <math.h>
#include <time.h>
const int N = 1000; /* number of terms in the reimann sum */
const float START = 0; /* the start of the integration interval */
const float END = 2; /* the end of the integration interval */
#ifndef DEG
#warning "DEG set to default value 4"
const int DEG = 4; /* degree of the polynomial proximation */
#endif
#include "../../reverse.h"
/* the function to approximate */
float f(float x) {
  if (x == 0) return 0;
  return exp(-1 / (x*x));
var t poly eval(var t P[DEG+1], float x) {
  var t val = P[0];
  float X = x;
  for (size_t i = 1; i < DEG+1; i++) {
    val = val + P[i] * var_create(X);
    X *= x;
  return val;
void poly init(var t P[DEG+1]) {
  for (size t i = 0; i < DEG+1; ++i) {
    P[i] = var create(i+1);
var_t reimann_integral(var_t P[DEG+1]) {
  var_t loss = var_create(0);
  float step size = (END-START)/N;
  for (size t j = 0; j < N; ++j) {
    float x = START + j*step_size;
    var t delta = poly eval(P, x) - var create(f(x));
    loss = loss + (delta*delta) * var_create(step_size);
  return loss;
int main() {
  size t runs = 10;
  float start time, end time;
  start_time = (float) clock() / CLOCKS_PER_SEC;
```

```
2
```

```
for (size_t i = 0; i < 10; ++i) {
  var_t P[DEG+1];
  tape_t tape = tape_create(64);
  tape_load(tape);
  poly_init(P);
  var_t loss = reimann_integral(P);
  tape_destroy(tape);
}
end_time = (float) clock() / CLOCKS_PER_SEC;

/* print average runtime in milliseconds */
  printf("%f", (end_time - start_time) / runs * 1000);
  return 0;</pre>
```

./benchmark/polynomial_approximation/primal.cpp

```
#include <stdlib.h>
#include <stdio.h>
#include <strings.h>
#include <math.h>
#include <time.h>
const int N = 1000; /* number of terms in the reimann sum */
const float START = 0; /* the start of the integration interval */
const float END = 2; /* the end of the integration interval */
#ifndef DEG
#warning "DEG set to default value 4"
const int DEG = 4; /* degree of the polynomial proximation */
#endif
/* the function to approximate */
float f(float x) {
  if (x == 0) return 0;
  return exp(-1 / (x*x));
float poly_eval(float P[DEG+1], float x) {
  float val = P[0];
  float X = x:
  for (size t i = 1; i < DEG+1; i++) {
    val = val + P[i] * X;
    X *= x;
  return val;
void poly_init(float P[DEG+1]) {
  for (size_t i = 0; i < DEG+1; ++i) {
    P[i] = i+1;
float reimann integral(float P[DEG+1]) {
  float loss = 0;
  float step_size = (END-START)/N;
  for (size_t j = 0; j < N; ++j) {
    float x = START + i*step size:
    float delta = poly eval(P, x) - f(x);
    loss = loss + (delta*delta) * step size;
  return loss;
int main() {
  size_t runs = 10;
  float start_time, end_time;
  start time = (float) clock() / CLOCKS PER SEC;
  for (size_t i = 0; i < runs; ++i) {
    float P[DEG+1];
```

```
poly_init(P);
  float loss = reimann_integral(P);
}
end_time = (float) clock() / CLOCKS_PER_SEC;
/* print average runtime in milliseconds */
printf("%f", (end_time - start_time) / runs * 1000);
return 0;
```

```
all: build
CC := clang
CFLAGS := -std = c + +11 -02 - lm
primal: primal.cpp
        $(if $(DEG),,$(error Must set DEG))
        $(CC) $(CFLAGS) -DDEG=$(DEG) primal.cpp -o primal_build_$(DEG)
reverse: reverse.cpp
        $(if $(DEG),,$(error Must set DEG))
        $(CC) $(CFLAGS) -DDEG=$(DEG) reverse.cpp -o reverse_build_$(DEG)
forward: forward.cpp
        $(if $(DEG),,$(error Must set DEG))
# I renamed GRADLEN to GL to avoid the overriding of GRADLEN
        $(eval GL := $(shell echo ${DEG}+1 | bc))
        $(CC) $(CFLAGS) -DDEG=$(DEG) -DGRADLEN=$(GL) forward.cpp -o forward_build_$(DEG)
forward_novec: forward.cpp
        $(if $(DEG),,$(error Must set DEG))
# I renamed GRADLEN to GL to avoid the overriding of GRADLEN
        $(eval GL := $(shell echo ${DEG}+1 | bc))
        $(CC) $(CFLAGS) -fno-vectorize -fno-slp-vectorize -DDEG=$(DEG) -DGRADLEN=$(GL) forward.cpp -o forward build novec $(DEG)
forward_gradlen: forward.cpp
        $(if $(DEG),,$(error Must set DEG))
        $(if $(GRADLEN),,$(error Must set GRADLEN))
        $(CC) $(CFLAGS) -DDEG=$(DEG) -DGRADLEN=$(GRADLEN) forward.cpp -o forward build gradlen $(DEG) $(GRADLEN)
parallel: forward_parallel.cpp
        $(if $(DEG),,$(error Must set DEG))
        $(CC) $(CFLAGS) -DDEG=$(DEG) forward parallel.cpp -o parallel build $(DEG)
parallel workers: forward parallel.cpp
        $(if $(DEG),,$(error Must set DEG))
        $(if $(WORKERS),,$(error Must set WORKERS))
        $(CC) $(CFLAGS) -DDEG=$(DEG) -DRI_WORKERS=$(WORKERS) forward_parallel.cpp -o parallel_build_workers_$(DEG)_$(WORKERS)
# use -j option to run build in parallel
build: reverse forward forward novec forward gradlen parallel
clean:
        rm primal build * forward build * reverse build * parallel build *
```

```
#include <stdlib.h>
#include <stdio.h>
#include <strings.h>
#include <math.h>
#include <time.h>
const int N = 1000; /* number of terms in the reimann sum */
const float START = 0; /* the start of the integration interval */
const float END = 2; /* the end of the integration interval */
#ifndef DEG
#warning "DEG set to default value 4"
const int DEG = 4; /* degree of the polynomial proximation */
#endif
#ifndef GRADLEN
#warning "GRADLEN set to default value 8"
#define GRADLEN 8
#endif
#include "../../forward.h"
/* the function to approximate */
float f(float x) {
  if (x == 0) return 0:
  return exp(-1 / (x*x));
var_t poly_eval(var_t P[DEG+1], float x) {
  var t val = P[0];
  float X = x;
  for (size t i = 1; i < DEG+1; i++) {
    val += P[i] * X;
    X *= x;
  return val;
void poly_init(var_t P[DEG+1], size_t grad_start, size_t grad_end) {
  for (size_t i = 0; i < DEG+1; ++i) {
    P[i] = \{.grad = \{0\}, .value = (float) i+1\};
    if (i >= grad_start && i < grad_end) {</pre>
      P[i].grad[i - grad_start] = 1;
 }
var_t reimann_integral(var_t P[DEG+1]) {
  var t loss = \{0\};
  float step size = (END-START) / N;
  for (size_t j = 0; j < N; ++j) {
    float x = START + j*step_size;
    var t delta = poly eval(P, x) - f(x);
    loss = loss + (delta*delta) * step size;
  return loss;
```

./benchmark/polynomial_approximation/forward.cpp

```
int main() {
    size_t runs = 10;
    float start_time, end_time;

start_time = (float) clock() / CLOCKS_PER_SEC;
for (size_t i = 0; i < runs; ++i) {
    for (size_t grad_start = 0; grad_start < DEG+1; grad_start += GRADLEN) {
        var_t P[DEG+1];
        poly_init(P, grad_start, grad_start + GRADLEN);
        var_t loss = reimann_integral(P);
    }
} end_time = (float) clock() / CLOCKS_PER_SEC;

/* print average runtime in milliseconds */
    printf("%f", (end_time - start_time) / runs * 1000);
    return 0;
}</pre>
```

```
#include <stdlib.h>
#include <stdio.h>
#include <strings.h>
#include <math.h>
#include <time.h>
#include <pthread.h>
#include <assert.h>
const int N = 1000; /* number of terms in the reimann sum */
const float START = 0; /* the start of the integration interval */
const float END = 2; /* the end of the integration interval */
const float ITERATIONS = 5000; /* number of gradient descent iterations */
const float ALPHA = 0.001; /* gradient descent speed */
#ifndef DEG
#warning "DEG set to default value 4"
const int DEG = 4; /* degree of the polynomial proximation */
#ifndef GRADLEN
#define GRADLEN 64
#endif
#include "../../forward.h"
/* the function to approximate */
float f(float x) {
  if (x == 0) return 0;
  return exp(-1 / (x*x));
var t poly eval(var t P[DEG+1], float x) {
  var_t val = P[0];
  float X = x;
  for (size t i = 1; i < DEG+1; i++) {
    val += P[i] * X;
    X *= x;
  return val;
void poly_print(float P[DEG+1]) {
  printf("polynomial: ");
  for (size t i = 0: i < DEG: ++i) {
    printf("%f, ", P[i]);
  printf("%f\n", P[DEG]);
#ifndef RI_WORKERS
#define RI WORKERS 2
const size_t RI_CHUNKS = ((DEG+1 + GRADLEN-1) / GRADLEN);
tvpedef struct {
  size t start chunk;
  size_t end_chunk;
  float *P;
```

./benchmark/polynomial_approximation/forward_parallel.cpp

```
float *grad;
  float value;
} ri_worker_param_t;
void *ri worker(void *param ptr) {
  ri_worker_param_t *param = (ri_worker_param_t *) param_ptr;
  for (size_t chunk_id = param->start_chunk; chunk_id < param->end_chunk; ++chunk_id) {
    var t loss = \{0\};
    var t P[DEG+1] = \{0\};
    for (size_t i = 0; i < DEG+1; ++i) {
      P[i].value = param->P[i];
      if (i >= chunk id * GRADLEN && i < chunk id * GRADLEN + GRADLEN) {
        P[i].grad[i - chunk id * GRADLEN] = 1;
    float step_size = (END-START) / N;
    for (size_t j = 0; j < N; ++j) {
      float x = START + j*step_size;
      var t delta = poly eval(\overline{P}, x) - f(x);
      loss = loss + (delta*delta) * step_size;
      if (loss.value != loss.value) {
      }
    if (chunk id == param->start chunk) {
      param->value = loss.value;
    } else {
      /* assert(param->value == loss.value); */
    for (size t i = 0; i < GRADLEN && chunk id * GRADLEN + <math>i < DEG+1; ++i) {
      param->grad[chunk id * GRADLEN + i] = loss.grad[i];
  return NULL;
float reimann integral(float P[DEG+1], float grad[DEG+1]) {
  pthread t worker threads[RI WORKERS]:
  ri worker param t worker params[RI WORKERS];
  assert(RI CHUNKS > 0);
  int handled chunks = 0;
  for (size_t worker_id = 0; worker_id < RI_WORKERS; ++worker_id) {</pre>
    size_t start_chunk = (size_t) ((float) RI_CHUNKS / RI_WORKERS * worker id);
    size t end chunk = (size t) ((float) RI CHUNKS / RI WORKERS * (worker id+1));
    worker_params[worker_id] = {
      .start_chunk = start_chunk,
      .end chunk = end chunk,
      P = P
      .grad = grad,
    int err = pthread create(&worker threads[worker id], NULL, &ri worker, &worker params[worker id]);
```

./benchmark/polynomial approximation/forward parallel.cpp

```
if (err) {
      printf("pthread_create error %d", err);
      exit(1);
      return 0;
    handled chunks += end chunk - start chunk;
  assert(handled_chunks == RI_CHUNKS);
  for (size_t worker_id = 0; worker_id < RI_WORKERS; ++worker_id) {</pre>
    int err = pthread_join(worker_threads[worker_id], NULL);
    if (err) {
      printf("pthread join error %d", err);
      exit(1);
      return err;
  float value;
  for (size_t worker_id = 1; worker_id < RI_WORKERS; ++worker_id) {</pre>
    if (worker_params[worker_id].end_chunk - worker_params[worker_id].start chunk > 0) {
      value = worker_params[worker_id].value;
    }
  return value;
int main() {
  size_t runs = 10;
  float start_time, end_time;
  start_time = (float) clock() / CLOCKS_PER_SEC;
  for (size t i = 0; i < runs; ++i) {
    float P[DEG+1];
    float loss grad[DEG+1];
    float loss = reimann_integral(P, loss_grad);
  end_time = (float) clock() / CLOCKS_PER_SEC;
  /* print average runtime in milliseconds */
  printf("%f", (end_time - start_time) / runs * 1000);
  return 0:
```

```
#!/bin/bash
gradlen=64

bench() {
    reverse=$(./reverse_build_"$1")
    forward=$(./forward_build_novec_"$1")
    forward_gradlen=$(./forward_build_gradlen_"$1"_"$gradlen")
    echo "$1","$reverse","$forward","$forward_novec","$forward_gradlen"
}

deg=(1 2 $(seq 4 4 512))
for d in ${deg[@]}; do
    make -j build DEG=$d GRADLEN=$gradlen > /dev/null &
done
wait

for d in ${deg[@]}; do
    bench $d
done
```

```
#!/bin/bash
d=500
bench() {
   parallel=$(./parallel_build_workers_"$d"_"$1")
   echo "$1","$parallel"
}
workers=$(seq 1 1 12)
for w in ${workers[@]}; do
   make -j parallel_workers DEG=$d WORKERS=$w > /dev/null &
done
wait
for w in ${workers[@]}; do
   bench $w
done
make clean
```

```
1
```

```
#!/bin/bash
bench() {
    reverse=$(./reverse_build_"$1")
    echo "$d","$reverse"
}

deg=(4 8 $(seq 4 16 512))
for d in ${deg[@]}; do
    make -j reverse DEG=$d > /dev/null &
done
wait

for d in ${deg[@]}; do
    bench $d
done
make clean
```

```
#!/bin/bash
bench() {
    parallel=$(./parallel_build_"$1")
    echo "$d","$parallel"
}

deg=(4 8 $(seq 4 16 512))
for d in ${deg[@]}; do
    make -j parallel DEG=$d > /dev/null &
done
wait

for d in ${deg[@]}; do
    bench $d
done

make clean
```

```
#!/bin/bash

deg=300

bench() {
    forward_gradlen=$(./forward_build_gradlen_"$deg"_"$1")
    echo "$1","$forward_gradlen"
}

gradlen=($(seq 4 1 512))
for gl in ${gradlen[@]}; do
    make -j forward_gradlen DEG=$deg GRADLEN=$gl > /dev/null &
done
wait

for gl in ${gradlen[@]}; do
    bench $gl
done

make clean
```