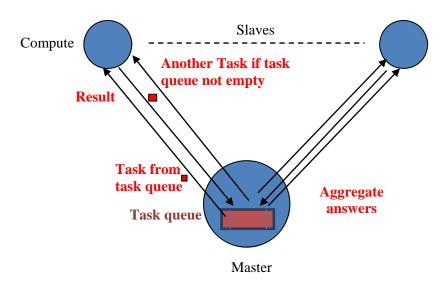
SUZAKU Pattern Programming Framework Specification

2 - Workpool Pattern Version 1

B. Wilkinson, March 17, 2016

2.1 Workpool

The workpool pattern is like a master-slave pattern but has a task queue that provides load balancing, as shown below. Individual tasks are given to the slaves. When a slave finishes a task and returns the result, it is given another task from the task queue, until the task queue is empty. At that point, the master waits until all



outstanding results are returned. The termination condition is the task queue empty and all result collected.

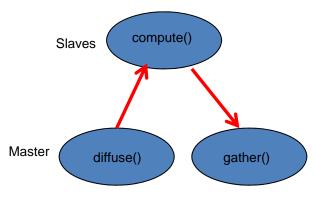
Workpool with a task queue

Algorithm. In the implementation of the workpool described here (version 1), the data items being sent between the master process and slave processes are limited to 1-D arrays. The programmer deposes the problem into **T** tasks. Each task consists of a 1-D array of **D** doubles with an associated task ID. Each slave result for a task consists of a 1-D array of **R** doubles with the associated task ID. The master sends out tasks to slaves. Slaves return results and are given new tasks, or a terminator message if there are no more tasks, i.e. if the number of tasks sent reaches **T**. The number of tasks can be less than number of slaves, equal to the number of slaves, or greater than the number of slaves. If the number of tasks is the same as the number of slaves, the workpool becomes essentially a master-slave pattern.

Programmer-written routines. The Suzaku workpool interface is modeled on the Seeds framework. The programmer must implement four routines:

- init() Sets values for the number of tasks (T), the number of data items in each task (D), and the number of data items in each result (R). Called once by all processes at the beginning of the computation.
- **diffuse**() Generates the next task when called by the master
- compute() Executed by a slave, takes a task generated by diffuse and generates the corresponding result

gather() Accepts a slave result and uses it to develop the final answer



Message passing done by framework

diffuse, compute, and gather routines

diffuse(), compute(), and gather() are dependent upon the application, and often very short.

Workpool code. The workpool itself is implemented by the provided routine **SZ_Workpool()** placed within a parallel section. **init()**, **diffuse()**, **compute()**, and **gather()** are called by **SZ_Workpool()** and given as input parameters. They can be re-named to accommodate for example multiple workpools in a single program.

2.2 Program structure

The program structure is shown below and consists of the four programmer routines and the Suzaku routines.

```
#include <stdio.h>
#include <string.h>
#include "suzaku.h"
void init(int *T, int *D, int *R) {
        return;
void diffuse(int *taskID,double output[D]) {
        return;
}
void compute(int taskID, double input[D], double output[R]) {
        return;
}
void gather(int taskID, double input[R]) {
        return;
}
int main(int argc, char *argv[]) {
                                               // number of processes
        int P;
        SZ_Init(P);
                                               // initialize MPI message-passing environment
```

```
SZ_Parallel_begin

SZ_Workpool(init, diffuse, compute, gather);

SZ_Parallel_end;

printf("Workpool results\n ... ", ....); // print out workpool results ...

SZ_Finalize();

return 0;
}
```

Workpool program structure

2.3 Signatures of Programmer-Written Routines

Init

```
void init(int *tasks, int * data_items, int *result_items)
```

This routine will be called at the beginning of the workpool by all processes. At the very least, it must set values for number of tasks (T), number of data items in each task (D), and number of data items in each result (R). The routine may be used for other initialization purposes. There is no implicit synchronization.

Parameters (pointers to integers):

```
    int *tasks
    int *data_items
    int parameter for the number of data items (doubles) in each task
    int *result items
    Input parameter for the number of data items (doubles) in result of each task
```

Limitations: The number of tasks, \mathbf{T} , must be equal or less than $INT_MAX - 1$, but this is very unlikely to be an issue. For \mathbf{D} and \mathbf{R} , the limiting factor is the maximum size of dynamic arrays on the platform.

A typical coding sequence would be:

Any names can be used in the formal parameter list.

Diffuse

The signature of this routine is:

void diffuse(int taskID, double output[D])

This routine generates the next task when called by the master.

Parameters:

int taskID Input parameter for the task ID for the associated task

double output[D] Output parameter for the task data, given as an array of **D** doubles

Notes: **taskID** is provided by the framework, from 0 to **T** - 1. Each time diffuse is called, **taskID** is incremented. **taskID** is carried with the task throughout the workpool. **taskID** provides a mechanism to do specific actions with particular tasks or results. When used by the programmer, **taskID** corresponds to a segment number in Seeds.

Compute

The signature of this routine is:

void compute(int taskID, double input[D], double output[R])

This routine is executed by a slave. It takes a task generated by diffuse and generates the corresponding result.

Parameters:

int taskID Input parameter for the task ID for the associated task

double input[D] Input parameter for the task data, given as an array of D doublesdouble output[R] Output parameter for the result, given as an array of R doubles

Gather

The signature of this routine is:

void gather(int taskID, double input[R])

This routine accepts a slave result and develops the final answer. Called by the master.

Parameters:

int taskID Input parameter for the task ID for the associated task

double input[R] Input parameter for the slave result, given as an array of R doubles

Notes: **gather()** is used to aggregate the answer from the task results during the workpool operation, and programmers are free to do this any way they like and whatever the application dictates. To be able to reach the answer from outside gather, the final answer can be declared globally at the top of the program outside main.

2.4 Signature of Suzaku Workpool Routine

This routine is provided and implements the workpool. It calls **init()**, **diffuse()**, **compute()**, and **gather()**. **SZ Workpool()** must be called within a Suzaku parallel section. The signature of the routine is:

Parameters:

*init Function pointer to init function
 *diffuse Function pointer to diffuse function
 *compute Function pointer to compute function
 *compute Function pointer to gather function

Notes: The function pointers could have been eliminated if their names were standardized (as in Seeds), e.g. init, diffuse, compute, and gather, but specifying the function names makes it more obvious which functions are used by the workpool, and also allows multiple workpools each using different function pointers. No data arrays need to be declared for Suzaku. These are generated by Suzaku dynamically.

The programmer can implement routines outside the workpool as the need arises. Results from the gather need to be used to create the final answer and variables declared outside main to be reachable from gather and other routines.

2.5 Compilation and Execution

Workpool code: The workpool routine **SZ_Workpool()** is implemented in **suzaku.c.** It can be compiled with:

mpicc -c -o suzaku.o suzaku.c

to create an object file **suzaku.o** (note the -c option). This avoids having to recompile **suzaku.c** every time you compile application code.

Application code: **SZ_Workpool**() does not use **suzaku.h** itself but since a workpool needs to be within a parallel section, the application code must include **suzaku.h**. For the commands below, the two files:

suzaku.h suzaku.o

must be placed in the same directory as the source file. To compile an application workpool program **prog1.c**, issue the command:

mpicc -o prog1 prog1.c suzaku.o

A make file is provided for the sample programs.

Instead of pre-compiling **suzaku.c** into **suzaku.o**, one could also compile both **suzaku.c** and **prog1.c** together with:

mpicc -o prog1 prog1.c suzaku.c

To execute **prog1**, issue the command:

mpiexec –n <no of processes> prog1

where **<no_of processes>** is the number of processes you wish to use. The workpool needs at least two processes, master and one slave. Note the master does not act as one slave as in the master-slave pattern in Part 1 because collective routines are not used.

2.6 Debug Messages

A version of the **SZ_Workpool**() routine is provided that includes print statements to see how the tasks are allocated slaves and results returned. This version is called **SZ_Workpool_debug**() and can be found in **suzaku.c**. To rename **SZ_Workpool**() in the application code to **SZ_Workpool_debug**() and recompile.



to

use,

Sample output with debug messages

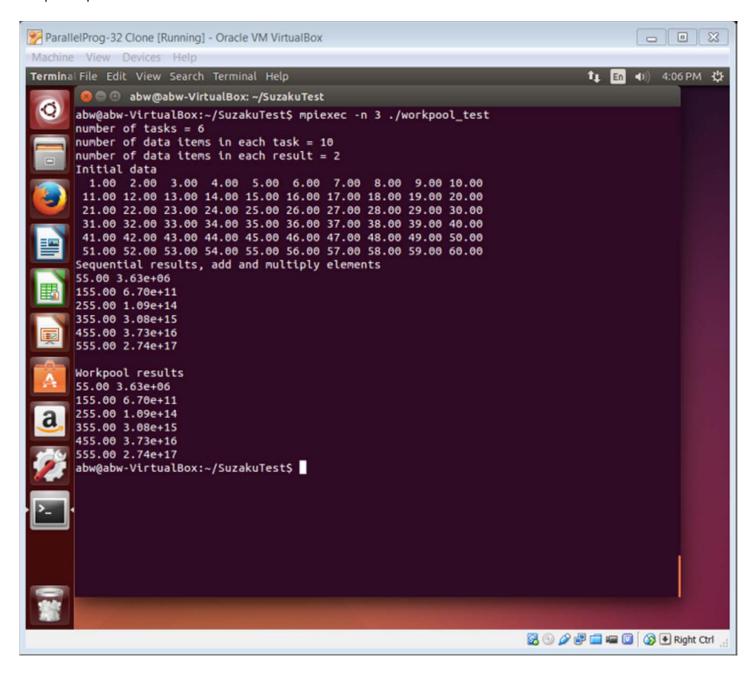
Sample programs

1. Adding and multiplying numbers

```
// Suzaku Workpool pattern version 1 Application: Adding and multiplying numbers. B. Wilkinson April 3, 2015
#include <stdio.h>
#include <string.h>
#include "suzaku.h"
                          // number of tasks, max = INT_MAX - 1
#define T 6
#define D 10
                          // number of data items in each task, doubles only
#define R 2
                          // number of data items in result of each task, doubles only
// No arrays need to be declared for Suzaku. Following used in this particular application, not required in general
double workpool result[T][R];
                                            // Final results computed by workpool, in gather()
double task[T][D];
                                            // Initial data, T tasks, each task D elements, created by initialize() for testing
// workpool functions to be provided by programmer:
void init(int *tasks, int *data_items, int *result_items) {
         *tasks = T;
         *data_items = D;
         *result_items = R;
         return;
}
void diffuse(int *taskID,double output[D]) {
                                                     // taskID not used
         static int temp = 0;
                                                     // only initialized first time function called
         for (j = 0; j < D; j++)
                 output[j] = ++temp;
                                                     // set elements to consecutive data values
}
void compute(int taskID, double input[D], double output[R]) {
                                   // function done by slaves -- simply adding the numbers together, and multiply them
         output[0] = 0;
         output[1] = 1;
         int i:
         for (i = 0; i < D; i++) {
                 output[0] += input[i];
                 output[1] *= input[i];
         return;
}
void gather(int taskID, double input[R]) {
                                                     // function done by master collecting slave results
                                                     // Final results computed by master, uses taskID
         int j;
         for (j = 0; j < R; j++) {
                 workpool_result[taskID][j] = input[j];
        }
}
// additional routines used in this application
void initialize() {
                                            // create initial data for sequential testing, not used by workpool
         int i,j;
         int temp = 0;
         for (i = 0; i < T; i++) {
                                            // initialize data
                 for (j = 0; j < D; j++)
                                                     // set elements to consecutive data values
                          task[i][j] = ++temp;
         printf("Initial data\n");
                                                              // print out data
         for (i = 0; i < T; i++) {
                 for (j = 0; j < D; j++)
```

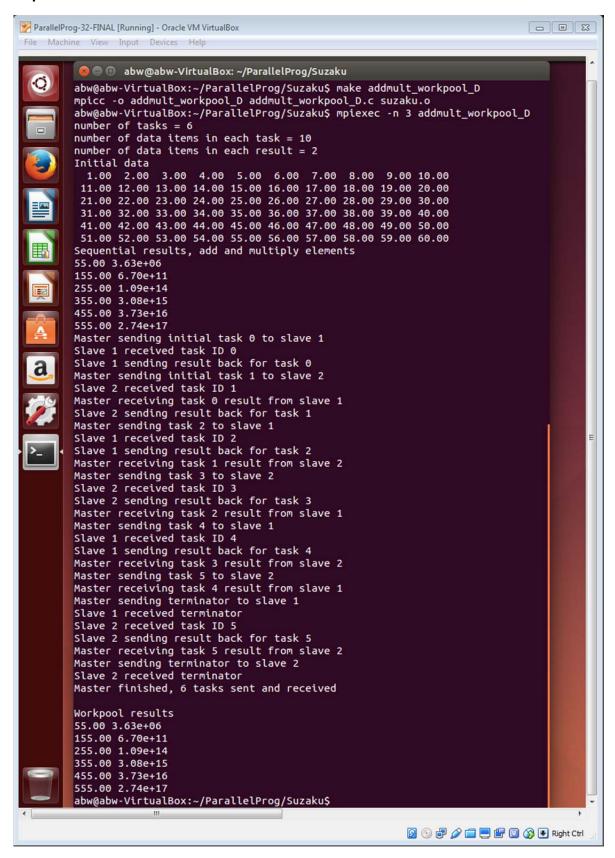
```
printf(" %5.2f",task[i][j]);
                  printf("\n");
         }
}
void compute_seq() {
                                                        // Compute results sequentially and print out
         int i,j;
                                              //** Final results computed sequentially
         double seq_result[T][D];
         printf("Sequential results, add and multiply elements\n"); // print out results
         for (i = 0; i < T; i++) {
                  seq_result[i][0] = 0;
                  seq_result[i][1] = 1;
                  for (j = 0; j < D; j++) {
                            seq_result[i][0] += task[i][j]; // add up all numbers in task result in [0]
                            seq_result[i][1] *= task[i][j]; // multiply up all numbers in task result in [1]
                  printf("%5.2f ", seq_result[i][0]);
                                                                 // print result 0
                  printf("%5.2e", seq_result[i][1]);
printf("\n");
                                                                 // print result 1
         }
}
int main(int argc, char *argv[]) {
                                                        // All variables declared here are in every process
         int i;
                                                        // number of processes, set by SZ_Init(P)
         int P;
         SZ_Init(P);
                                                        // initialize MPI message-passing, sets P to be number of processes
         printf("number of tasks = %d\n",T);
         printf("number of data items in each task = %d\n",D);
         printf("number of data items in each result = %d\n",R);
         initialize();
                                                        // create initial data for sequential testing, not used by workpool
                                                        // compute results sequentially and print out
         compute_seq();
         SZ_Parallel_begin
         SZ_Workpool(init,diffuse,compute,gather);
         SZ Parallel end;
                                                        // end of parallel
         printf("\nWorkpool results\n");
                                                        // print out workpool results
         for (i = 0; i < T; i++) {
                  printf("%5.2f ", workpool_result[i][0]);
printf("%5.2e", workpool_result[i][1]);
printf("\n");
                                                                           // result
                                                                           // result
         SZ_Finalize();
         return 0;
}
```

Sample output



Version with debug messages

Sample output:



This version could be used for educational purposes.

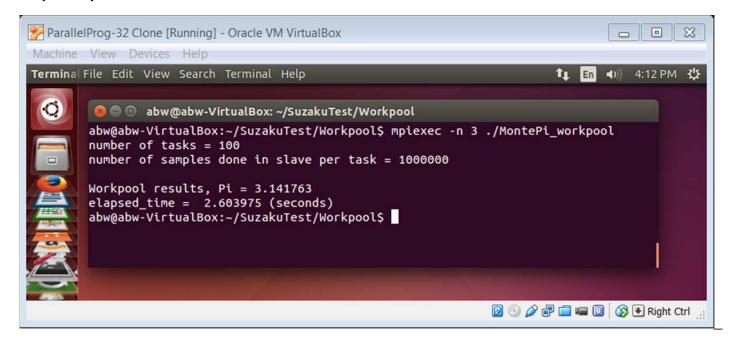
2. Monte Carlo Pi calculation

MontePi workpool.c:

```
// Suzaku Workpool pattern version 1 -- Application: Monte Carlo Pi. B. Wilkinson April 4, 2015
#include <stdio.h>
#include <string.h>
#include <stdlib.h>
#include <time.h>
#include "suzaku.h"
// required Suzaku constants
#define T 100
                          // number of tasks, max = INT_MAX - 1
#define D 1
                          // number of data items in each task, doubles only
#define R 1
                          // number of data items in result of each task, doubles only
// constant used in computation
                          // sample pts done in a slave
#define S 1000000
// gobal variable
double total = 0:
                          // final result
// required workpool functions
void init(int *tasks, int *data_items, int *result_items) {
         *tasks = T;
        *data_items = D;
        *result_items = R;
        return;
}
void diffuse(int *taskID,double output[D]) {
                                           // taskID not used in computation
        static int temp = 0;
                                           // only initialized first time function called
        output[0] = ++temp;
                                           // set seed to consecutive data value
}
void compute(int taskID, double input[D], double output[R]) {
        int i;
        double x, y;
        double inside = 0;
        srand(input[0]);
                                           // initialize random number generator
        for (i = 0; i < S; i++) {
                 x = rand() / (double) RAND_MAX;
                 y = rand() / (double) RAND_MAX;
                 if ((x * x + y * y) \le 1.0) inside++;
        output[0] = inside;
        return;
}
void gather(int taskID, double input[R]) {
        total += input[0];
                                           // aggregate answer
}
// additional routines used in this application
double get_pi() {
        double pi;
        pi = 4 * total / (S*T);
        printf("\nWorkpool results, Pi = %f\n",pi);
                                                             // print out workpool results
```

```
}
int main(int argc, char *argv[]) {
                                  // All variables declared here are in every process
        int i:
        int P;
                                  // number of processes, set by SZ_Init(P)
        clock_t time1, time2; // use clock for timing
        SZ Init(P);
                                  // initialize MPI environment, sets P to number of processes
        printf("number of tasks = %d\n",T);
        printf("number of samples done in slave per task = %d\n",S);
        time1 = clock(); // record time stamp
                                  // start of parallel section
        SZ_Parallel_begin
        SZ_Workpool(init,diffuse,compute,gather);
        SZ Parallel end;
                                  // end of parallel
        time2 = clock(); // record time stamp
                         // calculate final result
        get_pi();
        printf("elapsed_time =\t%lf (seconds)\n", (double)(time2 - time1)/CLOCKS_PER_SEC);
        SZ_Finalize();
        return 0;
}
```

Sample output



3 Matrix multiplication

matrixmult_workpool.c

```
// Suzaku Workpool pattern version 1 -- Application: Matrix Multiplication. B. Wilkinson April 5, 2015
#include <stdio.h>
#include <string.h>
#include <stdlib.h>
#include <time.h>
#include "suzaku.h"
// required Suzaku constants
#define T 9
                          // number of tasks, max = INT_MAX - 1
#define D 6
                          // number of data items in each task, 3 elements of row A and 3 elements of column B
#define R 1
                          // number of data items in result of each task
#define N 3
                          // size of arrays
double A[N][N], B[N][N], C[N][N], Cseq[N][N];
// required workpool functions
void init(int *tasks, int *data_items, int *result_items) {
         *tasks = T;
        *data_items = D;
        *result_items = R;
        return;
}
void diffuse(int taskID,double output[D]) {
                                                    // uses same approach as Seeds sample but inefficient copying arrays
                                                    // taskID used in computation
        int i;
        int a, b;
        a = taskID / N;
        b = taskID % N;
        for (i = 0; i < N; i++) {
                                                    //Copy one row of A and one column of B into output
                 output[i] = A[a][i];
                 output[i+N] = B[i][b];
        return;
}
void compute(int taskID, double input[D], double output[R]) {
        int i:
        output[0] = 0;
        for (i = 0; i < N; i++) {
                 output[0] += input[i] * input[i+N];
        return;
}
void gather(int taskID, double input[R]) {
        int a,b;
        a = taskID / 3;
        b = taskID % 3;
        C[a][b]= input[0];
        return;
}
// additional routines used in this application
```

```
void initialize() { // initialize arrays
         int i,j;
         for (i = 0; i < N; i++){
                  for(j = 0; j < N; j++) {
                            A[i][j] = i + N * j + 1;
                            B[i][j] = j + N * i + 1;
                  }
         }
         return;
}
void seq_matrix_mult(double A[N][N], double B[N][N], double C[N][N]) {
         int i,j,k;
         for (i = 0; i < N; i++)
                  for (j = 0; j < N; j++) {
                            C[i][j] = 0;
                           for (k = 0; k < N; k++)
                                    C[i][j] += A[i][k] * B[k][j];
         return;
}
void print_array(double array[N][N]) { // print out an array
         int i,j;
         for (i = 0; i < N; i++){
                  printf("\n");
                  for(j = 0; j < N; j++) {
                           printf("%5.2f ", array[i][j]);
                  }
         printf("\n");
         return;
}
int main(int argc, char *argv[]) {
                                    // All variables declared here are in every process
         int i;
         int P;
                                    // number of processes, set by SZ_Init(P)
         clock_t time1, time2; // use clock for timing
         SZ_Init(P);
                                    // initialize MPI environment, sets P to number of processes
         initialize();
                                    // initialize input arrays
         printf("Array A");
         print_array(A);
         printf("Array B");
         print_array(B);
         seq_matrix_mult(A,B,Cseq);
         printf("Multiplication sequentially");
         print_array(Cseq);
         time1 = clock(); // record time stamp
         SZ_Parallel_begin
                                    // start of parallel section
         SZ_Workpool(init,diffuse,compute,gather);
         SZ Parallel end;
                                    // end of parallel
         time2 = clock(); // record time stamp
         printf("Workpool results");
         print_array(C);
                                    // print final result
         printf("Elapsed_time =\t%lf (seconds)\n", (double)(time2 - time1)/CLOCKS_PER_SEC);
```

```
SZ_Finalize();
return 0;
}
```

Sample output (Note number of processes does not have to be the same as number of elements

