Sheet 07

PS Parallel Programming

Patrick Wintner

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1 Parallelizing Loops

The dependencies and parallelisation posibilities of code snippets are analysized.

1.1

1.1.1 Serial

```
1 for (int i=0; i < n-1; i++) {
2          x[i] = (y[i] + x[i+1]) / 7; // S
3 }</pre>
```

Statement S anti-depends (Write-After-Read) on itself: $S\delta^{-1}S$. Anti-dependencies can be eliminated through variable renaming.

1.1.2 Parallel

```
#pragma omp parallel for schedule(static)
for (int i=0; i < n-1; i++) {
            x2[i] = (y[i] + x[i+1]) / 7;
}</pre>
```

1.2

1.2.1 Serial

```
for (int i=0; i < n; i++) {
    a = (x[i] + y[i]) / (i+1); // S1
    z[i] = a; // S2
}
f = sqrt(a + k); // S3</pre>
```

Statement S2 truly depends (Read-After-Write) on S1 and S3 truly depends on the last instance of S1: $S1\delta S2, S2\delta S3$. The depency is obviously not loop-carried, therefore the loop can be parallelized by making 'a' private within the loop.

1.2.2 Parallel

1.3

1.3.1 Serial

```
1 for (int i=0; i < n; i++) {
2             x[i] = y[i] * 2 + b * i; // S1
3       }
4             for (int i=0; i < n; i++) {
6                  y[i] = x[i] + a / (i+1); // S2
7       }</pre>
```

Statement S2 both truly and anti-depends on S1: $S1\delta S2$, $S1\delta^{-1}S2$. There is no dependence within the loops, therefore the loops themselves can be parallelized.

1.3.2 Parallel

2 Parallelizing more Loops

The dependencies of code snippets are analysized and the code snippets themselves are parallelized. The wall time of both the serial and parallel versions is measured.

2.1 Measurement Method

All measurements were done on the LCC3 cluster by calling sbatch job.sh <executable> 3, e. g. sbatch job.sh a_ser 3 (3 is the number of measurements) with the number of loop iterations set to 100000000.

The following scripts are involved in the experiment.

2.1.1 SLURM Script

```
#!/bin/bash
# usage: sbatch [slurm_options] <executable> <number_of_measurements>
# Execute job in the partition "lva" unless you have special requirements.
# SBATCH --partition=lva
# Name your job to be able to identify it later
```

2.1.2 Benchmark Script

```
#!/bin/bash
  # Usage: ./benchmark.sh <executable> <number_of_measurements>
  ITER=100000000 # number of loop iterations
5 results=$1".dat"
6 echo "x y ey" > $results # create header
7 make toTable
8 make $1
  for i in {1,4,8,12} # number of threads
10
           measurements=$i"_"$1".log"
11
           export OMP_NUM_THREADS=$i
12
           for j in $(seq 1 $2) # repeat measurement £2 times
13
14
                    ./$1 $ITER >> $measurements # store measurement results in <executable>.
15
           done
16
           ./toTable $measurements $results $2 $i #store table in <executable>.dat
           rm $measurements
19 done
20 make clean
```

2.2.1 Serial

```
#ifndef A_SER_H
#define A_SER_H

double factor = 1; // S1

for (int i=0; i < n; i++) {
            x[i] = factor * y[i]; // S2
            factor = factor / 2; // S3
}
#endif</pre>
```

The zero-th instance of both statements S2 and S3 truly depends on S1: $S1\delta S2$, $S1\delta S3$. Furthermore, S2 has true loop-carried dependence on S3 and S3 has a true loop-carried dependence on itself: $S3\delta S2$, $S3\delta S3$, $S3\delta^{-1}S3$.

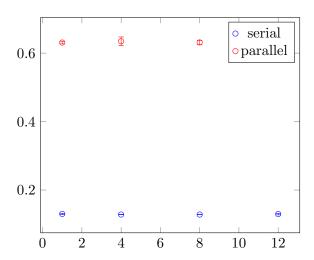
2.2.2 Parallel

```
#ifndef A_PAR_H
   #define A_PAR_H
2
   double *factor = malloc(n*sizeof(*factor));
   if(!factor) {
            free(x);
            free(y);
            free(z);
            perror("malloc");
            return EXIT_FAILURE;
10
   }
11
   factor[0] = 1;
12
   for(int i=0; i < n-1; i++) {
13
            factor[i+1] = factor[i]/2;
14
15
   #pragma omp parallel for schedule(static)
   for (int i=0; i < n; i++) {
17
           x[i] = factor[i] * y[i];
18
   }
19
```

```
20 free(factor);
21 #endif
```

It is used that the different instances of statement S2 do not depend on each other.

2.2.3 Experiment Results



2.3

2.3.1 Serial

```
#ifndef B_SER_H
#define B_SER_H

for (int i=1; i<n; i++) {
            x[i] =(x[i] + y[i-1]) / 2;
            y[i] = y[i] +z[i] * 3;
}

#endif</pre>
```

2.3.2 Parallel

```
#ifndef B_PAR_H
#define B_PAR_H
```

2.3.3 Experiment Results

2.4

2.4.1 **Serial**

2.4.2 Parallel

```
#ifndef C_PAR_H
#define C_PAR_H

x[0] = x[0] + 5 * y[0];
for (int i = 1; i<n; i++) {
            x[i] = x[i] + 5 * y[i];
            if (twice) {
                 x[i-1] = 2 * x[i-1];
            }

#endif</pre>
```

2.4.3 Experiment Results

3 ... And One More Loop to Parallelize

3.1 Serial

```
for (int i = 0; i < 4; ++i) {
    for (int j = 1; j < 4; ++j) {
        a[i + 2][j - 1] = b * a[i][j] + 4;
}
}</pre>
```

3.1.1 Distance and Direction Vectors

i	j	S: a[i+2][j-1]	S: a[i][j]
0	1	a[2][0]=	=a[0][1]
0	2	a[2][1] =	=a[0][2]
0	3	${f a[2][2]} =$	=a[0][3]
1	1	a[3][0]=	=a[1][1]
1	2	a[3][1] =	=a[1][2]
1	3	a[3][2] =	=a[1][3]
2	1	a[4][0] =	=a[2][1]
2	2	a[4][1]=	=a[2][2]
2	3	a[4][2] =	=a[2][3]
3	1	a[5][0] =	=a[3][1]
3	2	a[5][1] =	=a[3][2]
3	3	a[5][2]=	=a[3][3]

dependence relation	array element	distance vector	direction vector
$S[0][2]\delta S[2][1]$	a[2][1]	(2,-1)	(<,>)
$S[0][3]\delta S[2][2]$	a[2][2]	(2,-1)	(<,>)
$S[1][2]\delta S[3][1]$	a[3][1]	(2,-1)	(<,>)
$S[1][3]\delta S[3][2]$	a[3][2]	(2,-1)	(<,>)

There are four loop-carried true data dependencies in the code snippet. All write accesses happen before all read accesses. Therefore the loop can be divided into two loops, so that both loops themselves can be executed in parallel. This solution only works because the number of iterations is low enough. A solution that scales is to only parallelize the inner loop.

3.2 Parallel

```
#pragma omp parallel for schedule(static)
   for (int i = 0; i < 2; ++i) {
           for (int j = 1; j < 4; ++j) {
                   a[i + 2][j - 1] = b * a[i][j] + 4;
           }
   }
6
  #pragma omp parallel for schedule(static)
8
  for(int i = 2; i < 4; ++i) {
           for (int j = 1; j < 4; ++j) {
10
                   a[i+2][j-1] = b * a[i][j] + 4;
11
           }
12
13
  }
```