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The README.md and report.pdf are exactly the same infact to generate the report the readme was converted to tex using pandoc. So prefer reading the markdown

Analysis of the CPU

Theoretical limitations of CPU

For the cpu running lscpu gives

```
Architecture: x86_64

CPU op-mode(s): 32-bit, 64-bit

Address sizes: 46 bits physical, 48 bits virtual

Byte Order: Little Endian

CPU(s): 24

On-line CPU(s) list: 0-23

Vendor ID: GenuineIntel

Model name: 13th Gen Intel(R) Core(TM) i7-13700K

CPU family: 6

Model: 183

Thread(s) per core: 2

Core(s) per socket: 16

Socket(s): 1

Stepping: 1

CPU max MHz: 5400.0000

CPU min MHz: 800.0000

BogoMIPS: 6835.20

Flags: avx avx2 avx_vnni fma ...

Caches (sum of all):

L1d: 640 KiB (16 instances)

L2: 24 MiB (10 instances)

L3: 30 MiB (1 instance)

NUMA:

NUMA node(S): 1

NUMA node(CPU(s): 0-23
```

It shows that the cpu is 13th Gen Intel(R) Core(TM) i7-13700K and has 16 cores and 24 threads, looking more into the intel doc of i7-13700K it has 8 performant cores (with hyper threading of 2) and 8 efficient cores. For the P-cores it has a base frequency of 3.4GHz and turbo frequency of 5.3 GHz while for the E-cores it has a base frequency of 2.5GHz and turbo frequency of 4.2GHz

The i7-13700K operates at a default frequency of 3.4GHz. Since it supports avx2 and each p cores can do 2 inst/cycle on ports 0 and 1 we can achieve 2×8 (8 sp per 256bit vector) $\times 2$ (2 flops per fma) = 32 flops/cycle. For 8 p cores we can achieve at max $8 \times 32 \times 5.3$ GHz = 1356.8 GFLOP/s at turbo frequency

For each e core we achieve 16 flops/cycle and thus we can achieve at max 8 × 16 × 4.2 GHz = 537.6 GFLOP/s at turbo frequency

Total we can achieve a max flops of 1894.4 GFLOP/s

Doing a similar analysis for double precision the number of doubles we can fit in avx256 would be effectively halved and we can fidn the theoretical maximum flops for doubles to be around 947.2 GFLOP/s

PPC Project 9a Report

- [1] 13th Gen Intel (R) Core(TM) i7-13700K
- [2] Default clock freq of i7-13700K

Practical limitations of CPU

For evaluating the practical limitations of the cpu we will load an 8 element vector a into an avx256 register a8 and broadcast a scalar b into another avx256 register b8. Then we perform fused multiply and add operation d8 [k] = a8 * b* + d8 [k] for each of the 8 lanes while repeating this MN times in parallel. This code does a total of \$128MN + 64N\$ flops, since M is of the order of 1e5 we will ignore the 64N. This gives us a total of 128MN flops. Here is the snippet of the main code

```
#pragma omp parallel for schedule(dynamic, 1)
for (uint64_t i = 0; i < M; i++) {
    __m256 d8[8] = {};
    for(int k=0; k<8; k++) {
        // randomly initialize d8[k]
        alignas(32) float v[8] = {};
        for (int j = 0; j < 8; j++) v[j] = rand() / float(RAND_MAX) * 2.0f - 1.0f;
        d8[k] = _mm256_load_ps(v);
    }
    for (uint64_t j = 0; j < N; j++) {
        #pragma unroll
        for(int k=0; k<8; k++) {
            d8[k] = _mm256_fmadd_ps(a8, b8, d8[k]);
        }
    }
}

// Since N is much larger the flops of this is negligible
    __m256 s = _mm256_setzero_ps();
    for(int i=0; i<8; i++) {
        s = _mm256_add_ps(s, d8[i]);
    }
    out[i] += s[0];
}</pre>
```

We time how long it takes to run this using std::chrono::high_resolution_clock and we will also measure the clock cycles during it's execution using __rdtscp intrinsics this will help us get the measured clock freq. We take an average of 5 runs for reporting the wall clock time and the cpu cycles and thus the clock freq and the measured GSLOP/s. The entire code is at cpu_flops.cpp, we will run this with the following flags enabled

```
export OMP_NUM_THREADS=24

g++ -fopenmp -g -03 -march=native cpu_flops.cc -o main

perf stat ./main
```

Running it prints the following

```
Avg Wall time : 3.50322 s

Total FLOPs : 6400 GFLOP

Avg Achieved FLOP : 1826.89 GFLOP/s

Cycles elapsed : 11972383729

Measured CPU freq : 3.41753 GHz

sum : 2.05752

Performance counter stats for './main':

80,781.54 msec task-clock # 23.051 CPUs utilized

2,372 context-switches # 29.363 /sec

159 cpu-migrations # 1.968 /sec
```

```
393 page-faults # 4.865 /sec
389,859,205,718 cycles # 4.826 GHz
500,717,583,488 instructions # 1.28 insn per cycle
50,156,378,434 branches # 620.889 M/sec
1,403,191 branch-misses # 0.00% of all branches

3.504431292 seconds time elapsed

80.759760000 seconds user
0.023995000 seconds sys
```

We also print the sum of output after the exectution so the compiler does not optimize away everything. The main things to see is we achieve 1826.89 GFLOP/s and measured a clock freq of 4.826 GHz, there is also a huge difference between the clock frequencies measured by perf and __rdtscp. The reason could be __rdtscp does not vary according to the current core's frequency. Infact after reading more about it from Time_Stamp_Counter I came to realize it's use is highly discouraged. We achieved 96.43% of the theoretical peak flops. I think this is the practical best performance we can achieve as we never get the clock to run at turbo freq for all p and e core for all the avx256 instruction and there will be some dips in the clock freq degrading performance

Only running the cpu_flops.cpp on the p cores using

```
export OMP_NUM_THREADS=16
export OMP_PLACES="{0,1,2,3,4,5,6,7,8,9,10,11,12,13,14,15}"
export OMP_PROC_BIND=close
g++ -fopenmp -g -O3 -march=native cpu_flops.cc -o main
./main
```

gives us

```
Avg Wall time : 4.79077 s

Total FLOPs : 6400 GFLOP

Avg Achieved FLOP : 1335.9 GFLOP/s

Cycles elapsed : 16372601962

Measured CPU freq : 3.41753 GHz

sum : 2.05752
```

Thus the p cores can reach 99.93% of their peak flops. while running it only on the e cores using

```
export OMP_NUM_THREADS=8
export OMP_PLACES="{16,17,18,19,20,21,22,23}"
export OMP_PROC_BIND=close
g++ -fopenmp -g -03 -march=native cpu_flops.cc -o main
./main
```

gives us

```
Avg Wall time : 11.9549 s

Total FLOPs : 6400 GFLOP

Avg Achieved FLOP : 535.346 GFLOP/s

Cycles elapsed : 40856183393

Measured CPU freq : 3.41753 GHz

sum : 2.05752
```

which is also 99.50% of the theoretical max that the e cores can achieve, I don't have any definite answers as to why only individually e cores and p cores are able to achieve max flops while together they fall short by 2-3% but maybe they cannot be sustained at turbo freq all the time under avx256

Looking at generated assembly code from cpu_flops.s we see that the compiler has generated the 8 fma256 instruction for us

```
.L4:

vfmadd231ps %ymm0, %ymm1, %ymm2

vfmadd231ps %ymm0, %ymm1, %ymm9

vfmadd231ps %ymm0, %ymm1, %ymm7

vfmadd231ps %ymm0, %ymm1, %ymm6

vfmadd231ps %ymm0, %ymm1, %ymm5

vfmadd231ps %ymm0, %ymm1, %ymm5

vfmadd231ps %ymm0, %ymm1, %ymm4

vfmadd231ps %ymm0, %ymm1, %ymm4

vfmadd231ps %ymm0, %ymm1, %ymm3

subq $1, %rax

jne .L4
```

Comparison with CP3B Solution

For comparison with my fastest cp3b solution, we use the exact same code as the 137707 submission with replacing the avx512 instructions with avx256 and adding code for timing measurements and clock freq measurements. We set nx and ny to be 14000. On running

```
export OMP_NUM_THREADS=24
g++ -fopenmp -g -O3 -march=native cp3b.cc -o main
perf stat ./main
```

we see the following output

```
Wall time : 1.73274 s
Total FLOPs : 2744 GFLOP
Achieved FLOPS : 1583.62 GFLOP/s
Cycles elapsed : 5921721988
Measured CPU freq : 3.41756 GHz
Checksum : -0.00906668

Performance counter stats for './main':

45,399.47 msec task-clock # 11.634 CPUs utilized
689 context-switches # 15.176 /sec
25 cpu-migrations # 0.551 /sec
683,659 page-faults # 15.059 K/sec
206,358,697,147 cycles # 4.545 GHz
386,717,416,506 instructions # 1.87 insn per cycle
15,785,912,832 branches # 347.711 M/sec
62,951,893 branch-misses # 0.40% of all branches

3.902154516 seconds time elapsed

44.625873000 seconds user
0.776241000 seconds sys
```

We achieved 83.59% of the theoretical peak flops

Now again inspecting the assembly at cp3b.s we see the compiler did the right thing and generated the vectorized code for us. The assembly code for the innermost loop of the kernel looks like this

```
.L115:

vmovaps (%r8), %ymm1

vbroadcastss (%rdx), %ymm2

addq $32, %rdx

vmovaps (%r8, %rcx, 4), %ymm0

addq $32, %re

vfmadd231ps %ymm2, %ymm1, %ymm4

vfmadd231ps 192(%rsp), %ymm0, %ymm2

vmovaps %ymm2, 192(%rsp),

vbroadcastss -28(%rdx), %ymm2

vfmadd231ps 126(%rsp), %ymm0, %ymm2

vmovaps %ymm2, 128(%rsp)

vbroadcastss -24(%rdx), %ymm2

vfmadd231ps %ymm2, %ymm1, %ymm9

vfmadd231ps %ymm2, %ymm1, %ymm8

vfmadd231ps %ymm2, %ymm0, %ymm15

vbroadcastss -20(%rdx), %ymm2

vfmadd231ps %ymm2, %ymm0, %ymm15

vbroadcastss -16(%rdx), %ymm2

vfmadd231ps %ymm2, %ymm0, %ymm1

vbroadcastss -16(%rdx), %ymm2

vfmadd231ps %ymm2, %ymm0, %ymm10

vbroadcastss -12(%rdx), %ymm2

vfmadd231ps %ymm2, %ymm0, %ymm10

vbroadcastss -12(%rdx), %ymm2

vfmadd231ps %ymm2, %ymm0, %ymm10

vbroadcastss -8(%rdx), %ymm2

vfmadd231ps %ymm2, %ymm0, %ymm14

vbroadcastss -8(%rdx), %ymm2

vfmadd231ps %ymm2, %ymm0, %ymm13

vbroadcastss -4(%rdx), %ymm1

vfmadd231ps %ymm2, %ymm0, %ymm13

vbroadcastss -4(%rdx), %ymm1

vfmadd231ps %ymm2, %ymm0, %ymm12

vfmadd231ps %ymm2, %ymm0, %ymm12

vfmadd231ps %ymm2, %ymm0, %ymm12

vfmadd231ps %ymm1, fol(%rsp)

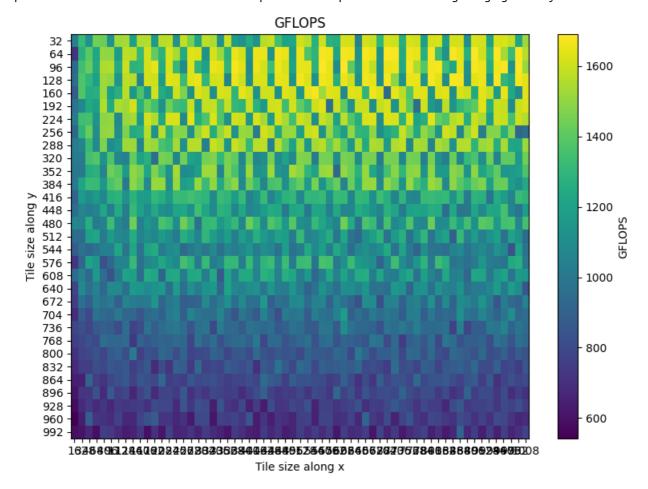
cmpq %rsi, %rdx

jne .L115
```

There are 2 instructions for loop counters, 2 for pointer arithematic, 8 for memory access, 8 for the boradcast operation and 16 for the actual fma operations

When on experimentation with different tile sizes we find that a tile size of Mc=Nc=96 and Kc=560 gives us 1690.02 GFLOP/s, which is 89.92% of the theoretical peak flops and 6% increase from the untuned matmul kernel, plot the heatmap of tile size along x and y shows a pattern that when one a tile can fit in the L2 cache (MC \times KC \times 4 bytes per float<= 24MB/10 (we have 10 instances)) we get high

performance but as soon as the tile exceeds the per core L2 the performance starts degrading significantly



The data from tuning is at tuning_results.csv and the python script for generating the benchmarks and plotting are at autotune.py and plot_heatmap.py

Analysis of the GPU

Theoretical limitations of GPU

For the GPU running nvcc device_query.cpp -o main && ./main gives

```
Name: NVIDIA GeForce RTX 2050
Compute Capability: 8.6
Clock Rate: 1.155 GHz
Max threads per block: 1024
Max threads per multiprocessor: 1536
Threads per warp: 32
Max registers per block: 65536
Max registers per multiprocessor: 65536
Total global memory: 3897 MB
Max shared mem per block: 48 KB
Shared mem per multiprocessor: 102400 B
Multiprocessor count: 16
Max warps per multiprocessor: 48
```

Thus we have the Nvidia GeForce RTX 2050 as our GPU for this analysis which is based on turing architecture. It has 16 streaming multiprocessors (SM) and from the wiki of GeForce RTX 2050 it has 2048 CUDA cores and 64 Tensor Cores. It also has a bandwidth of 112 GB/s and a default clock freq of 1.155 GHz and boost of 1.477 GHz. Each core can execute 1 fma i.e 2 FLOP per cycle, we can then calculate the theoretical max flops as follows

 $\$ \begin{aligned} \text{SP peak @ base} &= 2048 \times 2 \times 1.155!\times!10^9 \approx 4730.88;\mathrm{GFLOP/s}\\text{SP peak @ boost} &= 2048 \times 1.477!\times!10^9 \approx 6049.79;\mathrm{GFLOP/s} \end{aligned} \$\$

This matches with the processing power @ boost in the wiki of GeForce 2050

- [1] Nvidia GeForce RTX datasheet
- [2] RTX Series comparison

Practical limitations of GPU

For doing max computation on gpu we will let each thread execute $N\$ fma i.e each thread will compute a = a * b + a. Each thread then does a total of 2 * N flops which will allow the entire kernel to do blocks * threads * 2 * N flops. Here is our very simple and tiny kernel, the entire code is available at max_flops.cu

```
extern "C" __global__
void gpu_kernel(
    float *out,
    int N, float seed
) {
    int tid = threadIdx.x;
    float a = seed + tid;
    float b = seed - tid;

    for (int i = 0; i < N; i++) {
        a = a * b + a;
    }
    out[blockIdx.x * blockDim.x + tid] = a;
}

gpu_kernel<<<1024, 256>>>(
        d_out, iters, seed
);
```

We run this with nvcc max flops.cu -O3 -arch=sm 86 -o main && ./main with N set to 1e7 and it gives the following output

```
Elapsed Time : 1132.182 ms (1.132 s)
Total FLOP : 5.243e+03 GFLOP
Performance : 4630.773 GFLOP/s
```

We thus achieve 97.88% of the theoretical maximum flops. Profiling it with sudo nou ./main gives us

We can see that our kernel achieved a compute capacity of 99.18% which is all thanks to ~0% memory utilization. We also achieved 95.84% of max occupancy with 46 active warps out of 48

Looking at the ptx code with $nvcc --ptx -arch=sm_86 max_flops.cu -o main.ptx$ we see that the compiler directly generated fma f32 instructions while unrolling the inner loop by 4 for us

```
$L_BB0_3:
fma.rn.f32 %f13, %f2, %f20, %f20;
fma.rn.f32 %f14, %f2, %f13, %f13;
fma.rn.f32 %f15, %f2, %f14, %f14;
fma.rn.f32 %f20, %f2, %f15, %f15;
add.s32 %r13, %r13, -4;
setp.ne.s32 %p3, %r13, 0;
@%p3 bra $L_BB0_3;
```

The executable can be found at max_flops and the ptx code at max_flops.ptx

Comparison with CP5 Solution

Now for comparison with the fastest cp5.cu solution we again benchmark using nvcc cp5.cu -03 -arch=sm_86 -0 main && ./main. It follows the exact same code as submitted to cp5 without any modification and nx and ny set to 14000, the executable can be found at cp5

```
Elapsed Time: 1050.230 ms (1.050 s)
Total FLOP: 2.744e+03 GFLOP
Performance: 2612.948 GFLOP/s
```

It executated $2\times 14\times 14 = 2744$ GFLOP (as we are only computing the upper triangular matrix) in 1.050s achieving 2612.948 GFLOP/s which is 55.23% of the theoretical peak flops @ base. Profiling it with sudo ncu ./main reveals

WRN This kernel's theoretical occupancy (33.3%) is limited by the number of required registers. See the CUDA Best

Practices Guide (https://docs.nvidia.com/cuda/cuda-c-best-practices-

```
guide/index.html#occupancy) for more
details on optimizing occupancy.
```

This time we achieved 70% of max compute throughput and only 33% max theoretical occupancy which is due to high register use of 100 per thread and 16 active warps per block possible and also our memory throughput is 55% of peak. Thus we are register bound

Looking at the ptx code from cp5.ptx we see that the compiler issues vectorized loads from shared memory this time

```
fma.rn.f32 %f701, %f636, %f617, %f605;
fma.rn.f32 %f702, %f636, %f618, %f606;
fma.rn.f32 %f703, %f636, %f619, %f607;
fma.rn.f32 %f704, %f636, %f620, %f608;
ld.shared.v4.f32 {%f705, %f706, %f707, %f708}, [%r74+2048];
ld.shared.v4.f32 {%f713, %f714, %f715, %f716}, [%r74+2304];
ld.shared.v4.f32 {%f721, %f722, %f723, %f724}, [%r77+2048];
ld.shared.v4.f32 {%f729, %f730, %f731, %f732}, [%r77+2304];
fma.rn.f32 %f737, %f721, %f705, %f641;
fma.rn.f32 %f738, %f721, %f706, %f642;
fma.rn.f32 %f739, %f721, %f707, %f643;
```

not only that we can also see that the compiler also generated code for vectorized load from global to shared memory