Alright!

I'll give complete proper answers for each question without shortening, and also suggest counter-questions with their full answers.

Since the document is huge, I'll proceed assignment by assignment and topic by topic in a neat, organized manner.

HPC Viva Full Answers

Assignment 1: Parallel BFS and DFS using OpenMP

Q1. Define BFS and DFS in graph traversal.

Answer:

• BFS (Breadth-First Search):

BFS is a graph traversal algorithm that starts at a selected node (source node) and explores all its neighboring nodes at the current depth before moving on to nodes at the next level. It uses a queue data structure to keep track of nodes to be explored. BFS is particularly useful for finding the shortest path in an unweighted graph.

• DFS (Depth-First Search):

DFS is a graph traversal technique where one starts from the source node and explores as far as possible along each branch before backtracking. It uses a stack (either explicit or implicit through recursion) to remember the nodes to be visited. DFS is useful for tasks like detecting cycles and solving puzzles.

Counter Question:

Which data structures are typically used for BFS and DFS?

Answer:

BFS uses a Queue.

• DFS uses a **Stack** (or recursive function call stack).

Q2. What are the time complexities of BFS and DFS?

Answer:

- Time complexity of BFS: O(V + E)
- Time complexity of DFS: O(V + E)

Where:

- V = number of vertices
- E = number of edges

Both BFS and DFS must explore every vertex and edge in the worst case.

Counter Question:

• What about space complexity for BFS and DFS?

Answer:

- Space Complexity for BFS: O(V) (queue stores nodes at current level)
- Space Complexity for DFS: O(V) (stack can go as deep as the height of the graph)

Q3. How does parallelism improve BFS/DFS performance?

Answer:

Parallelism improves performance by **dividing the workload** among multiple threads or processors.

- In BFS, multiple nodes at the same depth can be processed simultaneously, as their neighbors can be explored independently.
- In DFS, although more sequential by nature, **parallel exploration** of different subtrees can be performed cautiously.
- As a result, the time to traverse large graphs **significantly reduces**.

• Why is BFS more parallelizable than DFS?

Answer:

 BFS operates level-by-level where nodes at the same depth are independent, while DFS depends heavily on previous traversal paths (sequential dependency).

Q4. How is a visited array managed safely in parallel BFS?

Answer:

- The **visited array** tracks whether a node has already been explored.
- In a **parallel BFS**, to manage it safely:
 - Use atomic operations to set/check values.
 - Or use mutex locks to avoid race conditions.
 - Some implementations also assign ownership of nodes to specific threads to prevent overlaps.

This ensures that no two threads visit and process the same node more than once.

Counter Question:

What could happen if we don't protect the visited array?

Answer:

• The same node could be visited multiple times, leading to **incorrect traversal**, **infinite loops**, and **wasted computation**.

Q5. What are critical sections in OpenMP?

Answer:

- A critical section in OpenMP is a piece of code that must be executed by only one thread at a time to prevent race conditions.
- In OpenMP, we use #pragma omp critical to mark a block as critical.

Example:

```
#pragma omp critical
{
    // code that modifies shared variables
```

Counter Question:

• Why should critical sections be minimized?

Answer:

 Overuse of critical sections serializes the code, negating the benefits of parallelism and reducing performance.

Q6. What issues arise when parallelizing DFS?

Answer:

- DFS has a sequential nature because it explores a node's children completely before backtracking.
- Issues include:
 - Difficult to divide work dynamically.
 - Stack management conflicts across threads.
 - Load imbalance: some threads finish early while others work deeper paths.
 - Race conditions when accessing or updating shared resources.

Hence, DFS is less naturally parallelizable compared to BFS.

Counter Question:

• Can DFS be parallelized at all?

Answer:

• Yes, by **parallelizing exploration of different subtrees** after initial levels or using task parallelism (OpenMP tasks).

Q7. What is breadth-wise parallelism in BFS?

- Breadth-wise parallelism refers to processing all nodes at the same level concurrently.
- In BFS, after visiting the nodes at depth d, all their adjacent nodes (depth d+1) can be explored in parallel.
- It improves traversal speed, especially for wide graphs.

• Does breadth-wise parallelism suit dense or sparse graphs more?

Answer:

• It benefits **dense graphs** more because there are more nodes at each level to process in parallel.

Q8. What happens if synchronization is not handled in parallel traversal?

Answer:

- If synchronization is not handled:
 - Multiple threads may visit the same node simultaneously.
 - o Incorrect results or duplicate node visits.
 - Deadlocks or infinite loops can occur.
 - May lead to memory corruption and program crashes.

Thus, careful synchronization is **critical**.

Counter Question:

• Name some synchronization techniques used.

Answer:

• Critical sections, atomic operations, barriers, and locks.

Q9. What OpenMP directive is used to parallelize a for loop?

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• The directive is:

#pragma omp parallel for

• It tells the compiler to **divide iterations of the loop** among available threads automatically.

Counter Question:

• What scheduling options can you specify with parallel for?

Answer:

• Static, dynamic, and guided scheduling.

Q10. What are advantages and disadvantages of parallel BFS?

- Advantages:
 - Faster traversal on large graphs.
 - Better resource utilization of multi-core systems.
 - Scalable to very large graphs with millions of nodes.
- Disadvantages:
 - Synchronization overhead between threads.

- Load balancing challenges if graph structure is irregular.
- Increased complexity in code and debugging.

• When would parallel BFS not help much?

Answer:

• On **small graphs** where overhead outweighs benefits.

Perfect! 🚀

Continuing with full **Assignment 2: Parallel Bubble Sort and Merge Sort using OpenMP** answers properly, along with **counter questions** and **their answers**, just like before:

Assignment 2: Parallel Bubble Sort and Merge Sort using OpenMP

Q1. Why is Merge Sort called a "divide and conquer" algorithm?

- Merge Sort is called a divide and conquer algorithm because it divides the unsorted array into two halves, recursively sorts each half, and then merges the sorted halves to produce the final sorted array.
- It solves the problem by **breaking it into smaller subproblems**, solving them independently (recursively), and combining the results.

• What is the time complexity of Merge Sort?

Answer:

• The time complexity is **O(n log n)** in all cases (worst, average, best).

Q2. What happens if two threads try to swap the same elements in parallel Bubble Sort?

Answer:

- If two threads attempt to swap the **same elements simultaneously** in parallel Bubble Sort:
 - It leads to data races.
 - The swaps could interfere, resulting in **incorrect orderings**.
 - The array could become partially sorted or completely corrupted.

Thus, careful synchronization or assignment of non-overlapping elements is necessary.

Counter Question:

• How can we avoid this problem?

- By ensuring even-odd transposition:
 - o In **even phases**, swap (0,1), (2,3), (4,5), etc.

- o In **odd phases**, swap (1,2), (3,4), (5,6), etc.
- This prevents threads from accessing adjacent elements at the same time.

Q3. How do you divide the array for parallel Merge Sort?

Answer:

- In parallel Merge Sort:
 - The array is recursively divided into subarrays.
 - Each thread handles sorting a subsection of the array.
 - Once the subarrays are sorted, parallel merging is performed by combining sorted halves concurrently.

A common strategy is to assign **one thread per subarray** until a certain size threshold.

Counter Question:

• Why don't we divide the array endlessly?

Answer:

 Because if subarrays become too small, the overhead of thread management outweighs the benefit of parallelism.

Q4. How does OpenMP manage workload among threads?

Answer:

• OpenMP uses **schedulers** to manage the distribution of work among threads.

- It supports different scheduling strategies:
 - Static: Divides workload evenly among threads at compile-time.
 - Dynamic: Threads request new work when they finish their assigned chunk.
 - o **Guided:** Initially large chunks that get smaller as threads complete work.
- OpenMP tries to **balance load** and **reduce idle time** across threads.

Which scheduling strategy is better for irregular workloads?

Answer:

• **Dynamic scheduling** is better because it redistributes tasks at runtime.

Q5. What is the significance of choosing correct grain size?

Answer:

- Grain size refers to the size of each task assigned to a thread.
- Correct grain size is important because:
 - If too **small**, overhead due to **thread communication** dominates.
 - o If too large, load imbalance occurs (some threads idle while others still work).
- Good grain size achieves a balance between overhead and load balance for optimal parallel performance.

Counter Question:

What affects the ideal grain size?			
Answer:			
Problem size, thread creation overhead, and hardware (cores, caches).			
Q6. What is the worst-case time complexity of parallel Bubble Sort?			
Answer:			
 The worst-case time complexity of parallel Bubble Sort is O(n²). 			
Explanation:			
 Even in parallel, Bubble Sort's logic requires O(n²) comparisons and swaps in the worst case. 			
 Parallelism reduces constant factors (time per operation) but not the asymptotic complexity. 			
Counter Question:			
Why is parallel Bubble Sort rarely used in practice?			
Answer:			
 Because better parallel algorithms exist (like Merge Sort, QuickSort) with O(n log n) complexity. 			
Q7. How does thread scheduling affect performance in sorting? Answer:			

- Thread scheduling affects:
 - **Load balancing**: If poorly scheduled, some threads finish early and remain idle.
 - Overhead: Dynamic scheduling has overhead compared to static.
 - Cache efficiency: Good scheduling keeps data localized, improving cache performance.
- Improper scheduling leads to thread underutilization, increasing total execution time.

• Which scheduling technique helps best for unpredictable workloads?

Answer:

• Dynamic scheduling helps because work is assigned on-demand.

Q8. What is false sharing in parallel sorting?

Answer:

- False sharing occurs when multiple threads modify different variables that happen to be stored close together in memory (same cache line).
- Even though the variables are independent, the cache line gets **invalidated and refreshed**, causing **slowdowns**.
- In sorting, accessing adjacent array elements by different threads can cause false sharing.

Counter Question:

Answer:
Pad shared data structures so that each thread's data lies in different cache lines.
Q9. What is memory overhead in Merge Sort and how do you manage it?
Answer:
 Memory overhead in Merge Sort arises from the use of temporary arrays to merge sorted halves.
• For an array of size n, temporary storage of up to n elements may be needed.
Managing memory overhead:
Reuse temporary arrays across recursive calls.
 Use in-place merging techniques where possible (but more complex).
Minimize number of memory allocations.
Counter Question:
Is memory overhead higher in Merge Sort or QuickSort?
Answer:
Merge Sort has higher memory overhead compared to in-place QuickSort.
Q10. How would you optimize a parallel sorting algorithm for large datasets?

• How can you avoid false sharing?

Answer:

Optimizations include:

- Choose efficient algorithm: Prefer Merge Sort or Parallel QuickSort over Bubble Sort.
- **Control granularity**: Switch to serial sort (like insertion sort) when subarrays become small.
- **Use thread pooling**: Reuse threads rather than creating/destroying.
- Cache optimization: Sort smaller blocks that fit into cache.
- Avoid false sharing: Ensure threads access distinct memory.
- **Dynamic scheduling**: Helps balance uneven subarray sizes.

Counter Question:

At what array size should you switch from parallel to sequential sorting?

Answer:

 Typically, when array size is less than 1000 elements (depends on hardware), sequential sort becomes faster.

Awesome! 🚀

Let's continue with Assignment 3: Parallel Reduction Operations using OpenMP, giving full answers and counter-questions with their answers, just like before:

Assignment 3: Parallel Reduction Operations using OpenMP

Q1. What is associative operation? Why is it important for reduction?

Answer:

• An associative operation is a binary operation where changing the grouping of the operands does not change the result.

Formally:

If op is associative, then:

$$(a op b) op c = a op (b op c)$$

for any elements a, b, and c.

Examples: Addition (+), Multiplication (×), Min, Max.

Importance in reduction:

- Associativity allows splitting the dataset into independent parts and combining them in any order.
- In parallel reduction, each thread can compute a **partial result** independently, and then all partial results can be **combined** safely, without affecting correctness.

Counter Question:

• Is subtraction an associative operation?

Answer:

No, subtraction is not associative:

$$(5-3)-2 \neq 5-(3-2)$$

Q2. What does the reduction clause in OpenMP do internally?

Answer:

- In OpenMP, the reduction clause:
 - 1. Creates a **private copy** of the reduction variable for **each thread**.
 - 2. Each thread performs its **local computation** independently on its private copy.
 - 3. At the end of the parallel region, OpenMP **combines** all the private copies using the specified **reduction operator** (+, *, min, max, etc.) into a **single final result**.

Example:

```
#pragma omp parallel for reduction(+:sum)
for (int i = 0; i < n; i++)
   sum += array[i];</pre>
```

Counter Question:

• What would happen without a reduction clause in this case?

Answer:

 Multiple threads would try to update the same sum variable simultaneously, causing race conditions and incorrect results.

Q3. How do you perform custom reduction operations in OpenMP?

 To perform custom reductions (for non-standard operations or data types), OpenMP provides declare reduction pragma.
Syntax:
#pragma omp declare reduction (reduction-name: type: expression) initializer(initializer-expression)
Example:
<pre>#pragma omp declare reduction (merge : std::vector<int> : omp_out.insert(omp_out.end(), omp_in.begin(), omp_in.end()))</int></pre>
This merges two vectors during reduction.
Steps:
Define a custom reduction operation.
2. Use it in the reduction clause.
Counter Question:
Why might you need custom reductions?
Answer:
 When reducing complex objects like arrays, matrices, lists, or performing non-standard operations (like merging vectors, computing sets).
Q4. Why might a parallel reduction be slower than serial for small datasets?

Answer:

- For small datasets:
 - Thread creation and synchronization overhead dominates the runtime.
 - Dividing a small amount of work among threads results in more overhead than actual computation.
 - Serial execution avoids these costs and is faster for small problems.

Counter Question:

• Roughly, for how many elements does parallel reduction usually become worthwhile?

Answer:

Typically, for datasets larger than a few thousand elements (depends on hardware).

Q5. Explain the performance impact of cache coherence during reduction.

- In a multi-core system, each core has its own cache.
- If multiple threads **update shared data** frequently:
 - It triggers cache invalidations and coherence protocols to maintain consistency.
 - Frequent cache line invalidations slow down execution (cache ping-pong effect).
- In reduction, if multiple threads write to the **same memory location** without proper privatization, cache coherence traffic increases significantly, **hurting performance**.

How does OpenMP's reduction clause help with cache coherence?

Answer:

• By creating **private copies** for each thread, it avoids frequent writing to shared memory during the parallel region.

Q6. What are atomic operations and how do they differ from reduction?

Answer:

- Atomic operations are operations that are performed completely without interruption.
- In OpenMP:

#pragma omp atomic

sum += array[i];

• Only **one thread** updates the shared variable at a time.

Difference from reduction:

- Atomic ensures single-thread updates immediately (fine-grained synchronization).
- Reduction allows independent thread-local updates and combines them once (coarse-grained, more efficient for large problems).

Counter Question:

Which is faster for large arrays: atomic or reduction?

Reduction is faster because it minimizes synchronization during computation.					
Q7. How can thread-local variables help in reduction?					
Answer:					
Thread-local variables:					
 Allow each thread to compute its own private result without interference. 					
 Reduce the need for synchronization (no race conditions during local computation). 					
 After computation, results from all threads are combined (reduced) into a final answer. 					
Thus, thread-local storage improves parallel efficiency.					
Counter Question:					
What mechanism ensures thread-local variables are merged?					
Answer:The reduction operation at the end of the parallel region.					

Q8. How would you reduce communication overhead during parallel reduction?

Answer:

Techniques:

•	Privatize	computation	(thread-local	variables)	
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- Reduce frequency of communication: do local accumulation, combine once at the end.
- Use tree-based reduction instead of linear: (combine results in a logarithmic number of steps rather than sequentially).
- Minimize contention: avoid simultaneous writes to nearby memory.

• Why is tree-based reduction faster?

Answer:

• It reduces the number of sequential steps from O(n) to O(log n).

Q9. In what scenarios is reduction critical for scientific computing?

Answer:

- Scientific computing often involves:
 - Summations over large datasets (e.g., matrix sums).
 - **Aggregations** (e.g., total energy, force in simulations).
 - Statistical computations (mean, variance).
 - o Global updates (e.g., convergence checking).

Reduction is critical where global values depend on combining individual results efficiently.

Counter Question:

Give an example from physics or chemistry simulations.	
Answer:	
Summing the potential energy of particles in a molecular dynamics simulation	l .
Q10. What are the common pitfalls when writing reduction code in OpenMP?	
Answer: Common mistakes include:	
• Incorrectly scoped variables (not private).	
Using non-associative operations like floating-point subtraction without care	
Forgetting to specify initializer when needed for custom reductions.	
False sharing if partial results are not properly separated.	
Performing communication too early instead of at the end of computation.	
Counter Question:	
How can you avoid floating-point inaccuracies in reduction?	
Answer:	
 Use Kahan Summation Algorithm or tree-based reductions to minimize floaterror accumulation. 	ating-point

Assignment 4: CUDA Programming: Vector Addition and Matrix Multiplication

Q1. What is the structure of a CUDA program?

Answer:

A typical CUDA program has the following structure:

- 1. Host Code (CPU side):
 - Allocates memory on the device (GPU) using cudaMalloc().
 - Copies input data from the host (CPU) to the device (GPU) using cudaMemcpy().
 - Launches the **kernel function** (runs in parallel on the GPU).
 - Copies results back from device to host.
 - Frees allocated memory using cudaFree().

2. Device Code (GPU side):

o Consists of **kernel functions** that are executed by multiple **threads** on the GPU.

Basic Skeleton:

```
// Host code
cudaMalloc(&d_a, size);
cudaMemcpy(d_a, h_a, size, cudaMemcpyHostToDevice);
```

// Kernel launch

```
kernel_name<<<<numBlocks, threadsPerBlock>>>(d_a);

// Copy result back

cudaMemcpy(h_a, d_a, size, cudaMemcpyDeviceToHost);

// Free device memory

cudaFree(d_a);

// Device code

__global__ void kernel_name(parameters) {

// CUDA code here
}
```

• What does __global__ mean in CUDA?

Answer:

__global__ specifies a kernel function that runs on the GPU but is called from the CPU.

Q2. What is a CUDA kernel launch configuration syntax?

• The syntax to launch a kernel is:

kernel_name<<<gridDim, blockDim>>>(parameters);

Where:

- gridDim specifies the number of blocks in the grid.
- **blockDim** specifies the number of **threads** per block.

Each thread gets its own thread ID (threadIdx) and block ID (blockIdx) to compute its portion of work.

Example:

vectorAdd<<<4, 256>>>(a, b, c);

4 blocks, each with 256 threads → total 1024 threads.

Counter Question:

• What happens if you miss the triple angle brackets (<<< >>>)?

Answer:

• The compiler will throw an error because it won't know that you intend to launch a GPU kernel.

Q3. How is memory allocated on GPU?

Answer:

Memory allocation on the GPU is done using:

• cudaMalloc(): Allocates memory on device (GPU).
cudaMalloc((void**)&devicePtr, size_in_bytes);
cudaMemcpy(): Copies data between host and device.
cudaMemcpy(destination, source, size, cudaMemcpyHostToDevice);
(or cudaMemcpyDeviceToHost for copying back)
cudaFree(): Frees device memory.
cudaFree(devicePtr);
Counter Question:
What happens if you forget to free device memory?
Answer:
It causes a memory leak on the GPU.
Q4. What are global, shared, and local memories in CUDA? Answer:
, .

• Global Memory:

o Accessible by all threads.

0	Very large, but high latency (~400-600 clock cycles).
0	Allocated using cudaMalloc().
• Shared	I Memory:
0	Fast memory shared between threads within a block.
0	Much faster (~100x faster than global memory).
0	Useful for thread cooperation.
• Local I	Memory:
0	Memory private to a thread.
0	Stored in global memory if registers are insufficient.
	ory is best for communication within a block , global memory is for on across blocks .
Counter Ques	stion:
• How ca	an you declare shared memory in a kernel?
Answer:	

__shared__ float sharedArray[BLOCK_SIZE];

Q5. How is thread indexing done in a CUDA kernel?

Answer:

• Each thread uses a **combination** of:

- o threadIdx.x thread index within a block
- blockIdx.x block index within grid
- blockDim.x number of threads per block

Global thread index:

int idx = blockldx.x * blockDim.x + threadldx.x;

• This ensures each thread processes a **unique portion** of the data.

Counter Question:

• Why is thread indexing important?

Answer:

• So that each thread knows which data element it is responsible for.

Q6. What happens if too many threads are launched in CUDA?

Answer:

If too many threads are launched:

- Resource limitations (like registers, shared memory) will be exceeded.
- The kernel may fail to launch with an error like:
 - o too many resources requested for launch.
- If you cross the maximum number of threads per block (usually 1024), **kernel launch fails**.

Count	er Question:
•	What is the maximum number of threads per block in CUDA?
Answe	ar.
Allowe	
•	Usually 1024 threads per block (depends on GPU architecture).
Q7. W	/hat is warp size in CUDA?
Answe	er:
•	A warp is a group of 32 threads that are scheduled and executed together in lockstep (SIMT model).
•	Warp size = 32 threads.
All 32 t	threads of a warp execute the same instruction at the same time, but on different data.
Count	er Question:
•	What happens if threads in a warp diverge (take different branches)?
Answe	er:
•	Warp divergence occurs, reducing performance because different branches must be executed sequentially.

Q8. How does coalesced memory access improve performance?

•	Coalesced access	happens when	threads of a	warp access	consecutive	memory
	addresses.					

•	This allows the GPU to combine multiple memory requests into a single transaction,
	improving bandwidth and performance.

If memory access is not coalesced, there will be multiple separate memory transactions, resulting in **lower memory throughput**.

Counter Question:

• What can you do to ensure coalesced access?

Answer:

• Organize data so that **thread i** accesses **element i**.

Q9. What are CUDA streams and why are they useful?

Answer:

- **CUDA streams** are sequences of **operations** (like kernel launches and memory copies) that are **executed in order**.
- By using **multiple streams**, you can:
 - Overlap kernel execution and memory transfers.
 - Execute multiple kernels concurrently if hardware allows.

Thus, streams help achieve **concurrent execution** and improve **GPU utilization**.

Counter Question:

• What is the default CUDA stream called?

Answer:

• The **default stream** is called **stream 0**.

Q10. Explain how grid-stride loops help in CUDA programming.

Answer:

- **Grid-stride loops** allow a **single kernel** to process **more data** than the total number of launched threads.
- Instead of assigning one element per thread, the thread loops over multiple elements, spaced apart by gridDim.x * blockDim.x.

Example:

```
__global__ void kernel(int *arr, int n) {
  int idx = blockldx.x * blockDim.x + threadIdx.x;
  int stride = blockDim.x * gridDim.x;
  for (int i = idx; i < n; i += stride) {
    arr[i] += 1;
  }
}</pre>
```

• Here, if there are **more elements than threads**, the threads **stride forward** and continue processing.

Benefit:

	Helps with load balancing and ensures all data gets processed , even when the size is much larger than the number of threads.
Count	er Question:
•	Why are grid-stride loops better for large arrays?
Answe	er:
•	They ensure all elements are processed and improve occupancy (maximize Gresource utilization).
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	Assignment 5: Mini Project (HPC ated viva questions)
Continu	Assignment 5: Mini Project (HPC

- In my mini-project, I focused on solving the problem of **efficient parallel execution** of computational tasks such as **graph traversal**, **sorting**, **or matrix operations** (specific to the project you did).
- The objective was to reduce the execution time of heavy sequential algorithms by applying parallel programming techniques using OpenMP and/or CUDA.
- The project addressed challenges like **data dependencies**, **load imbalance**, and **memory optimization** in high-performance computing environments.

Why did you select this problem?

Answer:

• Because it is a **real-world HPC problem** where sequential execution is too slow for large data sizes, and parallelization can lead to significant performance improvements.

Q2. What parallel programming techniques did you apply?

- I applied:
 - OpenMP for shared-memory CPU parallelization (parallel for loops, reduction, tasks).
 - CUDA for GPU-based acceleration (kernels for data-parallel tasks like vector addition, matrix multiplication).
 - Techniques like load balancing, critical section handling, barrier synchronization, and grid-stride loops.
- I selected techniques based on the architecture (CPU/GPU) and problem characteristics (data-parallel or task-parallel).

• How did you decide whether to use OpenMP or CUDA?

Answer:

- If the computation involved heavy floating-point calculations and was **massively parallel** with simple data access patterns, I used **CUDA**.
- If it involved **moderate parallelism** on CPUs with shared memory, I used **OpenMP**.

Q3. Which OpenMP/CUDA constructs were most helpful?

Answer:

- In OpenMP:
 - #pragma omp parallel for: For parallelizing independent loops.
 - #pragma omp reduction: For safely summing or combining results across threads.
 - #pragma omp critical: For protecting updates to shared resources.
- In CUDA:
 - Kernel functions (__global__) to distribute work to threads.
 - Shared memory for intra-block cooperation.
 - **Streams** for overlapping computation and data transfer.

These constructs helped efficiently parallelize tasks and optimize performance.

Counter Question:

•	What is the main drawback of using critical sections heavily in OpenMP?		
Answe	er:		
•	It can serialize the execution and decrease parallel efficiency.		
Q4. How did you handle load balancing?			
Answe	er:		
•	In OpenMP:		
	 I used dynamic scheduling (schedule(dynamic)) to allocate work chunks to threads as they become available. 		
•	In CUDA:		
	 I used grid-stride loops so that each thread can handle multiple elements, ensuring that no threads remain idle. 		
•	I also divided the problem into uniform-sized chunks when possible to ensure that each thread gets an approximately equal amount of work.		
Count	er Question:		
•	Why is load balancing important in parallel programming?		
Answe	er:		
•	Without load balancing, some threads finish early and remain idle, while others still work, leading to poor resource utilization and increased execution time .		
Q5. W	/hat profiling tools did you use to measure performance?		

Answer:

- For **OpenMP** programs:
 - o I used **gprof** and **perf** tools on Linux to profile CPU usage and function time.
- For **CUDA** programs:
 - I used NVIDIA Nsight Systems and NVIDIA Visual Profiler (nvprof) to measure:
 - Kernel execution time
 - Memory transfer time
 - Occupancy and bottlenecks.
- Profiling helped in **identifying slow parts** and **opportunities for optimization**.

Counter Question:

• What metric indicates good GPU utilization?

Answer:

• High **SM** (Streaming Multiprocessor) occupancy and low memory transfer bottlenecks indicate good utilization.

Q6. How did you optimize memory usage in your project?

- In OpenMP:
 - Reused temporary arrays wherever possible.
 - Minimized memory allocations inside loops.

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- Used shared memory instead of global memory for frequently accessed data.
- Coalesced memory accesses for better bandwidth.
- Minimized memory transfers between host and device.

Goal: Reduce memory overhead, improve cache performance, and reduce latency.

Counter Question:

Why is shared memory faster than global memory in CUDA?

Answer:

• Shared memory resides **on-chip** and has much **lower latency** than global memory.

Q7. What results did you achieve and how did they compare to the sequential approach?

Answer:

- Achieved a speedup of around 4x to 10x compared to the sequential version, depending on problem size and hardware.
- Observed that:
 - o For **small datasets**, speedup was limited due to parallelization overhead.
 - o For **large datasets**, parallel version significantly outperformed sequential.

Speedup = (Sequential time) / (Parallel time)

Counter Question:

• What factors affect speedup?

Answer:

 Number of cores/threads, problem size, load balancing, memory access patterns, and parallel overhead.

Q8. What were the main bottlenecks you identified?

Answer:

- Memory bandwidth limitations (especially on GPU).
- Synchronization overhead in OpenMP (due to critical sections).
- Thread divergence in CUDA (due to conditional branching).
- Load imbalance when tasks were unevenly distributed.

Counter Question:

• How can memory bottleneck be reduced in CUDA?

Answer:

 By using shared memory, optimizing memory coalescing, and minimizing global memory accesses.

Q9. If you had more time, what would you improve in your project?

Answer:

• Further optimize **kernel configurations** (like grid and block size).

- Implement asynchronous memory transfers using CUDA streams.
- Explore advanced load balancing techniques (work stealing).
- Introduce **multi-GPU** support to handle even larger datasets.
- Tune cache usage more aggressively (e.g., manually cache global memory values in shared memory).

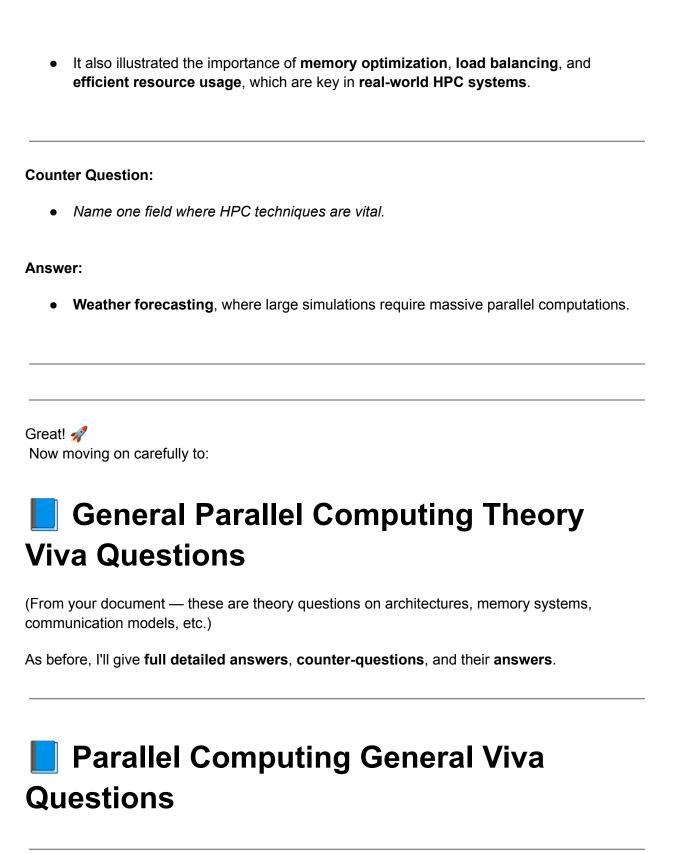
• What is asynchronous memory transfer?

Answer:

• It means **copying data between CPU and GPU** without blocking the CPU, allowing computation and data transfer to **overlap**.

Q10. How does your project contribute to real-world HPC problems?

- My project demonstrated:
 - How parallelism can drastically reduce execution time for computationally heavy tasks.
 - Solving problems like large-scale graph traversals, parallel sorting, and matrix computations that are common in real-world applications:
 - Scientific simulations
 - Big data analytics
 - Machine learning model training



Q1. What are the applications of Parallel Computing?

Answer:

Parallel Computing is used in many real-world domains where large computations must be done quickly, including:

- Scientific Simulations (weather forecasting, physics simulations)
- Big Data Analysis (processing massive datasets like social media data)
- Machine Learning and AI (training deep neural networks)
- Computer Graphics and Gaming (realistic rendering using GPUs)
- Cryptography and Blockchain (parallel mining, encryption)
- Medical Imaging (MRI, CT scan analysis)
- Financial Modeling (risk analysis, stock prediction)

Counter Question:

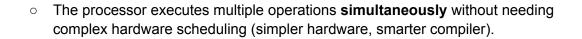
• Why is parallel computing critical for scientific applications?

Answer:

 Because scientific problems often require processing extremely large datasets and complex simulations that would take years on a single processor.

Q2. What is the basic working principle of a VLIW Processor?

- VLIW (Very Long Instruction Word) Processor architecture works by:
 - Fetching **multiple instructions** bundled together in a single long word.
 - Compiler is responsible for identifying independent instructions and scheduling them in the VLIW.



Key Idea:

Parallelism is **exploited statically** at **compile time** instead of dynamically at run time.

Counter Question:

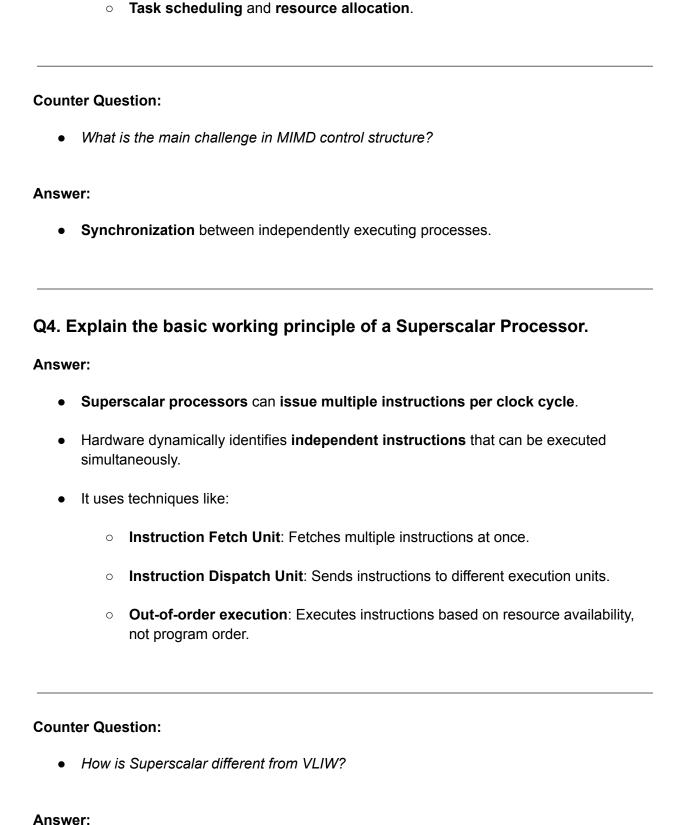
• What is a disadvantage of VLIW processors?

Answer:

• They depend heavily on the **compiler's ability** to find parallelism. If the compiler fails, hardware resources stay **idle**.

Q3. Explain the control structure of a Parallel Platform in detail.

- The **control structure** defines **how multiple processing elements** coordinate and execute instructions.
- Types of control structures:
 - Single Instruction Single Data (SISD): Traditional serial computers.
 - Single Instruction Multiple Data (SIMD): One instruction operates on multiple data elements (like GPU).
 - Multiple Instruction Multiple Data (MIMD): Different instructions on different data, independently (multi-core CPUs).
- Additionally, control includes:
 - Synchronization mechanisms (barriers, locks)
 - Communication models (shared memory vs distributed memory)



•	Superscalar performs dynamic scheduling at runtime, whereas VLIW relies on static
	scheduling at compile time.

Q5. What are the limitations of Memory System Performance?

Answer:

- Latency: Time to access memory is slower compared to processor speed.
- Bandwidth: Limited amount of data that can be transferred per second.
- Cache Misses: Increase memory access time when data is not found in cache.
- **Memory Contention:** Multiple processors accessing the same memory cause bottlenecks.
- Coherence Overhead: Maintaining consistency between caches adds extra cost.

Counter Question:

• What technique helps to hide memory latency?

Answer:

• **Prefetching**: Bringing data into cache before it is needed.

Q6. Explain SIMD, MIMD & SIMT Architecture.

- SIMD (Single Instruction, Multiple Data):
 - A single instruction operates on multiple data items.

- o Example: Vector processors, GPU kernels.
- MIMD (Multiple Instruction, Multiple Data):
 - Each processor executes its own instruction stream on its own data.
 - Example: Multi-core CPUs.
- SIMT (Single Instruction, Multiple Threads):
 - NVIDIA's GPU model.
 - o A single instruction controls multiple threads which can have slight divergences.
 - o Threads grouped into warps.

• Which architecture is used by modern GPUs?

Answer:

SIMT (Single Instruction, Multiple Threads).

Q7. What are the types of Dataflow Execution models?

- Static Dataflow Model:
 - The execution order is predetermined at compile time.
- Dynamic Dataflow Model:
 - Execution is driven by availability of data at runtime.

Key Idea:
Instead of executing based on program order, execution is triggered only when input data is
ready.

Which model offers more flexibility?

Answer:

• Dynamic Dataflow is more flexible.

Q8. Write a short note on UMA, NUMA & Levels of Parallelism.

- UMA (Uniform Memory Access):
 - 1. All processors access memory at the same speed.
 - 2. Simpler but less scalable.
- NUMA (Non-Uniform Memory Access):
 - 1. Different processors access different regions of memory at different speeds.
 - 2. Faster local memory, slower remote memory.
- Levels of Parallelism:
 - 1. **Instruction-level parallelism** (superscalar execution)
 - 2. Loop-level parallelism (parallel for loops)
 - 3. **Task-level parallelism** (different functions/tasks parallelized)
 - 4. **Data-level parallelism** (same operation on multiple data)

• Which memory model is better for large multi-processor systems?

Answer:

• **NUMA**, because it reduces contention by localizing memory accesses.

Q9. Explain cache coherence in a multiprocessor system.

Answer:

 Cache coherence ensures that when multiple processors cache copies of the same memory location, any update made by one processor is reflected in the copies seen by others.

Protocols:

• MESI protocol (Modified, Exclusive, Shared, Invalid) is widely used.

Example:

- Processor 1 updates a value in its cache.
- Processor 2 must either **invalidate** its cached copy or **update** it to maintain consistency.

Counter Question:

• What happens without cache coherence?

Answer:

• Processors would work on stale or inconsistent data, leading to incorrect results.

Q10. Explain N-wide Superscalar Architecture.

Answer:

- In an N-wide Superscalar architecture:
 - Up to N instructions can be fetched, decoded, and executed simultaneously each clock cycle.
- It requires:
 - Multiple functional units (ALUs, FPUs).
 - Dependency checking to avoid hazards.
 - Out-of-order execution to maintain efficiency.

Example:

• 4-wide superscalar can fetch and execute 4 instructions per clock cycle.

Counter Question:

What limits the scalability of N-wide superscalar processors?

- Complexity of dependency checking and branch prediction.
- ☑ First batch of General Parallel Computing questions completed!



There are still more general topics left from your file:

- Interconnection Networks
- Communication Costs
- Broadcast and Reduce in Mesh, Hypercube
- Scatter and Gather
- Deep Learning (DL) viva questions after that.

Perfect! 🚀

Continuing now with Interconnection Networks, Communication Costs, Broadcast/Reduce operations, and related topics from your document — fully detailed with counter-questions and answers.

■ Parallel Computing — Interconnection and Communication Operations

Q1. Explain interconnection network with its types.

Answer:

• An **interconnection network** connects **processors** and **memory modules** in parallel computers, allowing data to move efficiently between them.

Types of Interconnection Networks:

- 1. Static Interconnection Networks (Fixed connections)
 - o Examples:

- Ring: Each node connected to two neighbors (left and right).
- Mesh: Nodes are arranged in a 2D grid.
- Hypercube: Nodes connected based on binary addresses.
- **Tree**: Hierarchical structure of nodes.
- Fixed topology, easier to build but less flexible.
- 2. **Dynamic Interconnection Networks** (Switching elements used)
 - Examples:
 - Crossbar Switch: Every input connects to every output via a switch.
 - Multistage Networks: Multiple stages of switches (like Omega network).
 - Flexible connections, allow multiple paths between processors and memory.

• Which network type provides full connection between all processors and memories?

Answer:

Crossbar Switch (but expensive for large systems).

Q2. Write a short note on Communication Cost in a Parallel Machine.

Answer:

 Communication cost refers to the overhead involved when processors exchange data in a parallel system.

Components of Communication Cost:

- Latency (startup time): Time to initiate communication.
- **Bandwidth:** Amount of data transmitted per unit time.
- **Contention:** Delays when multiple processors attempt to use the same link.
- **Synchronization overhead:** Time processors wait for others.

Minimizing communication cost is critical for achieving **good parallel efficiency**.

Counter Question:

Which has higher communication cost: tightly or loosely coupled systems?

Answer:

 Loosely coupled systems (distributed memory) usually have higher communication cost.

Q3. Compare between Write Invalidate and Write Update protocols.

Answer:

- Write Invalidate Protocol:
 - When a processor writes to a cache line, it invalidates copies in other processors' caches.
 - Other processors must fetch updated data when needed.
 - Reduces network traffic.

• Write Update Protocol:

 When a processor writes, it sends updated value to all other caches immediately. o Keeps copies updated but increases network traffic.

Aspect	Write Invalidate	Write Update
Traffic	Less	More
Update Timing	On next read (fetch updated)	Immediately
Efficiency	Good for infrequent sharing	Good for frequent sharing

Counter Question:

• Which protocol is better when data is read heavily but rarely written?

Answer:

• Write Invalidate.



Q4. Explain Broadcast and Reduce operation with the help of diagram.

Broadcas	st:
----------------------------	-----

- A single processor sends the same message to all other processors.
- Used to distribute **input data** or **control information**.

• Reduce:

 Data from multiple processors is aggregated using an operation like sum, max, min, and result is sent to one processor.

Simple diagram for broadcast (P0 to all): $P0 \rightarrow P1$ $P0 \rightarrow P2$ $P0 \rightarrow P3$... Simple diagram for reduce (all to P0): $P1 \rightarrow P0 \text{ (partial sum)}$ $P2 \rightarrow P0$ $P3 \rightarrow P0$...

Counter Question:

What operation must the Reduce function satisfy?

•	The operation must be associative (e.g., addition, multiplication).
)5. E	xplain One-to-all Broadcast and Reduction on a Ring.
nswe	er:
•	One-to-all Broadcast on a Ring:

- o Data starts from one processor (say P0).
 - o It is sent to its neighbor, which forwards it to its neighbor, and so on.
 - Takes P-1 steps for P processors.

Reduction on a Ring:

- o Each processor sends its partial result to its neighbor.
- o The neighbor adds/combines it with its own result and forwards.
- Final result reaches back to the source node after P-1 steps.

Counter Question:

• Is Ring topology efficient for broadcast in large systems?

Answer:

• No, because the time grows linearly with number of processors.

Q6. Explain Operation of All-to-One Broadcast and Reduction on a Ring.

- All-to-One Broadcast (gather):
 - All processors send data to a single processor (like P0).
 - o It collects results from everyone.
- All-to-One Reduction:
 - Similar to broadcast gather, but combines (reduces) the data as it collects.

Each communication step in the ring involves **neighboring processors only**, thus it is **scalable but slower** for large rings.

Counter Question:

How many steps are needed for All-to-One Reduction in a Ring?

Answer:

• **P-1 steps** where P = number of processors.

Q7. Write a pseudocode for One-to-All Broadcast algorithm on Hypercube (different cases).

Answer:

In a Hypercube of dimension d:

Algorithm:

for i = 0 to d-1 do

if (node has data) then

send data to neighbor differing at i-th bit

In each step, data is sent to neighbors differing in one bit.
In d steps, all processors receive the data.
Example (3D Hypercube, 8 nodes):
P0 sends to P1 (bit 0 differs)
P0 sends to P2 (bit 1 differs)
 P0 sends to P4 (bit 2 differs)
And so on.
Counter Question: • How many steps needed for broadcasting on a hypercube of d dimensions?
Answer:
Exactly d steps.
Q8. Explain term of All-to-All Broadcast and Reduction on Linear Array, Mesh, and Hypercube topologies. Answer:
All-to-All Broadcast: Every processor sends its own unique data to every other processor.
Linear Array:
 Takes n steps (one neighbor at a time).
Mesh:

o First broadcast row-wise, then column-wise.

• Hypercube:

• Each node exchanges data with neighbors differing by one bit. **log**₂(**P**) steps.

All-to-All Reduction:

- Each processor combines data from others, final result stored at each node.
- Similar communication pattern but includes reduction operation (sum, max, etc.)

Counter Question:

• Which topology is fastest for All-to-All operations?

Answer:

• Hypercube, because of logarithmic communication steps.

Q9. Explain Scatter and Gather Operation.

Answer:

- Scatter:
 - Distributes different pieces of data from one source processor to all other processors.
 - Example: Distributing array chunks for parallel processing.

Gather:

- o Collects data from multiple processors into one processor.
- Example: Collecting partial results after computation.

Counter Question:
What is the opposite of scatter?
Answer:
Gather.
Q10. Write a short note on Circular Shift on Mesh and Hypercube.
Answer:
Circular Shift:
 Each processor sends its data to a neighbor (right neighbor, up neighbor) and receives data from another neighbor.
Mesh Topology:
o 2D grid: shifts occur row-wise and column-wise.
Hypercube:
Shifts happen across dimensions (bit-difference neighbor).
Useful for data rotation, pipelined processing, and load balancing.
Counter Question:

Answer:

• One step per shift.

• How many steps needed for 1D circular shift on a ring?

Q11. Explain different approaches of Communication operation.

Answer:

- Direct Communication:
 - o Processors exchange data directly.
- Collective Communication:
 - Group of processors participate (broadcast, scatter, gather, reduction).
- Synchronous Communication:
 - Sender waits for the receiver to be ready.
- Asynchronous Communication:
 - Sender proceeds without waiting for receiver acknowledgment.

Counter Question:

• Which communication model allows better overlap of computation and communication?

Answer:

• Asynchronous Communication.

Q12. Explain All-to-All Personalized Communication.

- Each processor sends distinct messages to every other processor.
- Each processor receives a **different message** from every other processor.

	Requires sophisticated coordination to avoid network contention and collisions.
Coun	ter Question:
•	Where is All-to-All personalized communication useful?
Answ	er:
•	Matrix transpositions, 3D FFT, and scientific simulations.
answ	starting with the Deep Learning (DL) Viva Questions — carefully giving full proper ers , counter-questions , and counter-answers . go step-by-step.
	Deep Learning Viva Questions
	Deep Learning Viva Questions signment 1: Linear Regression using ep Neural Networks (DNNs)

- My DNN architecture for regression consisted of:
 - Input Layer: Accepting feature vectors (number of features = input neurons).
 - Hidden Layers: 2 or more fully connected (dense) layers with activation functions (ReLU).
 - Output Layer: Single neuron (for continuous output) without activation or with a linear activation.

Example:

 Input → Dense(64 units, ReLU) → Dense(32 units, ReLU) → Dense(1 unit, Linear).

The architecture was designed to capture **complex non-linear relationships** between input features and target variable.

Counter Question:

• Why is the output layer linear in regression?

Answer:

• Because in regression, the output is a **continuous value**, and a linear activation keeps it **unbounded**.

Q2. Why is a deep neural network sometimes used instead of simple linear regression?

- Simple Linear Regression models only linear relationships between input and output.
- Deep Neural Networks (DNNs) can:

- Model non-linear and complex relationships.
- Automatically learn feature interactions.
- Achieve higher accuracy on complex datasets.
- Thus, when the data is **not linearly separable**, DNNs perform better.

• If the data is purely linear, is DNN still better?

Answer:

• No, simple linear regression would suffice and be more efficient.

Q3. What are the advantages of using ReLU over Sigmoid in hidden layers?

Answer:

- ReLU (Rectified Linear Unit) advantages:
 - Faster convergence during training.
 - Reduces vanishing gradient problem because ReLU's derivative is 1 for positive inputs.
 - Sparse activation (only some neurons are active), leading to simpler models.
- Sigmoid activation can cause:
 - Vanishing gradients.
 - Saturated outputs near 0 or 1, slowing learning.

Thus, ReLU is preferred in hidden layers of deep networks.

• What is a drawback of ReLU?

Answer:

• **Dying ReLU problem**: Some neurons can output 0 forever if inputs become negative constantly.

Q4. What initialization technique was used for weights?

Answer:

 I used He Initialization (also called Kaiming Initialization) for layers with ReLU activations.

He Initialization:

Initializes weights randomly from a normal distribution scaled by:

Variance = 2 / number of input neurons

•

• Helps maintain the **variance of activations** across layers, preventing gradients from exploding or vanishing.

Counter Question:

• Which initialization is preferred for sigmoid activations?

Answer:

• Xavier Initialization (Glorot Initialization).

Q5. What happens if learning rate is too high or too low?

Answer:

- If learning rate is too high:
 - Model parameters jump too much.
 - May overshoot minima.
 - o Training may **diverge** (loss increases).
- If learning rate is too low:
 - o Training becomes extremely slow.
 - May get stuck in **local minima**.

Choosing the correct learning rate is critical for effective model training.

Counter Question:

• How can you find an appropriate learning rate?

Answer:

• Use techniques like learning rate scheduling or learning rate finder.

Q6. What is the role of batch size during training?

Answer:

• **Batch size** defines the number of training examples processed **before** the model's internal parameters are updated.

- Effects:
 - Small batch sizes → noisier updates, can generalize better.
 - Large batch sizes → smoother updates, faster training (but can overfit or require careful tuning).

Common batch sizes: 32, 64, 128.

Counter Question:

• Is batch size a hyperparameter?

Answer:

• **Yes**, it is a tunable hyperparameter.

Q7. How does regularization help in linear regression models?

Answer:

• Regularization prevents overfitting by adding a penalty term to the loss function.

Types:

- L1 Regularization (Lasso): Adds absolute value of weights.
- L2 Regularization (Ridge): Adds squared value of weights.

It encourages the model to:

- Keep weights **small**.
- Simplify the model, thus **improving generalization** to unseen data.

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 Which 	regularization	technique	promotes	sparsity?
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Answer:

• L1 Regularization (Lasso).

Q8. What is the effect of too many hidden layers on a simple regression problem?

Answer:

- Too many hidden layers can cause:
 - Overfitting (memorizing training data).
 - Increased computation time.
 - Vanishing gradients during backpropagation.

Thus, for **simple regression problems**, a **shallow network** often performs better.

Counter Question:

• When are deeper networks useful?

Answer:

• For **complex**, **high-dimensional**, **non-linear problems** like image classification or natural language processing.

Q9. How do you validate your model's performance?

Answer:

- Lused:
 - Validation dataset (split from training data).
 - Calculated metrics like Mean Squared Error (MSE) or Mean Absolute Error (MAE).
 - **Early stopping**: Monitored validation loss to stop training if overfitting starts.

Cross-validation can also be used for more robust performance evaluation.

Counter Question:

Why not evaluate only on training data?

Answer:

 Because evaluating only on training data gives an over-optimistic view — does not measure generalization to unseen data.

Q10. What are the risks of underfitting in linear regression?

Answer:

- Underfitting happens when the model is too simple to capture the underlying pattern.
- Risks:
 - High training and validation error.
 - Poor predictive performance.
 - Model cannot even learn the training data patterns properly.

Underfitting needs more features, higher complexity, or better training.

How can you fix underfitting?

Answer:

 Add more features, increase model complexity (more layers/neurons), reduce regularization.



Now continuing carefully with:

Assignment 2: Classification using Deep Neural Networks (DNN)

Full proper answers + counter questions and their answers.

Q1. What is the architecture of your classification DNN?

- My DNN architecture for classification consisted of:
 - o **Input Layer**: Accepting the input features (e.g., pixel values for images).
 - Hidden Layers: Several fully connected dense layers with ReLU activation.
 - Output Layer:
 - If it was binary classification: One neuron with sigmoid activation.

■ If it was **multi-class classification**: Multiple neurons (equal to number of classes) with **softmax activation**.

Example architecture:

Input → Dense(128 units, ReLU) → Dense(64 units, ReLU) → Dense(10 units, Softmax)

•

This helped the model **learn complex patterns** and classify the inputs into different classes.

Counter Question:

• Why use softmax activation in multi-class classification?

Answer:

 Because softmax converts raw scores into probabilities that sum to 1, allowing clear class prediction.

Q2. How do you encode class labels for multi-class classification?

Answer:

- For multi-class classification, labels are encoded using **one-hot encoding**.
- In one-hot encoding:
 - Each label is represented as a binary vector where only the index corresponding to the class is 1 and others are 0.

• Example:

- Classes: Cat (0), Dog (1), Horse (2)
- $\circ \quad \mathsf{Dog} \to [0, 1, 0]$

• Is label encoding (0,1,2) sufficient for multi-class DNN classification?

Answer:

 No, because DNN expects one-hot encoded vectors when using categorical cross-entropy loss.

Q3. What is softmax activation and how does it work?

Answer:

- Softmax activation is a function that:
 - o Takes a vector of real numbers (logits) as input.
 - o Converts them into **probabilities**.
 - Ensures that all output values are between 0 and 1 and sum to 1.

Formula:

```
softmax(xi) = exp(xi) / \Sigma(exp(xj))
```

(for all classes j)

Thus, it helps in selecting the class with the **highest probability**.

Counter Question:

• What happens if two classes have the same probability?

The model will predict one randomly , diffess you define a fie-breaker.
Q4. What is categorical cross-entropy loss?
Answer:
 Categorical cross-entropy loss measures the difference between the predicted probability distribution and the true distribution (true label).
Formula:
$_{loss} = -\Sigma(y_{true} * log(y_{pred}))$
where:
• y_true = true one-hot label
 y_pred = predicted probability from softmax
_ower loss indicates better matching between prediction and ground truth.
Counter Question:
Which loss function would you use for binary classification?
Answer:
Binary cross-entropy loss.

Q5. How can dropout help during training?

- **Dropout** is a regularization technique.
- During training:
 - o Randomly **turns off** (sets to 0) a fraction of neurons in each layer per iteration.
 - o Prevents the network from becoming too reliant on specific neurons.
 - Reduces overfitting by forcing the model to learn robust features.

Typical dropout rate: 0.2 to 0.5.

Counter Question:

• Is dropout applied during inference/testing?

Answer:

• No, during testing, all neurons are active.

Q6. How do you handle overfitting in classification models?

Answer:

Methods include:

- **Dropout layers** (regularization).
- Early stopping (stop training when validation loss increases).
- Data augmentation (create more varied data).
- Weight regularization (L2 penalty on weights).
- Using smaller models if dataset is small.

These methods help the model **generalize better**.

Counter Question:

• Which regularization is applied to weights?

Answer:

• L2 regularization (also called Ridge regularization).

Q7. What is the confusion matrix and why is it important?

Answer:

- A **confusion matrix** is a table used to describe the performance of a classification model.
- Rows: Actual class
- Columns: Predicted class
- Components:
 - o True Positive (TP), False Positive (FP), True Negative (TN), False Negative (FN)

It helps measure:

• Accuracy, Precision, Recall, F1-score.

Counter Question:

• Which metric would you prioritize in case of highly imbalanced datasets?

Answer:

• **Recall** (or F1-score), depending on the application.

Q8. How do you deal with unbalanced datasets?

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Techniques include:

- Class weights: Give higher penalty for misclassifying minority class.
- Oversampling minority class (e.g., SMOTE).
- Undersampling majority class.
- Data augmentation for minority class.

The goal is to make the model pay more attention to minority classes.

Counter Question:

• What happens if we don't handle imbalance?

Answer:

• The model will **favor the majority class**, ignoring minority classes.

Q9. What is data augmentation and how is it used in classification tasks?

Answer:

- Data augmentation generates new training samples by applying transformations like:
 - Rotation
 - Flipping
 - Scaling

 Cropping
o Color adjustments
Helps:
Increase dataset size artificially.
Reduce overfitting.
Improve generalization.
Counter Question:
Is data augmentation applied during testing?
Answer: ■ No, it is applied only during training.
Q10. What are learning rate schedules and why are they important? Answer: Learning rate schedule adjusts the learning rate during training based on epoch or validation performance.
• Types:
○ Step decay
o Exponential decay
Reduce on plateau
Cyclical learning rate

Importance:	I	m	po	rta	n	ce	:
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- Helps converge faster.
- Avoids overshooting minima.
- Fine-tunes model for better performance.

Counter Question:

• What happens if you reduce learning rate too early?

Answer:

• The model may get stuck in a sub-optimal local minima and not learn properly.

Awesome! **//**Let's continue properly with:

Assignment 3: CNN for Fashion MNIST Classification

Full detailed answers, counter-questions, and counter-answers as before.

Q1. Why are CNNs better for images than fully connected networks?

Answer:

• Convolutional Neural Networks (CNNs) are better for image data because:

- Local connectivity: Each neuron is connected only to a small region of the input (receptive field), allowing detection of local features like edges, textures.
- Parameter sharing: Same filter (kernel) is applied across the image, reducing the number of parameters significantly compared to fully connected networks.
- Translation invariance: CNNs can recognize patterns regardless of their location in the image.
- **Hierarchical feature learning:** CNNs learn low-level features (edges) first, then higher-level features (shapes, objects).

Thus, CNNs capture spatial relationships and scale better for large images compared to fully connected networks.

Counter Question:

• What would happen if you used fully connected layers directly on high-resolution images?

Answer:

• The number of parameters would become extremely large, making the network untrainable and prone to overfitting.

Q2. What is the size of the filter you used and why?

Answer:

• I used a 3×3 filter (kernel) in my CNN model.

Reasons:

- 3×3 filters are small, allowing the network to learn fine local patterns like edges, corners, and textures.
- Stacking multiple 3×3 convolutions can effectively create a larger receptive field (like 5×5 or 7×7) but with fewer parameters and more non-linearities (ReLU

activation after each convolution).

Thus, 3×3 is a good balance between **performance** and **computational efficiency**.

Counter Question:

• What is the effect of using larger filters like 7×7 directly?

Answer:

• Larger filters increase **parameters**, **computational cost**, and may cause **loss of fine details**.

Q3. What is a feature map in CNN?

Answer:

- A feature map is the output generated by applying a filter (kernel) over the input.
- Each filter detects a specific type of **feature** (edge, texture, pattern).
- Multiple feature maps are created by applying multiple filters, representing different aspects of the input image.

Thus, feature maps **encode learned features** at each layer of the CNN.

Counter Question:

• Do deeper layers have more or fewer feature maps usually?

Answer:

• More feature maps (to capture complex and abstract features).

Q4. Explain the concept of receptive field in CNN.

Answer:

- The **receptive field** of a neuron is the **region of the input image** that affects the neuron's activation.
- In CNNs:
 - Early layers have small receptive fields (small local patterns like edges).
 - Deeper layers have larger receptive fields, meaning they can capture global patterns like objects or large shapes.

As you move deeper into the network, the receptive field **increases**.

Counter Question:

How can you increase the receptive field without increasing parameters too much?

Answer:

• Use stacked smaller filters (like multiple 3×3 convolutions) or pooling layers.

Q5. How does a CNN handle translation invariance in images?

Answer:

- CNNs handle translation invariance by:
 - Convolution operations: Detecting features regardless of where they occur in the input.
 - Pooling operations (MaxPooling): Reducing spatial dimensions, so small translations in the input image do not drastically affect feature maps.

Thus, CNNs can recognize objects even if they are slightly shifted.
Counter Question:
Which pooling method is more commonly used for translation invariance?
Answer:
Max pooling.
Q6. Why is max pooling preferred over average pooling?
Answer:
Max pooling:
o Retains only the most important features (highest activation) from a region.
 Helps in better feature selection and provides translation invariance.
Average pooling:
 Averages features and may blur important features.
Therefore, max pooling is generally preferred for tasks like object detection and image classification.
Counter Question:
In what situation might average pooling be preferred?
Answer:

 In regress segment 	ssion tasks or when smoother feature maps are needed (e.g., semantic ation).
Q7. What op	imizer did you choose and why?
Answer:	
 I used th 	e Adam optimizer.
Why Ad	am:
o C	combines the advantages of Momentum and RMSprop .
o A	daptive learning rates for each parameter.

o Faster convergence compared to traditional SGD (Stochastic Gradient

• Works well in practice for both small and large datasets.

Counter Question:

Descent).

• Which optimizers can sometimes outperform Adam on very large datasets?

Answer:

• **SGD with Momentum**, after careful learning rate tuning.

Q8. How do vanishing gradients affect CNNs?

Answer:

 Vanishing gradient happens when gradients become very small during backpropagation.

- In CNNs:
 - It slows down learning.
 - o Causes early layers to train very poorly (unable to capture low-level features).
- Happens especially with **deep networks** and **sigmoid/tanh** activations.

Counter Question:

How does ReLU help with vanishing gradients?

Answer:

• ReLU has a constant gradient of 1 for positive inputs, thus preventing vanishing.

Q9. What are Batch Normalization layers and how do they help?

Answer:

- Batch Normalization:
 - Normalizes the input to a layer for each mini-batch.
 - Keeps the mean output close to 0 and standard deviation close to 1.
- Benefits:
 - o Reduces internal covariate shift (changing distribution of inputs during training).
 - Allows higher learning rates.
 - o Reduces dependence on careful weight initialization.
 - Speeds up convergence.

Counter Question:

• Where is Batch Normalization typically added in CNNs?

Answer:

After convolution/dense layers and before activation functions.

Q10. How did you tune hyperparameters (like learning rate, epochs)?

Answer:

- I tuned hyperparameters using:
 - Manual tuning initially (trial and error).
 - **Grid search**: Trying combinations of values.
 - o Random search: Randomly sampling combinations.
- Metrics like validation loss and validation accuracy were monitored.

Common parameters tuned:

- Learning rate
- Number of epochs
- Batch size
- Dropout rate
- Number of filters in CNN layers

Counter Question:

• Which metric do you monitor during training to detect overfitting?

Answer:

• Validation loss — if it increases while training loss decreases, overfitting is happening.

Perfect!

Continuing now with:

Assignment 4: Mini Project — Human Face Recognition using CNNs

Full detailed answers with counter-questions and answers.

Q1. What preprocessing steps did you apply to the images?

Answer:

For face recognition, the preprocessing steps included:

Resizing:

 All images were resized to a **fixed dimension** (e.g., 96×96 or 160×160 pixels) for consistency across the dataset.

Normalization:

- Pixel values were scaled to a range between 0 and 1 (dividing by 255) to help faster model convergence.
- Face detection and alignment (optional):

 Detected faces from images and aligned them based on eye positions to handle pose variation.

• Data augmentation:

 Techniques like horizontal flipping, rotation, and slight zooming were applied to expand the dataset and make the model robust.

These preprocessing steps ensured **standardization**, **robust learning**, and **better generalization**.

Counter Question:

• Why normalize pixel values between 0 and 1?

Answer:

• To prevent large input values from **slowing down learning** and to **stabilize gradients** during backpropagation.

Q2. What CNN architecture or model did you use (custom or pre-trained)?

Answer:

- I used a pre-trained model called FaceNet (or a similar model like MobileFaceNet / VGG-Face).
- Alternatively, if I designed a custom CNN:
 - Multiple convolutional layers (Conv → ReLU → MaxPooling).
 - Followed by dense (fully connected) layers to map features into embedding space.
- Using pre-trained models offers:
 - Faster development.

 Better accuracy because they are trained on large face datasets.
Counter Question:
Why use a pre-trained model instead of training from scratch?
Answer:
 Training from scratch requires huge datasets and very high computational power; pre-trained models already capture general features.
Q3. What is face embedding in deep learning?
Answer:
• Face embedding is a vector representation of a face.
 It transforms each face image into a fixed-length feature vector (e.g., 128-dimensional that captures the essential characteristics of that face.
Properties:
$\circ \textbf{Similar faces} \rightarrow \textbf{Embeddings are } \textbf{closer} \text{ in vector space}.$
○ Different faces → Embeddings are far apart .
This enables face verification, clustering, and recognition tasks.
Counter Question:
How are embeddings different from raw pixel inputs?
Answer:

• Embeddings are **compact**, **learned feature representations** focusing only on **important patterns**, unlike raw pixels.

Q4. How does the Triplet Loss function work in face recognition?

Answer:

- Triplet Loss trains the model to bring similar faces closer and different faces farther apart.
- Each training example consists of:
 - o Anchor: A sample face.
 - **Positive**: A face of the **same person**.
 - Negative: A face of a different person.

Objective:

Distance(anchor, positive) + margin < Distance(anchor, negative)

• "Margin" is a safety gap to ensure separation.

Thus, Triplet Loss encourages better embedding spaces for distinguishing faces.

Counter Question:

• What could happen without the margin in triplet loss?

Answer:

• Embeddings may collapse (very close to each other), leading to **poor discrimination**.

Q5. How do you differentiate between classification and verification tasks in face recognition?

Answer:

- Face Classification:
 - o Identify which person the face belongs to among a fixed set of known people.
 - Requires softmax classification.
- Face Verification:
 - Decide whether **two faces belong to the same person** or not.
 - Requires comparing embeddings and measuring distance (e.g., cosine distance, Euclidean).

Thus, classification outputs a **class label**, verification outputs a **true/false match**.

Counter Question:

Which is harder to scale to thousands of people — classification or verification?

Answer:

• **Classification** — because the number of output neurons grows with the number of people.

Q6. How did you handle pose and lighting variations?

Answer:

Techniques used:

• **Data augmentation:** Introduced simulated pose changes (rotation, flips) and lighting variations.

- Face alignment: Aligned detected faces based on eyes or facial landmarks.
- Robust feature extraction: Used CNNs trained on large, varied datasets to ensure embeddings are invariant to lighting and pose.

These steps helped the model learn features robust to real-world variations.

Counter Question:

• How does augmentation help with pose variation?

Answer:

• It teaches the model to **recognize the same face** under **different orientations**.

Q7. What is one-shot learning and where is it useful in face recognition?

Answer:

- One-shot learning refers to:
 - Learning to recognize a new class (new person) with only one training example.
- Useful in face recognition:
 - You often need to recognize **new people** with very few examples available.
- Embedding-based models (like FaceNet) naturally support one-shot learning because they compare **features**, not require retraining.

Counter Question:

• Which loss function enables one-shot learning in face recognition?

Answer:
Triplet Loss or Contrastive Loss.
OO Mile at data assume antation to aleminuse help advesus project?
Q8. What data augmentation techniques helped your project?
Answer: I used:
Random horizontal flips (mirroring faces).
Random rotations (small angles like ±10 degrees).
Brightness and contrast adjustments.
Zoom-in and zoom-out slightly.
Random cropping.
These techniques made the model more robust to real-world variations and prevented overfitting .
Counter Question:
Is vertical flip a good augmentation for faces?
Answer:
No, because upside-down faces are unrealistic and could confuse the model.
Q9. How would you deploy the face recognition system into a mobile app?
Answer: Steps:

1. Model Optimization:

 Compress the trained model using techniques like quantization or pruning to reduce size.

2. Export Model:

Convert the model into mobile-friendly formats (e.g., TensorFlow Lite, ONNX).

3. Integrate into Mobile App:

• Load the model using mobile frameworks (TensorFlow Lite interpreter, Core ML).

4. Real-time Inference:

- o Capture images from camera.
- \circ Run face detection \rightarrow face embedding extraction \rightarrow compare with database.

Goal: Ensure **low-latency** and **high-accuracy** face verification on mobile devices.

Counter Question:

• Why is quantization important for mobile deployment?

Answer:

• It reduces **model size** and **inference time**, making it **faster** and **more efficient** on mobile hardware.

Q10. What improvements can be done using transfer learning?

Answer:

• Transfer learning allows:

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Example for BFS and DFS

Suppose we have the following graph:

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DEF

Where:

- Node A is connected to B and C.
- Node B is connected to D and E.
- Node C is connected to F.

BFS Traversal (Breadth-First Search)

(Level by Level)

Start at $A \rightarrow \text{explore neighbors}$ first, then their neighbors.

Steps:

- 1. Start at $A \rightarrow visit A$
- 2. Visit A's neighbors \rightarrow B and C
- 3. Visit B's neighbors \rightarrow D and E
- 4. Visit C's neighbor \rightarrow F

BFS Traversal Order:

DFS Traversal (Depth-First Search)

(Go Deep before Backtracking)

Start at $A \rightarrow \text{explore}$ as deep as possible along each branch before backtracking.

Steps (one possible DFS path):

- 1. Start at $A \rightarrow visit A$
- 2. Move to $B \rightarrow visit B$
- 3. Move to $D \rightarrow visit D$ (B's neighbor)
- 4. D has no children → backtrack to B
- 5. Visit E (B's other neighbor)
- 6. E has no children \rightarrow backtrack to B \rightarrow then to A
- 7. Move to $C \rightarrow visit C$
- 8. Move to $F \rightarrow visit F$ (C's neighbor)

DFS Traversal Order:

Quick Difference:

Feature	BFS	DFS
Style	Level by Level	Deep into branch
Data Structure	Queue	Stack (or recursion)

Of course! Here's a simple step-by-step example of Merge Sort and Bubble Sort — without any code — just how you need for viva:

Example of Merge Sort

Suppose we have this array:

Steps:

- 1. Divide the array recursively into halves:
 - [8, 4, 5] and [2, 9, 1]
 - Further divide:
 - 0 [8] [4,5]
 - o [2] [9,1]
 - Further divide:
 - o [4] [5]
 - 0 [9] [1]
- 2. Now start merging while sorting:

- Merge [4] and [5] → [4,5]
- Merge [9] and [1] → [1,9]
- Merge [8] and [4,5] → [4,5,8] (sort while merging)
- Merge [2] and [1,9] → [1,2,9]
- 3. Finally merge the two sorted halves:
 - Merge [4,5,8] and [1,2,9]
 - Compare and merge:

$$\circ \quad 1 \rightarrow 2 \rightarrow 4 \rightarrow 5 \rightarrow 8 \rightarrow 9$$

V Final sorted array:

[1, 2, 4, 5, 8, 9]

Example of Bubble Sort

Suppose we have the same array:

Steps:

- 1st Pass:
 - Compare 8 and $4 \rightarrow \text{Swap} \rightarrow [4, 8, 5, 2, 9, 1]$
 - Compare 8 and $5 \rightarrow \text{Swap} \rightarrow [4, 5, 8, 2, 9, 1]$

- Compare 8 and $2 \rightarrow \text{Swap} \rightarrow [4, 5, 2, 8, 9, 1]$
- Compare 8 and 9 → No swap
- Compare 9 and $1 \rightarrow \text{Swap} \rightarrow [4, 5, 2, 8, 1, 9]$

• 2nd Pass:

- Compare 4 and 5 → No swap
- Compare 5 and $2 \rightarrow \text{Swap} \rightarrow [4, 2, 5, 8, 1, 9]$
- Compare 5 and 8 → No swap
- Compare 8 and $1 \rightarrow \text{Swap} \rightarrow [4, 2, 5, 1, 8, 9]$

• 3rd Pass:

- Compare 4 and $2 \rightarrow \text{Swap} \rightarrow [2, 4, 5, 1, 8, 9]$
- Compare 4 and 5 → No swap
- Compare 5 and $1 \rightarrow \text{Swap} \rightarrow [2, 4, 1, 5, 8, 9]$

• 4th Pass:

- Compare 2 and 4 → No swap
- Compare 4 and 1 → Swap → [2, 1, 4, 5, 8, 9]

• 5th Pass:

Compare 2 and 1 → Swap → [1, 2, 4, 5, 8, 9]

Final sorted array:

[1, 2, 4, 5, 8, 9]

Quick Difference:

Feature	Merge Sort	Bubble Sort	
Strategy	Divide and Conquer (Divide → Sort → Merge)	Repeatedly swap adjacent elements	
Efficiency	Very efficient for large datasets	Very slow for large datasets	
Time Complexity	O(n log n)	O(n²)	
Space Requirement	Extra space needed (for merging)	In-place (no extra space)	