

# Characterization of Large Language Model Development in the Datacenter

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## Abstract

Large Language Models (LLMs) have presented impressive performance across several transformative tasks. However, it is non-trivial to efficiently utilize large-scale cluster resources to develop LLMs, often riddled with numerous challenges such as frequent hardware failures, intricate parallelization strategies, and imbalanced resource utilization. In this paper, we present an in-depth characterization study of a six-month LLM development workload trace collected from our GPU datacenter Acme. Specifically, we investigate discrepancies between LLMs and prior task-specific Deep Learning (DL) workloads, explore resource utilization patterns, and identify the impact of various job failures. Our analysis summarizes hurdles we encountered and uncovers potential opportunities to optimize systems tailored for LLMs. Furthermore, we introduce our system efforts: (1) *fault-tolerant pretraining*, which enhances fault tolerance through LLM-involved failure diagnosis and automatic recovery. (2) *decoupled scheduling for evaluation*, which achieves timely performance feedback via trial decomposition and scheduling optimization.

## 1 Introduction

Over the years, advances in LLMs have attracted significant attention from both academia and industry owing to their impressive performance and capabilities, such as ChatGPT [2] and GitHub Copilot [3]. However, due to their immense model sizes and extensive data demands, training such models necessitates a substantial computational infrastructure with thousands of accelerators [27, 69]. Hence, it is a common practice for tech companies and cloud providers to build large-scale GPU clusters to facilitate LLM development, especially after the popularity of ChatGPT. Nevertheless, it is non-trivial to perform efficient LLM development on such high-cost infrastructure. Developers often confront numerous issues and challenges, including frequent hardware failures [65, 97], intricate parallelization strategies [69, 114], unstable training progress [1, 111], long queuing delay [105], etc.

Developing LLMs is closely intertwined with the support of GPU clusters in various aspects. A thorough analysis of cluster workloads is essential for comprehending challenges and uncovering opportunities in designing systems tailored

for LLMs. However, many conclusions and implications from existing DL workloads analysis works [41, 47, 98], conducted before the rise of LLMs, are not applicable to LLM development. This is primarily due to the divergent characteristics and requirements of LLMs:

(1) **Paradigm Transition.** DL workloads generally follow a *task-specific* paradigm that trains the model on domain-specific data to tackle a particular task (e.g., translation [18]). In contrast, LLMs follow an emerging paradigm that performs self-supervised training on broad data to generate a *foundation model* [19] and further adapts to a wide range of downstream tasks. This shift signifies a substantial divergence in the model development pipeline (e.g., pretraining [86], alignment [40]) and workload characteristics from prior DL workloads (§2.1). (2) **Tailored Software Stack.** To accommodate the enormous model size of LLMs, a series of systems implement advanced techniques to optimize the execution of LLMs. For instance, Deepspeed [81], Megatron [69] and Alpa [114] accelerate the training via hybrid parallelism or state-sharding optimizer. As for model serving, Orca [105] and vLLM [52] improve throughput via iteration scheduling or memory management. (3) **Unified Architecture.** Prior DL workloads usually employ various model architectures (e.g., CNN [55], RNN [18]) to address diverse tasks. In contrast, LLMs commonly embrace the Transformer [94] architecture, like BERT [31], GPT-3 [20], LLaMA [92] and PaLM [27]. The architectural homogeneity suggests a high level of uniformity in the LLM development pipeline and similarity across different datacenters.

To bridge this gap, we present an in-depth study of our operational experiences in the datacenter Acme of Shanghai AI Laboratory. It houses two distinct clusters, Seren and Kalos, dedicated to LLM development and equipped with 4,704 A100 GPUs in total. Our analysis draws upon traces collected over a six-month period from March to August 2023, encompassing scheduler logs, infrastructure monitoring data, failure logs, and fine-grained profiling data. Our key findings and identified challenges can be summarized as follows:

- **Shorter Job Duration and Unfair Queuing Delay.** In contrast to the common stereotype that LLM workloads are usually long-term, the workloads in our datacenter exhibit 2.7~12.8× shorter average job duration compared to the DL workloads in previous traces [41, 47, 98]. This can be

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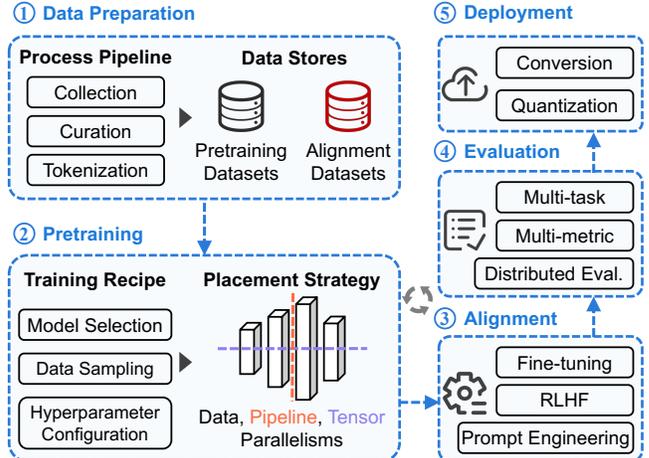
attributed to the presence of numerous short-term tasks such as evaluation. In terms of job queuing delay, our findings also diverge from previous DL traces that larger-scale jobs experience longer wait times. We observe that evaluation jobs, despite being short-term and small-scale, have the longest queuing delay. This discrepancy stems from reserving the majority of resources for pretraining jobs to minimize their queuing delays. Evaluation jobs are scheduled with a lower priority, utilizing the limited spare resources.

- **Imbalanced Resource Usage.** The imbalance is manifested in two aspects. Firstly, in terms of workload distribution, pretraining jobs only account for 3.2% of the total job count but consume 94.0% of the whole compute resource (i.e., GPU time) in Kalos. Conversely, evaluation jobs, despite constituting 92.9% of all jobs, only utilize a meager 0.8% of resources. Secondly, when looking at infrastructure utilization, we find that associated resources including CPU, host memory, and network, are frequently underutilized. In contrast, the GPU, as the primary resource, shows high utilization rates. Both GPU memory and GPU utilization exhibit substantially higher median values at 75% (60GB) and 99% respectively in Kalos, as opposed to the 10% and 4% observed in PAI [98]. These observations corroborate that LLMs are computationally and memory intensive. It also implies that GPU-sharing-based techniques [42, 99, 100, 107] may not be suitable for LLMs.

- **Long GPU Idle Time in Evaluation Workload.** Our profiling of evaluation workloads reveals substantial underutilization of GPU resources at various stages. For example, the evaluation job on HumanEval consumes 29.5% of its time for model loading and data preprocessing, and an additional 19.0% is spent conducting synthesized program correctness tests. As a result, only half of the time is dedicated to GPU inference, leading to long queuing delays in evaluation trials.

- **Frequent Job Failures.** We find various errors primarily occur at the beginning of LLM workloads, leading to fast job termination. However, infrastructure failures, which are common in long-term pretraining jobs, significantly impede training efficiency. Therefore, prompt diagnosis and recovery from these failures are crucial to enhance training efficiency.

Based on our characterization study, we identify several challenges encountered during the LLM development, such as unstable training progress, remote storage bottleneck and delayed feedback on model performance. To tackle these issues, we consolidate the insights gained from our operational experience and build two systems that are integrated into our LLM framework to improve development robustness and efficiency. *Firstly*, to mitigate the frequent failure problem, we establish a system to achieve *fault-tolerant pretraining*. It incorporates three key designs: (1) achieving frequent model saving through asynchronous checkpointing, (2) identifying the root causes of various failures through a combination of heuristic rules and LLM, (3) employing a holistic detection toolkit to pinpoint fault nodes and automatically restart training from properly saved checkpoint. It accelerates checkpoint-



**Figure 1:** Overview of the LLM development pipeline.

ing by  $3.6\sim 58.7\times$  and significantly reduces manual intervention. *Secondly*, we develop a system that performs *decoupled scheduling for evaluation* to provide developers with timely feedback on model quality. It not only resolves the remote model loading contention issue via decoupled model retrieval but also minimizes GPU idle time via decoupling the metric computation process. It further leverages the prior knowledge and flexibility of datasets to balance workload across all GPUs. Our experiment shows that it can reduce the evaluation makespan by up to  $1.8\times$ .

We believe the observations and insights derived from our datacenter do not stand in isolation. Our traces are publicly available at <https://github.com/InternLM/AcmeTrace>. We also release our system code and models (Appendix C). We hope these resources and lessons can benefit researchers in optimizing LLM systems as well as GPU cluster management.

## 2 Background

### 2.1 LLM Development Pipeline

Distinguished from task-specific DL models, LLMs follow an emerging paradigm that performs self-supervised training on broad data and further adapts to a wide range of downstream tasks [19]. The development of LLMs necessitates the use of extensive computational infrastructure due to their substantial model size (comprising billions of parameters) and the vast amount of training data involved. Figure 1 depicts the comprehensive LLM development pipeline, encompassing five distinct stages (blue blocks) that span from scratch to service (follow blue arrows). The grey circular arrow indicates that the pretraining stage enables periodical alignment and evaluation to assess intermediate models and adjust configuration on the fly. We explain each stage as follows:

**Data Preparation.** The initial stage involves gathering and preprocessing the training data, which can be categorized into two parts: (1) *pretraining data*, consisting of extensive unlabeled corpora obtained from public or private sources and curated through processes like detoxification and dedu-

Cluster	#CPUs	#GPUs	Mem(GB)	Network	#Nodes
<b>Seren</b>	128	8	1,024	1×200Gb/s	286
<b>Kalos</b>			2,048	5×200Gb/s	302

**Table 1:** Summary of per-node specification and cluster scale for two independent LLM clusters in Acme.

plication; (2) *alignment data*, comprising a smaller set of high-quality labeled corpora used to align the model with specific tasks. This data is typically acquired through expensive human annotation or labeling [73]. Besides, all the data must be tokenized to ensure compatibility with the model’s input.

**Pretraining.** It involves self-supervised training on large-scale curated data, demanding a majority of resources within the overall development workflow. Training LLMs efficiently at scale necessitates various system innovations, such as state-sharding optimizers [81], meticulous model placement using data, pipeline, and tensor parallelisms [68, 69, 114].

**Alignment.** This stage aims to adapt LLMs with user intent on a wide range of downstream tasks. Two primary aligning paradigms are commonly used: (1) *prompt engineering*, specifying prompts (i.e., inputs) without modifying model parameters. For example, in text summarization, appending a prompt *TL; DR* to the input article can improve model performance [80]; (2) *fine-tuning*, updating model parameters on a task-specific dataset to improve performance in a particular domain. Additionally, reinforcement learning from human feedback (RLHF) [73] further enhances the alignment effect, and parameter-efficient techniques like LoRA [40] have been proposed to reduce the cost of fine-tuning.

**Evaluation.** Given the vast application scenarios of LLM, it may be inaccurate to assess model quality solely based on a single metric like training loss. There are numerous factors to consider, such as accuracy, fairness, and toxicity [59]. Consequently, it is crucial to account for a diverse set of criteria and measure performance across multiple tasks [22]. Furthermore, regular evaluation is essential during the pretraining stage to provide timely feedback on model quality.

**Deployment.** To meet the strict cost and latency constraints of LLM applications, several advanced techniques have been developed to achieve efficient model serving, including quantization [30], distillation [84], CUDA kernel optimization [29, 45], model parallelism [58, 105] and memory management [52].

## 2.2 Acme Overview

Acme is our private GPU datacenter that empowers researchers and engineers to develop DL models across diverse domains. In this work, we focus on analyzing workloads within two clusters dedicated to developing LLMs: Seren and Kalos. We collect and analyze all jobs in these two clusters. Note that there are additional clusters within Acme that are designated for different fields, such as autonomous driving, and AI for scientific research. However, these clusters are excluded in this work as they are unrelated.

**Cluster Architecture.** Table 1 summarizes configurations of these two homogeneous LLM clusters. Seren and Kalos have 2,288 and 2,416 GPUs respectively. Each node is equipped with 8× NVIDIA A100-SXM 80GB GPUs [6] and 2× Intel Xeon Platinum 8358P CPUs (128 threads in total). GPUs are interconnected to each other by NVLink and NVSwitch [9], and inter-node communication is achieved via NVIDIA Mellanox 200Gbps HDR InfiniBand [4]. Compared to Seren, Kalos is a relatively newer cluster with an improved network configuration. Each node in the Kalos has a larger host memory (2TB) and is equipped with four InfiniBand HCAs specifically for application communication, along with an extra HCA dedicated to storage.

Besides, the distributed storage system is also critical for workload performance. Acme adopts an all-NVMe shared parallel file system for fast data access and storage. Moreover, as time has advanced, our resource scheduling system has evolved to support diverse cluster environments. Specifically, our scheduler on Seren and Kalos is built atop Slurm [103] and Kubernetes [21] respectively. In order to provide resource guarantees for large-scale pretraining jobs, our scheduler enables resource isolation and quota reservation. It further incorporates a best-effort job mechanism for higher utilization.

**LLM Workloads.** We develop a collection of LLMs<sup>1</sup> ranging from 7B to over 123B parameters. All of these models follow the transformer-based decoder-only architecture, similar to the GPT [20, 79, 80] and LLaMA [92, 93] series. Acme encompasses tasks in the aforementioned general LLM development pipeline (§2.1). Note that Acme does not involve any serving jobs, as our LLMs are deployed on a separate cluster specifically for serving purposes.

**Software Stack.** To support the training of billion-scale models across thousands of GPUs, we built a system InternEvo<sup>2</sup>, which integrates various system optimization techniques, such as FlashAttention [28, 29], 3D parallelism [69], zero redundancy optimization [81], mixed precision training [10], selective activation recomputation [51] and fine-grained communication overlap. Moreover, it accommodates additional tasks such as model fine-tuning and evaluation.

## 2.3 Traces from Acme

The optimization of LLM-tailored systems and datacenter management can significantly boost development efficiency and yield substantial financial benefits. Achieving this goal requires a profound understanding of the intrinsic characteristics of LLM workloads. Many insights in existing DL workloads analysis works [41, 47, 98] are not applicable to LLM workloads due to the unique attributes of LLMs. To fill this gap, we collected and analyzed workload traces from our datacenter Acme. Table 2 compares the specifications and trace information of Acme with prior trace analysis works conducted by Microsoft, SenseTime, and Alibaba. Unlike

<sup>1</sup>Model: <https://huggingface.co/internlm>

<sup>2</sup>System: <https://github.com/InternLM/InternEvo>

Datacenter	For Task-Specific DL Models			For LLMs
	Philly [47]	Helios [41]	PAI [98]	Acme
Year	2017	2020	2020	2023
Duration	3 months	6 months	2 months	6 months
#Jobs	113K	3.36M	1.26M	1.09M
Avg. #GPUs	1.9	3.7	0.7	6.3
GPU Model	12GB/24GB	1080Ti/V100	T4/P100/V100	A100
Total #GPUs	2,490	6,416	6,742	4,704

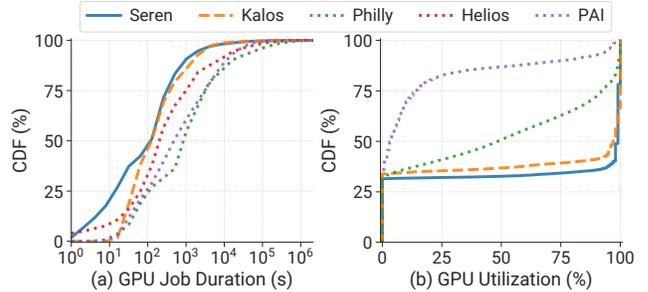
**Table 2:** Comparison between our datacenter Acme and GPU datacenters in prior trace analysis works: Microsoft Philly [47], SenseTime Helios [41], Alibaba PAI [98]. Philly only provides GPU memory sizes without clarifying GPU models. The average number of requested GPUs in PAI can be less than 1 (0.7), as it supports fractional ( $<1$ ) GPU requests.

Acme, which is solely dedicated to LLM development, these datacenters encompass a mixture of general DL workloads from various domains. For instance, Helios [41] consists of 4 clusters dedicated to training models in computer vision and reinforcement learning, while PAI [98] includes a diverse range of servers for training and serving jobs.

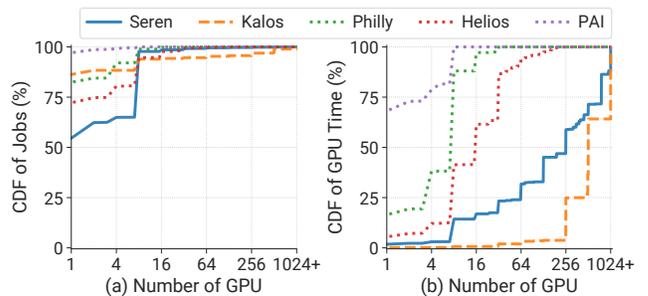
**Trace Source.** Our characterization study is based on traces collected from two LLM clusters in Acme. The traces span 6 months from March to August 2023. Seren contains 368K CPU jobs and 664K GPU jobs, while Kalos job trace consists of 42K CPU jobs and 20K GPU jobs. Additionally, we provide a summary of the data sources for the traces used in our study: (1) *Job Log*. We collect the job logs from our scheduler database, which contains detailed information for each job. This includes the job’s execution time (submission, start, and end), final status (completed, canceled, failed), requested resources (CPU, GPU, memory), work directory, and other relevant data. (2) *Hardware Monitor Data*. This encompasses long-term, multi-dimensional data obtained from various sources. We collect CPU, memory, and network usage data from Prometheus [77] database, GPU-related metrics from NVIDIA DCGM [7], and power-related data from IPMI [12]. The sampling interval for this data is set at 15 seconds. (3) *Runtime Log*. To conduct a precise job failure analysis, we capture stdout and stderr logs from LLM frameworks during job execution. (4) *Profiling Data*. For a subset of representative jobs, we delve deeper by performing fine-grained profiling using tools like DCGM. The synergy of these trace dimensions allows us to gain a holistic understanding of LLM job characteristics in datacenters.

### 3 Datacenter Characterization

In this section, we perform a thorough analysis of Acme, including comparing workload distribution between LLMs and previous DL workloads (§3.1), investigating different LLM workload types (§3.2), exploring resource utilization patterns (§3.3) and assessing environmental impacts (§3.4).



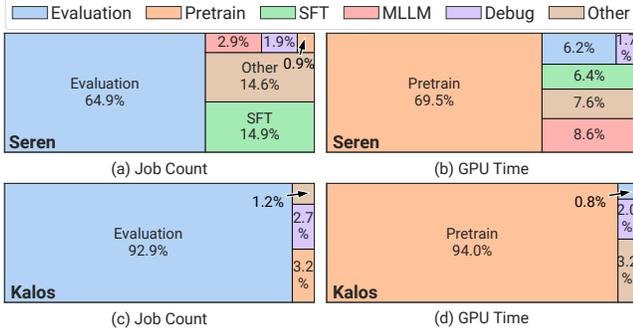
**Figure 2:** Overview of different datacenter characteristics. (a) *Workload*: CDF of the GPU job duration. (b) *Infrastructure*: CDF of GPU utilization, where Helios’ data is not available.



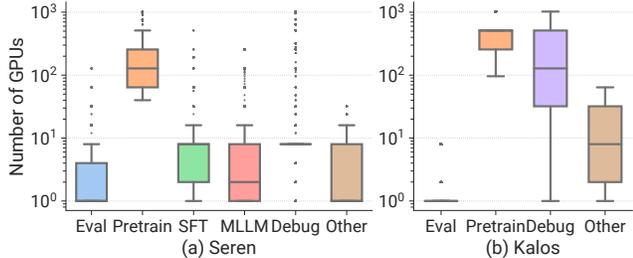
**Figure 3:** Comparison of workload distribution based on the number of requested GPUs. (a) CDF of job count. (b) CDF of GPU time (i.e., requested GPU number  $\times$  duration).

#### 3.1 LLMs versus Prior DL Workloads

**Shorter Job Duration.** As shown in Figure 2 (a), contrary to the prevailing stereotype that LLM-related jobs are typically long-running, we find the workloads in our clusters (blue and orange lines) exhibit shorter GPU job durations (i.e., job runtime, excluding queuing delay) compared to the DL workloads observed in previous job traces (dotted lines). Specifically, both the Seren and Kalos have a median job duration of 2 minutes, which is  $1.7\sim 7.2\times$  shorter than the median job durations of other clusters. Furthermore, it is evident that the more recent trace demonstrates a shorter job duration distribution. In particular, when considering the average job duration in the Philly cluster (collected in 2017), it is  $2.7\sim 3.8\times$  longer than Helios (2020) and PAI (2020), and  $12.8\times$  longer than Acme (2023). To provide an explanation for this observation, we outline three potential factors: (1) *Hardware upgrade*. The iteration of GPU and networking delivers substantial efficiency improvement. (2) *Abundant resources*. Users usually request more resources (as shown in Table 2), averaging 5.7 GPUs in the Seren and 26.8 GPUs in the Kalos. This can significantly accelerate the training process. (3) *Extensive associated workloads*: LLM development pipeline involves numerous small-scale associated jobs, such as evaluation. We will delve into this in §3.2. (4) *High incompleteness rate*: Approximately 40% of jobs fail, with completed jobs consuming only 20~30% of GPU resources. This highlights the urgent need for a fault-tolerant system. Further details can be found in Figure 17 and Appendix A.1.



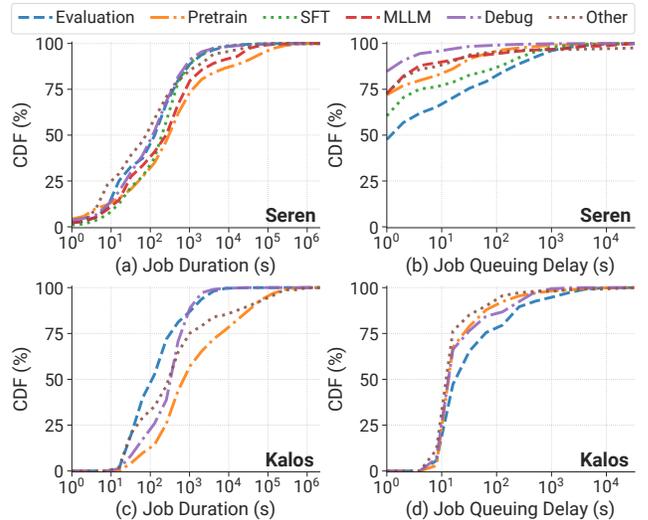
**Figure 4:** Distribution of different workload types in Seren (a, b) and Kalos (c, d). Note that CPU jobs are excluded. *SFT*: Supervised Fine-Tuning for model alignment. *MLLM*: Multimodal Large Language Model. *Other*: Unclassified jobs.



**Figure 5:** The boxplot of the distribution of GPU demand across different workload types in Seren (a) and Kalos (b).

**Polarized GPU Utilization** Figure 2 (b) shows cluster-wide GPU utilization distributions across various clusters. It is evident that the GPU utilization in our two clusters exhibits a polarized pattern, primarily concentrated in two distinct states: 0% and 100%. This polarization mainly stems from the fact that the workloads in our clusters share similar model architectures, i.e., transformer-based LLMs. In contrast, Philly and PAI encompass a broader range of utilization. Besides, when comparing the median GPU utilization, Seren and Kalos exhibit significantly higher values at 97% and 99%, respectively, in contrast to 48% and 4% observed in Philly and PAI. This observation aligns with the common understanding that LLMs are computationally intensive. It also implies that GPU-sharing-based scheduling techniques [42, 99, 100, 107] may not be suitable for LLM development. Note that ‘GPU utilization’ may sometimes be a weak utilization indicator [8, 95]. We provide a more precise utilization analysis in §3.3.

**High-skewed Workload Distribution.** We further investigate the CDF of GPU demands in relation to the number of jobs (Figure 3 (a)) and GPU time (Figure 3 (b)). For the number of jobs, all the clusters share a similar pattern in that the majority of jobs are single-GPU jobs and less than 7% of jobs request over 8 GPUs. However, when examining GPU time, single-GPU jobs only account for less than 2% resources in our two clusters, while taking over 68% GPU time in PAI. In stark contrast, large-scale jobs ( $\geq 256$  GPUs) dominated the GPU time in Kalos, occupying more than 96% of resources. The much steeper distribution poses substantial challenges for the design of cluster schedulers. A majority of resources



**Figure 6:** CDF of GPU job duration and queuing delay for different workload types in Seren (a, b) and Kalos (c, d).

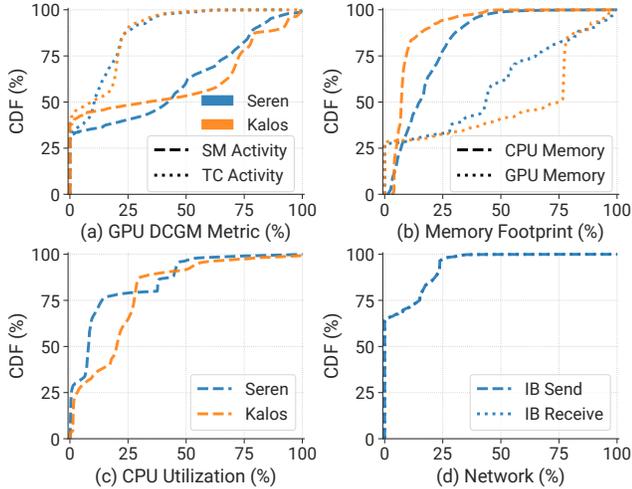
are allocated to a few pretraining jobs, potentially causing head-of-line blocking issue and resulting in severe queuing delay. Existing DL cluster schedulers [37, 62, 76, 99, 102] typically depend on preemption mechanism, however, the considerable recovery overhead makes them not applicable to LLM workloads. This highlights the critical need for a scheduling system tailored for LLM clusters, considering the workload features of the entire pipeline.

### 3.2 Workload Categories

To strive for a deeper understanding of the characteristics of different workloads in the LLM development pipeline (§2.1), we further categorize jobs into specific types according to their production division and metadata (e.g., job names).

**Irrelevance of Job Count and Resource Usage.** Figure 4 presents the distribution of job counts and GPU time across various workload types, where only Seren contains SFT and MLLM workloads. Besides, MLLM jobs incorporate their own development pipeline (e.g., pretraining) and adopt smaller model scales for exploration purposes. Our analysis primarily focuses on LLM jobs. It is obvious that evaluation jobs constitute the majority of the total job count in both clusters, yet they consume a relatively small portion of resources (0.8% in Kalos). In contrast, pretraining jobs only account for 0.9% and 3.2% of the total job count but consume 69.5% and 94.0% of the total GPU time in Seren and Kalos respectively.

**Job Type Correlates with GPU Demand.** We further depict GPU demand distribution across various workload types in Figure 5. Each box is framed by the first and third quartiles, while the median value is indicated by the black line within the box. Both whiskers are defined at  $1.5 \times$  the InterQuartile Range (IQR). Compared to evaluation jobs, which typically require less than 4 GPUs, pretraining jobs often require over 100 GPUs. This observation partially explains why evaluation jobs in Kalos consume only minimal resources in Figure 4(d). Additionally, we notice that debugging jobs have a wide range



**Figure 7: Infrastructure utilization.** CDF of various metrics: (a) active fractions of Streaming Multiprocessor (SM) and Tensor Core (TC), (b) memory footprints of host and GPU, (c) CPU utilization, (d) normalized InfiniBand (IB) HCA send and receive bandwidths (Seren only). Seren and Kalos are represented by blue and orange lines respectively.

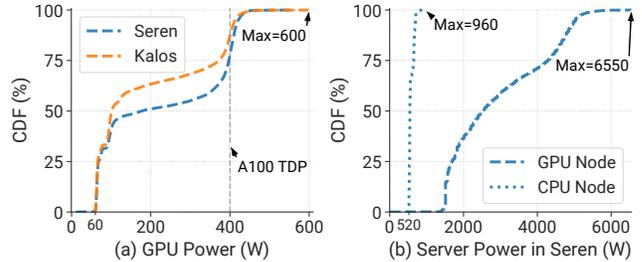
of GPU requests, which aligns with the fact that testing jobs are typically needed for various types of tasks.

**Similar Temporal Distribution.** Figure 6 shows the distribution of job duration and queuing delay across different workloads. In terms of job duration, although pretraining jobs have the longest duration, they surpass other workloads within an order of magnitude in the median, and less than 5% jobs last for over 1 day in both clusters. This can be attributed to frequent failures during pretraining, which will be further explored in §5. Regarding job queuing delay, contrary to previous reports [41, 47, 98] suggesting that larger-scale jobs experience longer wait times, we observe that evaluation jobs have the longest queuing delay despite having the lowest GPU demands and shortest job duration. This discrepancy is due to the majority of resources being reserved for pretraining jobs to minimize their queuing delays. Evaluation jobs are typically submitted as a batch simultaneously with lower priority, utilizing the limited spare resources.

### 3.3 Infrastructure

Beyond the workload characterization, we further conduct a comprehensive analysis of our infrastructure utilization.

**Higher GPU Utilization.** Given the critical role of GPUs in LLM development, as shown in Figure 7 (a, b), we collect fine-grained performance counter metrics from DCGM [7], including SM Activity (PROF\_SM\_ACTIVE), TC Activity (PROF\_PIPE\_TENSOR\_ACTIVE), and GPU memory footprint (DEV\_FB\_USED). In contrast to PAI [98], where a significant portion of GPU memory is underutilized (less than 25% memory), our observations in Kalos indicate that 50% of GPUs consume over 75% of GPU memory (60 GB). Furthermore, we observe that the median SM activity in both clusters is approximately 40%, which is twice the reported 20% in PAI.



**Figure 8: Power consumption.** CDF of (a) A100 GPU power, (b) server power in Seren. *TDP*: Thermal Design Power.



**Figure 9: Average power distribution of hardware modules in Seren GPU servers,** gathered from IPMI and DCGM.

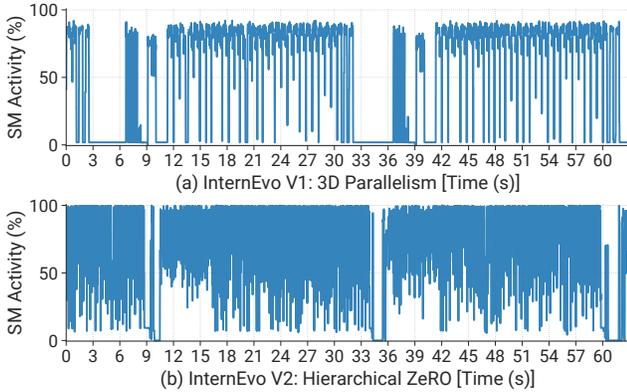
These findings align with the memory-intensive and compute-intensive natures of LLMs.

**Underutilized Associated Resources.** We also delve into the aspects of CPU, host memory, and network that are closely associated with LLM development. In Figure 7 (b), we compare the memory footprint on the host side and GPU side. It is evident that CPU memory utilization remains below 50%. Note that Kalos boasts twice the memory capacity (2TB) compared to Seren (Table 1). This demonstrates the significant underutilization of CPU memory. More detailed analysis is provided in Appendix A.2. Although the GPU memory offloading technique [82, 83] improves CPU memory utilization and alleviates GPU memory limitations, it also impedes training throughput due to limited PCIe bandwidth. Therefore, we do not employ the offloading mechanism. Additionally, due to a high CPU-to-GPU ratio (16 CPUs per GPU), CPUs are typically underutilized, as depicted in Figure 7 (c). Moreover, in Figure 7 (d), we measure the network send and receive bandwidths of IB in Seren. Two lines are well overlapped, as IB serves for symmetrical communication during LLM execution. We observe that NICs remain idle for over 60% of the time, and the active bandwidth rarely surpasses 25% of the maximum bandwidth provided by IB.

### 3.4 Environmental Impact

LLM development leads to substantial energy consumption and carbon emissions [74, 104]. We report our analysis of infrastructure power consumption patterns to inspire future datacenter designs that minimize environmental impact.

**GPUs Dominate Power Consumption.** Figure 8 (a) depicts the distribution of GPU power consumption. We observe that around 30% of GPUs are in an idle state and still need to consume 60W. Besides, due to intensive computation demand, we find that 22.1% and 12.5% of GPUs consume over 400W (TDP) in Seren and Kalos respectively, with some even reaching 600W. This may cause the risk of some metastable issues [43]. Figure 8 (b) presents the power consumption distribution of all GPU servers, along with an additional 6 CPU-



**Figure 10:** GPU SM utilization of pretraining a 123B LLM using different strategies of InternEvo [25] over 2048 GPUs.

only servers, in Seren. We find GPU servers consume  $5\times$  power than CPU servers on average. Additionally, Figure 9 demonstrates that GPUs account for approximately  $2/3$  of the total power consumption in GPU servers, while CPUs only contribute 11.2% and power supply units (PSUs) consume 9.6% of the energy during voltage conversion. These observations align with the understanding that GPUs are the primary power consumers in LLM development.

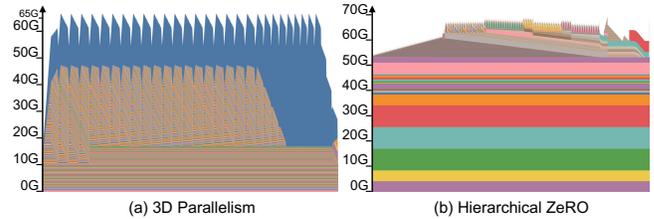
**Carbon Emission.** Our datacenter Acme has a PUE (Power Usage Effectiveness) of 1.25. Moreover, it operates on approximately 30.61% of carbon-free energy (statistics for the year 2022), which includes renewable sources like solar and wind power and achieves a carbon emissions footprint ( $\text{CO}_2\text{e}$ ) rate of 0.478  $\text{tCO}_2\text{e}/\text{MWh}$ . Based on our node-level power consumption data, we calculate that Seren consumed approximately 673 MWh electricity in May 2023, which leads to total effective emissions of 321.7  $\text{tCO}_2\text{e}$ . We believe that implementing advanced approaches like [14, 26, 104] can effectively reduce carbon emissions.

## 4 Workload Profiling

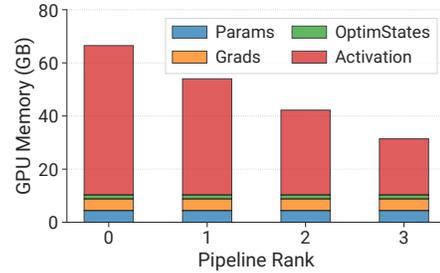
In this section, we conduct fine-grained analyses of resource utilization for representative tasks. Specifically, we focus on pretraining and evaluation jobs, as they are the most resource-intensive or quantity-intensive workloads.

### 4.1 Pretraining Workload

As aforementioned, pretraining LLMs requires substantial computational resources. To enhance training efficiency, our pretraining framework, InternEvo [25], undergoes continuous refinement and iteration in its system design. As presented in Figure 10, the initial version of InternEvo (adopted by our early jobs) is denoted as (a) primarily utilizes 3D parallelism akin to that of MegatronLM [69], and (b) employs a hierarchical ZeRO mechanism [25] that implements selective redundant sharding of model states. To provide a detailed example, we profile an LLM with 123 billion parameters across 2048 GPUs. We also provide the profiling results of 1024 GPUs in Appendix A.3. For (a) 3D parallelism approach, we adopt a configuration with pipeline parallelism= 4,



**Figure 11:** Memory snapshot under different pretraining strategies. Note that the extensive blue segment at the top of (a) is simplified and can be further broken down into massive fragments (memory allocations), similar to the lower part.

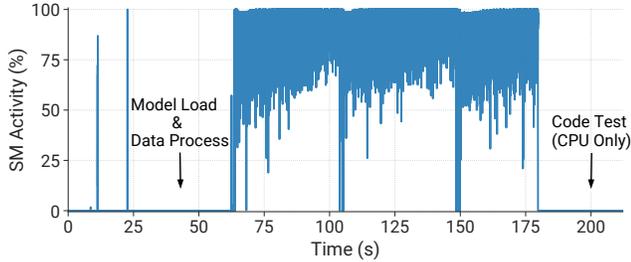


**Figure 12:** GPU memory consumption of different pipeline ranks employing the 1F1B [68] strategy in InternEvo V1.

tensor parallelism= 8. We sample the first GPU of the first pipeline rank for profiling. For (b) hierarchical ZeRO approach, we limit parameter sharding to subgroups of 64 GPUs each. We collect GPU performance counters like DCGM metrics at 1 ms intervals.

**GPU SM Utilization.** Figure 10 illustrates the GPU SM utilization for the same LLM under various training strategies. Both versions maintain the same global batch size and are optimized according to their respective configurations. It is evident that InternEvo V2 presents superior peak SM utilization and exhibits reduced idle periods compared to InternEvo V1, achieving around 16% acceleration. The relatively low utilization of 3D parallelism is mainly due to the impact of communication introduced by hybrid parallelism on the critical path, such as bubbles in pipeline parallelism. Note that the different inter- and intra-node communication hardware settings may lead to different optimal configurations.

**GPU Memory Footprint.** For a model comprising  $\Psi$  parameters, in the mainstream mixed precision training using Adam [49] optimizer, the memory footprint of the parameters, gradients, and optimizer states are  $2\Psi$ ,  $2\Psi$ , and  $12\Psi$ , respectively. To reduce memory cost, ZeRO [81] effectively shards redundant memory of these elements across global GPU workers. Figure 11 illustrates the actual GPU memory usage over time captured by the Pytorch memory snapshot tool [11]. The upper dynamic part represents activations and the lower static part represents the memory occupied by parameters, gradients and optimizer states. Note that only allocated memory is depicted, while reserved memory is not presented. Our analysis reveals that, in comparison to hierarchical ZeRO, the memory requirement for activations in 3D parallelism is substantially higher, even surpassing the sum of other components. This



**Figure 13:** GPU SM utilization for the entire evaluation workload on HumanEval [24] dataset using a 7B LLM.

observation underscores the importance of efficient activation memory management as a key factor for enhancing batch size and throughput in 3D parallelism.

**Imbalance in Activation Sizes.** When employing pipeline parallelism, each rank needs to hold a different quantity of activations since the diverse number of micro-batches pending backward computation across various pipeline ranks. Figure 12 illustrates this imbalance issue on different pipeline ranks. It suggests that we should employ a specialized partitioning mechanism to address the unbalanced memory usage among different ranks in pipeline parallelism, in order to achieve higher efficiency, such as recomputing activations.

## 4.2 Evaluation Workload

It is necessary to regularly evaluate the checkpoints produced during pretraining to guide the evolution of LLM pretraining. Therefore, the LLM evaluation jobs take the majority of jobs, each performing metric computation on different LLM benchmark datasets. We analyze the workflow of the entire evaluation and combine it with fine-grained resource usage information collection quantitatively, demonstrating two upcoming resource utilization issues. We will also discuss the corresponding solutions in §6.2.

**High Model Loading and Data Preprocessing Overhead.** During the initiation phase of evaluation jobs, it is imperative to load model checkpoints for each task. Additionally, the data preprocessing stage, particularly for tokenization, constitutes a significant time expenditure. These factors contribute to the underutilization of allocated GPU resources for a relatively long period. As illustrated in Figure 13, the evaluation task consumes over 1 minute prior to the actual GPU inference, accounting for 29.5% of the evaluation duration. This overhead is likely to increase with larger models or datasets. To address the preprocessing overhead, one effective strategy is to cache the tokenized data. Moreover, evaluation jobs are flexible, allowing for the consolidation of multiple evaluation tasks (datasets) into a single job. This consolidation can effectively reduce the relative time overhead of the model loading phase within the evaluation process.

**High Metric Computation Overhead.** The evaluation process can often involve complex and time-consuming metric computation. For example, synthesized program correctness tests need to be performed on coding datasets like Hu-

manEval [24] and MBPP [17]. Moreover, the OpenAI GPT-4 API is invoked to assess the performance of model conversations (e.g., Chatbot Arena [113]). These procedures can take up to 30 minutes, during which the GPU remains idle. Therefore, we can observe distinct stages of GPU usage, including stages that require GPU for inference and generation, and stages that do not require GPU for metric computation and verification. Taking the HumanEval benchmark as an example, as shown in Figure 13, the GPU is idle for the last 42 seconds, wasting about 19.0% of the total GPU time.

## 5 Failure Analysis

In this section, we conduct a comprehensive analysis of job failures, primarily relying on *runtime logs* and *hardware monitor data* from our two clusters. In KaLos, we gather logs from 32,500 tasks, which include 31,293 (96.3%) inference tasks, 647 (2.0%) pretraining tasks, and debugging tasks (1.7%). In Seren, we only collected logs from 675 pretraining tasks. Additionally, for pretraining tasks, we extract all pertinent information and metadata recorded in the logs, including actual training steps, cold-start overhead, recovery timestamp, etc. We hope our analysis can provide insights for future fault-tolerance research in the development of LLMs.

### 5.1 Failure Category

We employ a failure diagnosis system, leveraging a combination of rule-based and LLM techniques, to extract error information from the runtime logs. We provide detailed explanations of this system in §6.1. Besides, to ensure the accurate identification of the types and root causes of failures, manual checks and corrections are conducted. Table 3 provides a summary of common failures in Acme, including their occurrence frequency and restart time. Basically, they can be classified into three categories as follows. Note that these classifications may overlap, and the primary criterion for classifying a specific type of error is its most frequent occurrence.

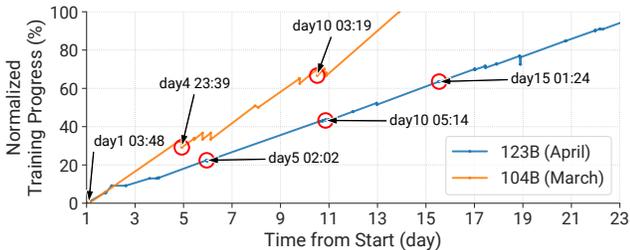
- **Infrastructure.** Infrastructure-related failures arise from issues within the underlying computational platform or remote storage. These failures mainly occur midway through the job execution process, especially in pretraining tasks. They severely impact the training progress due to laborious and time-consuming recovery process.
- **Framework.** Several types of runtime errors, such as *RuntimeError*, *ValueError*, and *AttributeError*, can be associated with tensor operations, shapes, data types, or unexpected behaviors. They are often observed in the initial phases of jobs and are typically resolved by fixing the configurations.
- **Script.** Script errors typically stem from programming errors or user oversights. They constitute the majority of failures and are often addressed by revising codes.

### 5.2 Failure Characterization

We highlight several key observations from our analysis: **Infrastructure Failures Cause Most Severe Impact.** As shown in Table 3, jobs that fail because of infrastructure is-

Category	Reason	Num	GPU Demand		Time to Failure (mins)		GPU Time (mins)		Time to Restart (mins)			Cluster
			Average	Median	Average	Median	Average	Total%	Average	Median	TR/TF%	
Infrastructure	NVLink Error	54	800	896	868.1	155.3	585683	30.25%	95.6	0.2	11.02%	S, K
	CUDA Error	21	847	1024	923.2	586.0	785099	15.77%	78.3	2.0	8.48%	S, K
	Node Failure	16	712	768	1288.8	535.8	934394	14.30%	102.8	21.5	7.98%	S
	ECC Error	12	680	512	1303.4	1192.3	958404	11.00%	2.8	1.8	0.21%	S, K
	Network Error	12	758	768	549.6	310.1	394821	4.53%	592.1	7.4	107.74%	S, K
	Connection Error	147	29	1	51.9	0.5	24492	3.44%	0.8	0.0	1.51%	S, K
	S3 Storage Error	10	422	256	2317.8	202.2	222151	2.12%	6.2	0.2	0.27%	S
	NCCL Timeout Error	6	596	512	159.7	48.1	86856	0.50%	66.7	43.6	41.78%	K
NCCL Remote Error	3	1152	1024	50.5	22.6	52419	0.15%	0.0	0.7	0.09%	K	
Framework	Dataloader Killed	6	445	508	1580.6	961.4	764170	4.38%	115.1	0.9	7.28%	K
	Attribute Error	67	228	8	67.8	1.2	60914	3.90%	2.4	0.0	3.58%	S, K
	Out of Memory Error	14	572	640	323.8	14.5	245278	3.28%	122.7	1.2	37.89%	S, K
	Runtime Error	65	441	352	66.4	3.9	27667	1.72%	10.9	1.5	16.41%	S, K
	Assertion Error	105	413	256	41.7	3.0	12315	1.24%	185.9	1.6	445.87%	S, K
	Value Error	33	387	256	9.9	3.7	5049	0.16%	27.4	0.6	276.74%	S, K
	Zero Division Error	5	499	256	14.5	15.6	5363	0.03%	2.5	1.1	17.31%	S, K
	Model Loading Error	104	8	8	2.6	2.6	20	0.00%	0.0	0.0	0.00%	K
Dataset Loading Error	5	1	1	1.6	1.6	1	0.00%	0.0	0.0	0.00%	K	
Script	File Not Found Error	568	21	1	14.2	0.4	5210	2.83%	0.4	0.0	2.58%	S, K
	OS Error	266	8	1	9.6	0.8	1098	0.28%	0.3	0.0	3.17%	S, K
	Type Error	620	18	4	0.9	0.3	97	0.06%	0.2	0.0	28.27%	S, K
	Name Error	18	247	24	3.2	0.5	947	0.02%	2.9	2.4	90.92%	S, K
	Permission Error	7	438	512	4.3	0.8	2131	0.01%	2.4	2.2	56.38%	S
	Import Error	111	93	8	1.1	0.4	74	0.01%	0.7	0.0	63.68%	S, K
	Key Error	260	7	0	3.0	1.6	55	0.01%	0.1	0.0	2.10%	S, K
	Syntax Error	10	391	384	0.7	0.6	348	0.00%	1.7	1.7	261.73%	S, K
	Argument Error	3	344	512	0.7	0.7	288	0.00%	2.7	0.7	408.47%	S
	Called Process Error	4	256	256	0.2	0.2	52	0.00%	11.7	10.9	5714.29%	S
Index Error	23	6	1	1.6	0.9	21	0.00%	0.8	0.0	49.73%	S, K	

**Table 3:** Job failure statistics. It is sorted based on **Total%** (i.e., the percentage of GPU time summation in different categories). **Num:** Number of Occurrence. **TF:** Time to Failure. **TR:** Time to Restart (i.e., Restart Timestamp – Failure Timestamp). **GPU Time:** TF×GPU Demand. **S/K:** Occurrence of errors in Seren/Kalos respectively.



**Figure 14:** The training progress of two LLMs, with special emphasis on the manual recovery at night.

sues often use a substantial number of GPUs (*GPU Demand*) and require considerable effort to restart (*Time to Restart*). They take over 82% GPU resources (*GPU Time*) with only 11% failed job quantity (*Num*). Most of these jobs are long-term pretraining tasks that can experience hardware failures multiple times, such as issues with GPU (e.g., *CUDAError*, *ECCError*), NVLink (*NVLinkError*), and network system (*NCCLRemoteError*, *S3StorageError*). Note that *NodeFailure* indicates uncategorized errors caused by unclear hardware issues. Addressing these infrastructure failures requires meticulous diagnostic efforts to pinpoint the source of the problems, often leading to the maintenance or replacement of defective hardware, which results in significant restart costs.

**Failures Caused by High Temperature.** Another notewor-

thy observation is that training 7B models in Kalos tends to result in GPU overheating, which can cause *NVLinkError* or *ECCError*. This phenomenon is largely due to the highly optimized communication cost, resulting in an exceptionally low GPU idle rate. We observe that the overall temperature in the cluster server room increased by approximately 5°C when training these models. Besides, we find most of these jobs occurred in July 2023, which is the hottest month on record [64]. This anomalous climate may be a potential cause of these failures, which is aligned with the finding recently reported by Microsoft [101]. We provide more detailed data on GPU temperature in Appendix A.4. Subsequently, our team enhanced the cooling capabilities of the cluster, leading to a significant reduction in the frequency of such failures.

**Many Failures Induced by Auxiliary Services.** In our pre-training framework, we connect to external components or services for metric reporting, logging, monitoring and alerting. These auxiliary services are vulnerable to network instabilities, potentially resulting in timeouts or failures that can decelerate or disrupt the training process. A significant number of *ConnectionError* and *NetworkError* incidents stem from these auxiliary services.

**Evaluation Jobs Rarely Encounter Errors.** In Kalos, only 6.7% of evaluation tasks encounter errors, and notably, there

are no recorded instances of GPU or NVLink failures. The low error rate may be attributed to their short duration and the resultant decreased stress on GPUs and NVLink connections. Consequently, this diminishes the chance of hardware and operational failures that are more frequent in pretraining jobs.

### 5.3 Failure Recovery

There are three scenarios where we should restart a job: (1) when an error occurs within the job, (2) when there are anomalies in training metrics such as a loss spike, and (3) when the training process is stuck. A ‘loss spike’ refers to a sudden increase in the loss that was previously decreasing normally, and does not recover over a certain period. Upon restarting, jobs revert to the last checkpoint, resulting in loss of training progress. Since existing LLM frameworks lack automatic recovery support, developers usually manually restart interrupted training jobs. Developers often need to be on call in turn to ensure timely completion of the pretraining model.

As shown in Figure 14, we select two pretraining jobs in the early stage (March to April) when we handle all failures manually. We extract information from the logs of two clusters’ large-scale model training processes, including the runtime duration of each submission, start and end times, and the initial and final iteration numbers of training. The 104B model is an early attempt when the framework is still under development. Consequently, the process of loading previous model checkpoints led to a substantial loss in the overall training process. Conversely, in the training of the 123B model a month later, we improved the framework and adopted smaller checkpoint save intervals. Moreover, we added a feature to gracefully terminate jobs, allowing for the preservation of current training results before ending the job. It is evident that the training process of the 123B model is more stable, with fewer losses incurred due to rollbacks. However, this progress came at a cost, as jobs that were interrupted at various times had to be rapidly restarted.

## 6 Deployed LLM Systems

As highlighted earlier, the development process of LLMs presents significant obstacles yet unveils viable strategies for overcoming these issues. This section will introduce our efforts in two stages: (1) *Pretraining*: enhancing fault tolerance through LLM-involved failure diagnosis and automatic recovery. (2) *Evaluation*: achieving prompt performance response via task decomposition.

### 6.1 Fault-tolerant Pretraining

**Motivation.** During LLM pretraining, failures are inevitable and frequently occur due to the substantial number of GPUs involved and the extensive duration of the training process [15, 46, 89, 97]. These failures can dramatically impede the training progress and lead to severe resource inefficiency (§5). Consequently, to minimize infrastructure downtime, it is common practice to assign on-call duties to address failures manually. This places a significant burden on engineers and re-

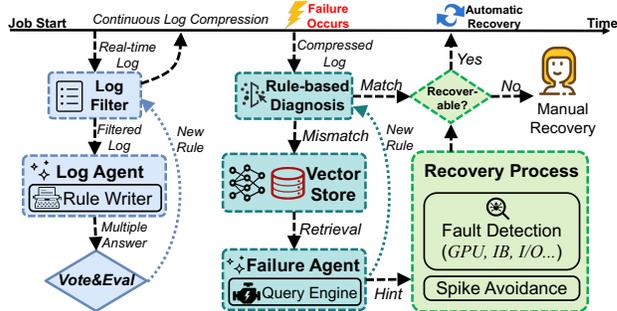


Figure 15: Workflow of failure diagnosis and model recovery.

searchers, as expressed in the complaints raised by the Meta OPT [111] and BigScience BLOOM [1] teams. Our team also faces this problem. To alleviate this burden and enhance hardware efficiency, we develop a system that automatically detects the root causes of faults and facilitates recovery.

**System Design.** Our fault tolerance system is seamlessly integrated into our LLM pretraining framework. It comprises three essential modules: (1) *Checkpointing*, achieving more frequent model saving to minimize the loss of training progress; (2) *Diagnosis*, using a combination of heuristic rules and LLM to identify the root cause of different failures accurately; (3) *Recovery*, employing a holistic detection toolkit to pinpoint fault nodes and automatically restart training from the properly saved checkpoint. We delve into them in detail.

**1. Asynchronous Checkpointing.** Frequent checkpointing efficiently mitigates the wasted time caused by unexpected faults [32]. However, as LLMs can produce TB-scale model states (referring to total model states across all GPUs), the process of saving checkpoints itself can introduce substantial overhead, resulting in training time slowdown up to 43% [61]. To tackle this problem, we adopt the asynchronous checkpointing strategy [65, 71], which effectively separates the checkpointing process from the training process. Our observations indicate that the CPU memory (refer to Figure 7 (b)) is capable of accommodating several checkpoints. By taking advantage of this, we can store the model state in memory and utilize a separate thread to regularly save these states to remote persistent storage. This simple strategy significantly reduces checkpointing overhead.

**2. Failure Diagnosis.** As we discussed in §5, failures can arise from numerous intricate factors, including errors from user script or framework, as well as issues with hardware subjected to high-stress conditions. It is crucial to determine whether a failure is recoverable for the purpose of automatic recovery. A common approach is to use a combination of heuristic rules to filter and conduct regular expression matching on the logs of faulty jobs [23, 47, 53, 54, 67]. However, this approach often proves inaccurate due to the wide-ranging diversity and complexity of error logs. There might not be a specific error statement in many cases, but multiple errors could coexist simultaneously. For example, a job might fail with messages that include *NCCLTimeoutError*, *CUDAError*, and multiple

kinds of *RuntimeError*, whereas the root cause is *CUDAError*. Trying to match every error scenario with a specific rule set can become impractical.

To address this challenge, we utilize the exceptional text understanding ability and extensive knowledge base of LLMs to identify the root causes of different failures automatically. As depicted in Figure 15, we incorporate an LLM with rule-based diagnosis to achieve efficient and accurate failure diagnosis. It mainly contains the following two steps:

► **Real-time Log Compression.** The extensive log files generated by pretraining jobs, primarily consisting of training metric records, can reach sizes of hundreds of MBs. To accelerate diagnosing and meet the context length limit of LLMs, log compression is conducted first. The system continuously updates a collection of regular expressions, termed as *Filter Rules*. These rules efficiently remove regular log outputs, such as initialization information, training metric records, framework outputs, and debug information. A vital component of the system, the LLM-based *Log Agent*, is responsible for analyzing real-time generated log segments and identifying lines that follow fixed patterns. By doing so, the LLM-based *Log Agent* dynamically writes regular expressions to update the *Filter Rules*, effectively minimizing the size of the log files. Additionally, the *Log Agent* forwards identified error messages to subsequent modules for diagnosis.

Furthermore, we employ the self-consistency [96] approach to ensure the robustness of the *Log Agent*'s results and to guarantee the formatting of these results. This involves processing each log segment multiple times and having another LLM vote on multiple results from the *Log Agent*, ensuring the accuracy of matches through regular expressions. Over time, the *Filter Rules* become more comprehensive for the current task, making the log filtering process more efficient. Furthermore, the system can utilize metadata from tasks to identify repetitive or similar tasks, directly applying existing *Filter Rules* for log filtering, thereby avoiding redundant work. This feature is particularly beneficial in large model cluster environments, where fewer tenants and task resubmissions are common.

► **LLM-assisted Automated Diagnosis.** The *Log Agent* efficiently compresses runtime logs, isolating critical error logs like *CUDAErrors* or runtime exceptions. Though logs are already compressed upon arrival at this module, error logs may still be lengthy. We apply a two-step approach to tackle this issue. First, the error logs are compared against a rule set that has been defined over time through the diagnosis of errors from past failed jobs. If the pre-defined rules fail to diagnose the issue, the compressed log is vectorized through an embedding model and stored in a vector store, serving as a retrieval repository. Then, the *Failure Agent* intervenes. It utilizes a *Query Engine* [56] to search through the vector store. Through this search, the *Failure Agent* can identify log lines that reflect the root cause of job interruption, extract the type of error, and indicate whether the error originated from user mistakes or infrastructure failures, providing a hint for the

recovery process. In addition, it also generates a mitigation suggestion for users or the operation team.

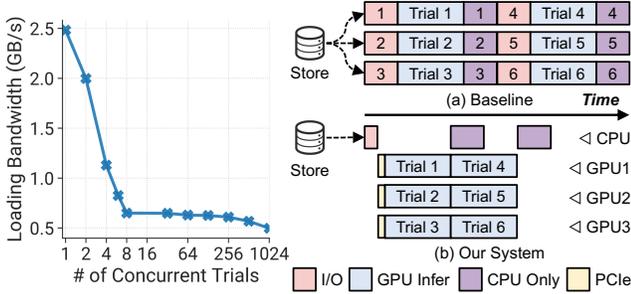
The *Failure Agent* also contributes to the continuous learning of the failure diagnosis system. For each new failure, once diagnosed and resolved, the *Failure Agent* writes a corresponding regular expression and adds it to the *Rule-based Diagnosis* module. This process is iterative and ensures that the *Failure Diagnosis* system evolves, becoming more adept at diagnosing and suggesting mitigation methods for failures. To achieve more robust performance, we currently utilize the GPT-4 [2] for diagnosis, with plans to transition to our LLMs.

**3. Fast Fault Detection and Recovery.** Based on the failure diagnosis result, if it belongs to one kind of infrastructure failure, we conduct a corresponding detection test to identify the problematic nodes. For instance, to promptly resolve the frequent *NVLinkError*, we employ a two-round NCCL test [5] approach. First, we divide all nodes into multiple two-node worlds and execute `allgather` task in each pair. If the total number of servers is odd, we leave one world size as three. If `allgather` task fails in a world, the nodes in that world are potentially faulty nodes. Then, in the second round, we pair potential faulty nodes with normal nodes to form new worlds. The nodes in each world continue to execute the `allgather` task, thus identifying the faulty nodes and then cordoning them off. On the other hand, if the failure is attributed to a sudden increase in loss (i.e., ‘loss spike’ [27, 111]), which is automatically triggered by our pretraining framework, we opt to an earlier healthy restart checkpoint and bypass subsequent data batches. This approach effectively maintains model quality.

**System Performance.** Our asynchronous checkpointing strategy offers a substantial reduction in checkpointing overheads, as the checkpointing process does not block the training process. The checkpoint time and overhead percentage of 7B and 123B size models are reduced by 3.6~58.7× (interval=30 mins), respectively. Note that the time taken for persisting to storage is not included in asynchronous checkpointing measurement. Moreover, our failure diagnosis system significantly reduces manual intervention by around 90% and thus reduces developers’ burden.

## 6.2 Decoupled Scheduling for Evaluation

**Motivation.** Evaluating the quality of LLMs based solely on a single metric, such as training loss, may not provide an accurate assessment [59]. Therefore, it is vital to incorporate a variety of criteria and evaluate performance across an array of tasks [22]. Our LLM framework conducts regular evaluations for every checkpoint during the pretraining phase in our data-center. This allows developers to track the progress of model training and identify the optimal model checkpoint. We aim for swift feedback to facilitate timely adjustment. However, as shown in Figure 6, evaluation jobs experience the longest queuing delay due to limited resources and concurrent submission of numerous trials. Despite these challenges, we identify



**Figure 16:** *Left:* Stress testing of model loading from remote storage in Seren. Each trial involves one GPU. *Right:* Scheduling evaluation trials. (a) Baseline: each dataset is treated as a trial. (b) Our system: decoupled scheduling.

several opportunities to expedite the evaluation process.

**System Design.** We develop a *trial coordinator* to harmonize the operations of the cluster scheduler and LLM framework. This design incorporates the following three key techniques aimed at enhancing the efficiency of the evaluation process.

**1. Decoupling Remote Model Loading.** Given the substantial size of LLMs, retrieving and loading them from remote storage can be a lengthy process. Furthermore, the concurrent execution of numerous evaluation tasks (around 60 datasets) can exacerbate this loading process due to increased contention. Figure 16 (Left) shows the average model loading speed on a range of concurrent evaluation trials within Seren. It reveals a huge decline in loading speed when increasing the number of single-GPU trials from 1 to 8 on a single node, due to the bandwidth limitation (25Gb/s) of our storage NIC. On the other hand, the loading speed stabilizes when the number of trials ranges from 8 to 256 GPUs. This observation inspires us to take a strategic approach. Rather than submitting each evaluation dataset as a separate trial, we separated the model loading process from the evaluation process, as depicted in Figure 16 (Right). Specifically, the trial coordinator initially retrieves the available node list from the cluster scheduler and then generates a series of precursor jobs for each node. These jobs load the model from remote storage to local shared memory. Following this, the coordinator submits the evaluation jobs to the scheduler, which loads the model via the high-bandwidth PCIe. This method effectively utilizes spare host memory. After the evaluation finishes, the coordinator clears the files.

**2. Decoupling Metric Computation.** As shown in Figure 13, the evaluation process can often involve complex and time-consuming metric computation. For example, synthesized program correctness tests must be performed on coding datasets like HumanEval [24] and MBPP [17]. To address this issue, we decouple the metric computation process from the evaluation trial. After the model inference is performed on the GPU, its output is quickly saved into files, terminating the inference workload. Given that the outputs are typically text-based and thus small in size, this file-dumping process is swift. We then generate CPU jobs to carry out the metric

computations. This approach effectively minimizes GPU idle time and accelerates the evaluation.

**3. Prior-based Elastic Scheduling.** In addition to the decoupling approach, we notice that our prior knowledge regarding the approximate trial runtime for each evaluation dataset is quite robust. Furthermore, these datasets are flexible, allowing us to batch multiple sets into one trial to circumvent model loading. We can also break down large datasets and decouple metric computation. As a result, the trial coordinator can maximize GPU occupancy through decomposition, balance each GPU’s workload using prior information, and employ a round-robin allocation strategy on sorted job queues. Moreover, we prioritize evaluation trials with lengthy CPU metric computations in the job queue to better overlap its computation. This approach not only enhances workload balance but also minimizes trial switch overhead.

**System Performance.** We conducted a representative test of the trial coordinator using a typical evaluation job on a 7B size LLM, which involved evaluating the workload across 63 datasets. We measured the makespan necessary to complete all evaluation trials under two different conditions: a single node (representing limited resources) and four nodes (representing relatively ample resources). The trial coordinator can reduce the makespan by  $1.3\times$  and  $1.8\times$  respectively.

## 7 Discussion

**Scope Limitations.** Despite our best efforts to analyze the workloads in Acme, it is an inescapable reality that we cannot cover all types of workloads. Limitations include: (1) Our analysis focuses on the developmental process preceding model serving and Acme does not encompass any serving jobs (i.e., workload in the deployment stage). (2) We concentrate our analysis predominantly on GPU jobs, providing limited room for CPU jobs. (3) We mainly characterize transformer-based, decoder-only architecture models (GPT-3 [20] and LLaMA 2 [93]). For newer model architecture, we provide a simple characterization of the Mixture of Experts (MoE) model [85] in Appendix A.5. Other model architectures like the Multimodal LLM [78] fall outside our scope of analysis.

## 8 Conclusion

In summary, we analyze LLM workloads and resource utilization in our datacenter Acme, revealing unique features and challenges of LLM development, such as resource inefficiencies and failure impacts. We also uncover potential opportunities to optimize systems tailored for LLMs and introduce our efforts for pretraining and evaluation workloads. We believe that our lessons and insights have broad applicability and can well benefit subsequent research.

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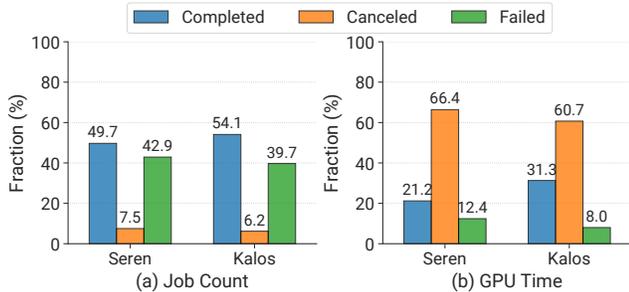
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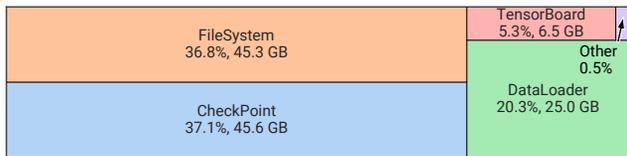
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**Figure 17:** Final statuses of jobs in terms of (a) quantity and (b) utilized GPU resources.



**Figure 18:** Distribution of host memory on a server in Seren during executing a pretraining job. Idle memory is not shown.

## A Supplementary Characterization

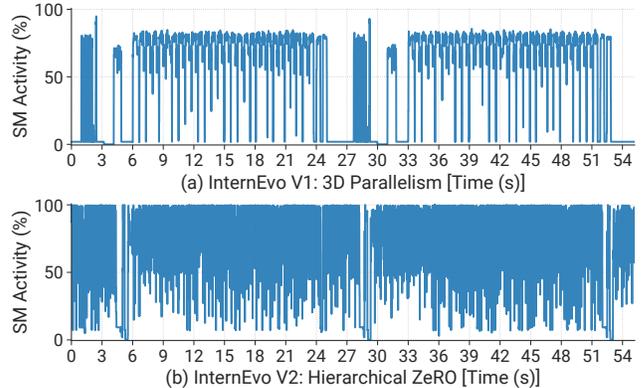
In this section, we provide additional analysis to further characterize the workload features during our LLM development.

### A.1 Job Final Statuses

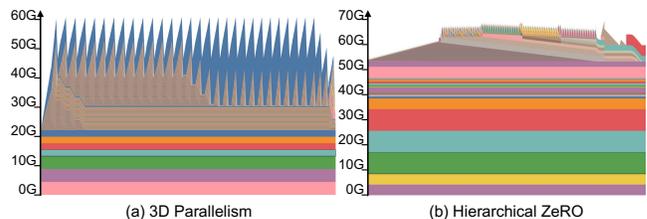
**High Incompletion Rate.** Figure 17 summarizes the distribution of three key final statuses across our two clusters, revealing a similar pattern. It is obvious that only approximately 20~30% of resources are consumed by jobs that finally complete. Besides, about 40% of jobs fail, utilizing 10% of GPU resources. This suggests that failures predominantly occur in the early stages of execution, aligning with the statistics presented in Table 3. Canceled jobs, while constituting only around 7% of the total job count, command over 60% of GPU resources. This pattern suggests a prevalence of large-scale pretraining jobs being canceled by users. Beyond the common cancellation motives cited in prior studies (e.g., achieving desired model performance sooner than expected, early recognition of poor hyperparameter configuration) [41, 75, 99], our experience with LLM pretraining has identified two additional frequent causes: (1) Users pausing jobs to adjust parameters in response to performance anomalies, such as loss spikes. (2) Jobs stalling due to infrastructure issues without throwing error messages, only to be addressed upon manual inspection by users, leading to significant resource wastage. These observations underscore the necessity for a failure-handling system that can autonomously detect and rectify faulty jobs, which is further elaborated in §5.3 and §6.1.

### A.2 Host Memory

**Memory Footprint Breakdown.** As depicted in Figure 18, we illustrate the distribution of active physical host memory within a compute node, which utilizes 123GB of the total 1TB available. This showcases a typical pattern of CPU memory

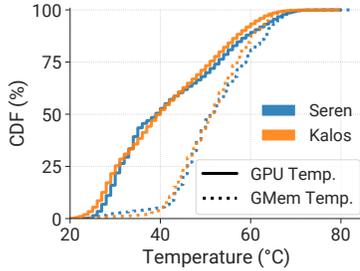


**Figure 19:** GPU SM utilization of pretraining a 123B LLM using different strategies of InternEvo [25] over 1024 GPUs.

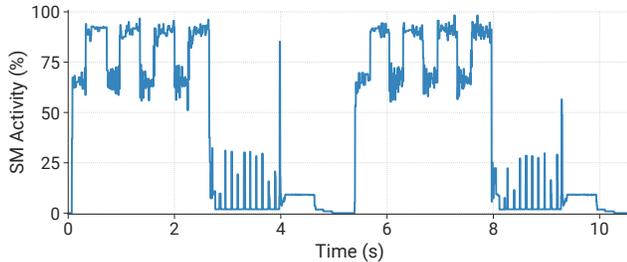


**Figure 20:** Memory snapshot under different pretraining strategies. Note that the extensive blue segment at the top of (a) is simplified and can be further broken down into massive fragments (memory allocations), similar to the lower part.

consumption for pretraining jobs. However, it is important to note that memory usage can significantly vary across different tasks. Specifically, the memory footprint of the dataloader can be considerably larger when employing Megatron-LM [69], which requires loading the metadata of the entire dataset. In contrast, our approach of loading data on-the-fly proves to be more memory-efficient without obviously impacting throughput. Furthermore, the memory requirements for asynchronous checkpointing (§6.1) largely depend on the model size and training configurations. The memory footprint depicted in this figure corresponds to the configuration outlined in Figure 10(a). In addition to the training processes, memory-intensive operations include TensorBoard [13] (6.5GB), the client daemon along the critical components (e.g., data and metadata caching) of the distributed file system (45.3GB). The remainder (0.6GB) encompasses Prometheus monitoring components, NVIDIA drivers, the Slurm scheduler daemon, and other system processes primarily related to sensor monitoring and system management. In general, there is a substantial amount of memory available, which can be leveraged for various purposes. Previous efforts [16, 66] have shown the potential for disaggregating CPU and memory usage from GPU allocations, suggesting that there may be additional optimization opportunities for LLMs, such as enhancing fault tolerance [97].



**Figure 21:** CDF of GPU core and GPU memory temperature.



**Figure 22:** GPU SM utilization of pretraining a MoE model Mistral 7B [48] with 1024 GPUs in Seren.

### A.3 Pretraining under Different Scale

Figures 19 and 20 provide supplementary pretraining profiling results for a 123B LLM across 1024 GPUs. These figures complement the profiling observations depicted in Figures 10 and 11. It is evident that they present very similar patterns to the 2048 GPUs, demonstrating the generalizability of our characterization.

### A.4 GPU Temperature

Figure 21 depicts the temperature distributions of GPU core and GPU memory. The GPU memory temperature is generally higher than the GPU core temperature. As corresponding to power consumption distribution shown in Figure 8, temperature presents a similar pattern in that some GPUs are under heavy load and have higher temperature (over 65°C). These suggest a need for enhancements in our cluster’s cooling system to address the issue of elevated temperatures.

### A.5 MoE Model

As shown in Figure 22, the MoE model presents much lower GPU utilization compared with the dense model shown in Figure 10. This is mainly due to the fact that the MoE model requires frequent all-to-all communication and necessitates high-quality internode communication, however, our single IB NIC server cannot efficiently handle such job. Here we directly use the official training configuration released by Mistral. On the other hand, InternEvo is still under development and we are working on performing tailored optimizations for MoE models.

## B Related Work

**DL Workload Characterization.** Prior works conduct DL workloads analysis from different companies. Philly [47] pro-

vides insights on the impact of job locality on queuing delay and resource utilization from Microsoft, in addition to identifying different failure reasons. Helios [41] from SenseTime illustrates the nature of cluster resource utilization and user disparity, evaluated from the perspectives of the cluster, job, and user. Alibaba’s PAI [98] contributes to this discourse by analyzing the challenges encountered within their clusters from both temporal and spatial viewpoints. Different from them, we focus on the characteristics of LLM workloads.

**Fault Tolerance Systems.** Fault tolerance is a crucial consideration across various disciplines. In the context of LLM systems, Varuna [15], Bamboo [89], and Oobleck [46] focus on the fast recovery from failures in cloud spot instance scenarios. Gemini [97] facilitates swift recovery through CPU memory checkpointing. In addition, a body of research works dedicated to GPU-related failure analysis [35, 47, 61, 72, 87, 90, 91, 110], while several deep learning schedulers [44, 57, 108] consider fault tolerance. Furthermore, a series of studies have profiled [33, 34, 50, 63, 70, 112] or diagnosed [36, 38, 39, 60, 88, 106, 109] potential performance bottlenecks within RDMA or intra-host network communication. We provide a thorough analysis of the failure events in LLM workloads and propose a LLM-involved diagnosis system.

## C Resource Links

### InternLM Links

**Project:** <https://github.com/InternLM>  
**Trace:** <https://github.com/InternLM/AcmeTrace>  
**System:** <https://github.com/InternLM/InternEvo>  
**Model:** <https://huggingface.co/internlm>