

Mycelium/Polymer

*Creating a pseudo-instruction-set for scalable distributed computing*

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## **Introduction**

Composability is the new hotness in the compute sphere. Admins find it convenient to isolate the finite resource groups of a computer, collect them in bulk, then slice them back up into VMs. Allocate rack space to high disk capacity. Create a section of the building for GPU compute. When the time comes, compose a virtual machine which has a slice of the storage and some of the compute. Every action is instant. A VM can be spun up or terminated in the scope of single moments or actions; capacity is scaled to match the demand of end users.

Companies like Liquid have brought us immeasurable levels of technological advancement. Their composable infrastructure offers unparalleled flexibility. Through bridging via systems' native PCI-Express interfaces, entire buildings of systems can communicate as if they were all within one server chassis. This approach strays from a more conventional system of hypervisor nodes, and is more rooted in hardware design and low-level resources sharing. Liquid has created a system of addressing hardware which transcends the boundaries of individual systems. It truly breaks a computing environment down into its raw resources. Then, provided only a simple BOM of required specs, a new computer can be conjured from the resource pool. This is the future of the server space—a future I hope home labs will soon share. In the meantime, I want for a legacy-compatible option. Something to run on my ever-growing collection of X86 dinosaurs.

Aging penguins rejoice. A solution may lie in an abstraction of the way we approach compute problems. Assuming an adequate instruction set is available, we can still break down and compile our applications into a distributed system of thinking. This theoretical instruction set could breath new life into compute clusters of old, who now lack some features of bleeding-edge systems.

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## Prior Developments

### Mainframes

A centralized paradigm, mainframe systems from IBM and other business partners allowed one large shared computer to be used by several employees or students. Each user could create a new session with the mainframe, share its resource, then log out when completed. As the cost of individual computing systems was vast, a large central system was one of the most affordable options.

Sharing resources in such a way has obvious drawbacks. If the mainframe had downtime or needed repairs, entire businesses or communities could be without access to their work. During certain times of the day, employees may find it difficult to access the mainframe. They'll have to wait until after peak usage time to upload their deliverables.



*(above: A modern IBM mainframe)*

### Personal Systems

To remediate some issues of having one large shared system, companies and schools opted for larger numbers of smaller and less powerful individual computers. These personal systems may still have shared resources (like company storage or printers)—or may be hooked into a managed company network—but each device feature its own capabilities and approaches problems from a much more individualistic view.

These devices could be any business laptops, school tablets, or other devices provided by an IT service to a student or employee.

In this case, as with all the others, issues will arise. Individual systems will age and need repair (then eventual replacement). Companies will often neglect their fleet, resulting in an entire network of outdated and insecure personal systems. Oversight of the fleet can also be time consuming.

While individual devices are convenient for employees and students, the upfront cost of registering a new device can be quite high. For an engineering firm, a new workstation and suite of engineering software can cost several thousands of dollars. Networked and mainframe solutions can offer better company-wide integration and stricter management of users. Personal systems can usually only be viable if connected to a shared network and closely maintained by IT personnel. With a centralized approach, users can be added and removed with improved ease and granularity.

## **Virtualization**

VMs, a generalized term for several virtualization and containerization technologies in server and user space, are a method of centralizing and consolidating the assets of a company's compute infrastructure. One central system serves the fleet of users. Many of the restraints of mainframes had been overcome through brute compute force. Machines can be spun up and terminated with the flow of demand from employees. And, as technologies improve, virtual hosts can be upgraded and improved to give users an improved experience.

For the purposes of this consideration, I have grouped containerization in with virtualization. Many will refer to containers as "light-weight" VMs. Usually this will manifest as a Linux hypervisor host (such as a VMWare or ProxMox instance) with Linux containers. Certain kernel features and resources of the hypervisor are shared with the containers to ensure they can run as lightweight as possible.

Via VMs and containers, tens or hundreds of systems can be run within one single server chassis. Several users can log in and access a seemingly dedicated machine. Nobody has to know they are sharing a single computer via virtualization and RDP.

## **Compute Clusters**

These clusters are aptly named. They are usually groups of peer machines or virtual machines. The group of peers can be referred to as a cluster. Sometimes each individual system or virtual system is called a node. For the purposes of this writing, I will refer to individual systems as nodes.

Compute clusters enable the technologies associated with high-availability (HA) systems. Since the compute solution is split across multiple nodes, there can exist failsafes and redundancies to allow near-perfect uptimes. When it comes time to service or upgrade a server node, changes can often be made with reduced downtime.

Individual nodes in a cluster may feature specialties. For example, some nodes may feature increased graphical compute capability; others may focus on providing storage or database resources. As demand for different resources changes, nodes with different specialties (or offering different services to users) can be added or removed.

## **Composability**

As previously mentioned, composability shares common use-cases with virtualization. The difference here is that composable infrastructures are bare metal. There needn't be the overhead of a virtual machine. Tech media poster-child Liquid's approach is to do everything over PCI-Express. This effectively creates a computer spanning the whole floor of a datacenter.

Due to the sprawling nature of a composed system, there exist implicit power and hardware redundancies which can make this infrastructure more reliable and maintainable.

Is this something I will be able to use in my home lab? Or is this a hyperconverged enterprise solution for the technically elite? I fear the latter. Hobbyists need a comparable solution to serve the home.



(above: Liqud's logo)

Similarly to how Liqud has abstracted computer hardware to be composable, I think it could be possible to abstract clusters of Linux nodes to perform distributed tasks. In order to approach an implementation, we need to abstract the compute interface of a Linux system.

## Resource Abstraction

### Logic Procedures

Due to volatility of logic values and computational complexity of business or application logic, operations relating to one application process should be consolidated to the same node. This will allow the most seamless and least latent communication between that application's tasks.

However, duplicate logical procedures can be created on separate nodes.

These procedures are the CPU-side workloads which drive the application as a whole. These processes will decide what compute tasks should be run, what data will be inputted, and where to send the processed data.

### Compute Tasks

The computations performed on nodes in the cluster must have a defined input and output schema. Compute tasks are provided as a collection of instruction sets contingent upon a provided input data range. Each instance of the compute execution will provide its output to the invoker.

A creator of a compute task should be able to define how their task is executed. For instance, if they know the task contains a large number of floating point operations, the programmer may choose to target systems with graphical compute acceleration. If a task is dependent on high memory capacity, that may affect deployability to graphics or CPU compute nodes.

Each task's instructions are defined as a "kernel," similar to how CUDA and OpenCL approach instances of compute. Each unit of work is a kernel; it will be duplicated in parallel across the number of nodes necessary to meet the task's demand.

## **Data Storage**

## **Communication**

## **Synchronization**

## **Optimization**

### **Out-of-order Execution**

Temporal coupling occurs within a computation when tasks must be completed in a certain specific order. If executed temporally incorrect, the output may result in a crash or incorrect data. The executable format of the abstract instruction set should calculate a topological graph of which tasks are reliant on the result of previous tasks, grouping mutually exclusive operations into frames which can be executed in parallel.

This can allow extra system resources to be used to accelerate a singly threaded operation. The singly threaded logic task can be broken down into sequences of buckets of mutually exclusive elements. Executed in order, the end result of the accelerated computation should be equal to the un-accelerated output. I recognize a similarity to AVX extensions, which use parallelized vector silicon to unroll loops in the code and accelerate them.

## **Speculation**