

[0.3] Optimization, Hyperparameters, and Scale

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ARENA

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 - ② Experiment tracking + hyperparameter sweeps with Weights & Biases.
 - ③ Distributed training — when one GPU isn't enough.

The training loop

- Same four steps every iteration: *forward, loss, backward, step*.

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for x, y in dataloader:
    y_hat = model(x)
    loss = loss_fn(y, y_hat)
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- `zero_grad()` before backward: gradients **accumulate** by default in PyTorch.

Loss landscapes & why this is hard

- Loss is a function $\mathcal{L} : \mathbb{R}^d \rightarrow \mathbb{R}$ over millions of parameters. We never see the full thing — only 1D/2D slices through it.

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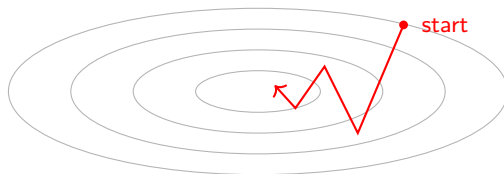
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 - **Plateaus:** gradient is tiny; nothing moves.



ill-conditioned “ravine” — SGD zigzags

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Gradient descent recap

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- In practice we don't compute $\nabla \mathcal{L}$ on the full dataset — we use **batches**. The gradient becomes a noisy estimate, but each step is much cheaper.

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- **Batch size:**
 - Larger $\rightarrow$ (+) lower-variance gradient, better GPU utilization
 - Larger $\rightarrow$ (-) more memory, fewer optimizer steps per epoch, often worse generalization at scale

SGD

$$\theta_{t+1} = \theta_t - \alpha g_t$$

```
class SGD(Optimizer):
    def __init__(self, params, lr):
        self.params = list(params)
        self.lr = lr

    @torch.no_grad()
    def step(self):
        for p in self.params:
            p = p - self.lr * p.grad  # SGD update

    def zero_grad(self):
        for p in self.params:
            p.grad.zero_()
```

Momentum

SGD with momentum

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- distill.pub/2017/momentum has a beautiful interactive on this.

RMSprop

$$s_{t+1} = \beta s_t + (1 - \beta) (\nabla_{\theta} \mathcal{L})^2$$
$$\theta_{t+1} = \theta_t - \alpha \frac{\nabla_{\theta} \mathcal{L}}{\sqrt{s_{t+1}} + \varepsilon}$$

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- **Memory cost:** s_t is parameter-sized $\Rightarrow 2\times$ memory vs SGD.

Adam

$m_{t+1} = \beta_1 m_t + (1 - \beta_1) \nabla \mathcal{L}$	(first moment)
$v_{t+1} = \beta_2 v_t + (1 - \beta_2) (\nabla \mathcal{L})^2$	(second moment)
$\hat{m}_{t+1} = m_{t+1} / (1 - \beta_1^{t+1})$	(bias correction)
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Adam vs AdamW: the weight-decay subtlety

- **Naive weight decay:** add $\frac{\lambda}{2} \|\theta\|^2$ to the loss. Looks innocent, but for adaptive methods, the L2 gradient $\lambda\theta$ gets divided by $\sqrt{\hat{v}_t}$ along with everything else — so parameters with large historical gradients are decayed less. Not what you want.

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AdamW: decoupled weight decay

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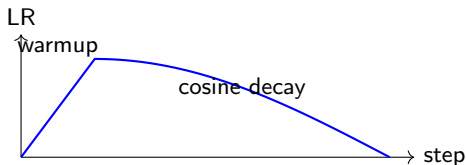
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- In practice: just use AdamW.

Schedules & parameter groups

- Fixed LR is fine for toy problems. Real training uses a **schedule**: warmup (linear ramp from 0) + decay (cosine or linear).

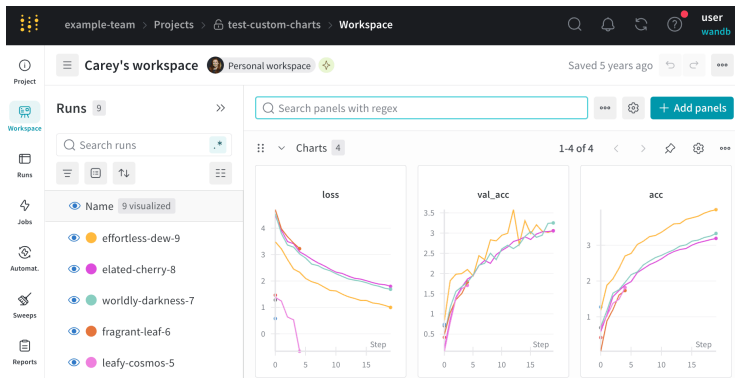


- Parameter groups:** different learning rates for different parts of the model. Common in fine-tuning (slow backbone, fast head):

```
optimizer = torch.optim.AdamW([  
    {"params": backbone.parameters(), "lr": 1e-5, "weight_decay": 0.01},  
    {"params": head.parameters(), "lr": 1e-3, "weight_decay": 0.0},  
)
```

Why track experiments? Weights & Biases

- WandB is the de-facto tool: can easily track and log loss, accuracy etc. while the model is training

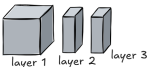


Parallel processes:



(each represents computations happening on a different GPU)

Model:



(each layer is a sequential step)

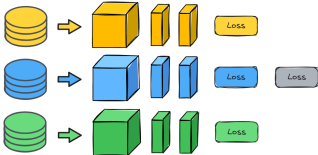
Pipeline Parallelism / Vertical Model Parallelism

- Splitting model into serialized stages, e.g. layers
- Requires careful management of inter-stage latency
- Natively supported by PyTorch: `torch.distributed.pipeline`



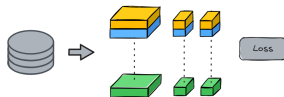
Data Parallelism

- Splitting batches of data across processes
- Easiest form to implement & manage; supported by PyTorch



Tensor Parallelism / Horizontal Model Parallelism

- Splitting model into parallel stages (often splitting parameters into shards)
- Most complex to implement (requires fine-grained control over computations)
- Experimentally supported by PyTorch: `torch.distributed.tensor.parallel`



- **Optimizer**: the rule for turning $\nabla \mathcal{L}$ into a parameter update.
- **Learning rate** α : step size. The most important dial.
- **Momentum** (μ): exponentially weighted average of past gradients.
- **Weight decay** (λ): shrink parameters toward zero each step.
- **Adam**: momentum + per-parameter adaptive LR + bias correction.
- **AdamW**: Adam with decoupled weight decay. Use by default.
- **LR schedule**: warmup then decay; pretty much always.
- **Parameter group**: subset of parameters with its own LR / weight decay.
- **WandB**: experiment-tracking dashboard
- **Sweep**: automated hyperparameter search