Rotational Equilibrium: How Weight Decay Balances Learning Across Neural Networks

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Abstract

Weight decay can significantly impact the optimization dynamics of deep neural networks. In certain situations the effects of weight decay and gradient updates on the magnitude of a parameter vector cancel out on average, forming a state known as equilibrium. This causes the expected rotation of the vector in each update to remain constant along with its magnitude. Importantly, equilibrium can arise independently for the weight vectors of different layers and neurons. These equilibria are highly homogeneous for some optimizer and normalization configurations, effectively balancing the average rotation—a proxy for the effective learning rate—across network components. In this work we explore the equilibrium states of multiple optimizers including AdamW and SGD with momentum, providing insights into interactions between the learning rate, weight decay, initialization, normalization and learning rate schedule. We show how rotational equilibrium can be enforced throughout training, eliminating the chaotic transient phase corresponding to the transition towards equilibrium, thus simplifying the training dynamics. Finally, we show that rotational behavior may play a key role in the effectiveness of AdamW compared to Adam with L_2-regularization, the performance of different normalization layers, and the need for learning rate warmup.

1 Introduction

Efficient training of deep neural networks intuitively requires the learning of different components (layers, neurons etc.) to be balanced in some sense. A layer that is updated slowly, perhaps barely changing through the training process, is unlikely to contribute optimally to the final model resulting in worse performance and wasted compute. Conversely, a rapidly changing layer may cause instability, limiting the maximum stable learning rate and preventing other layers from learning effectively. This suggests that sufficiently imbalanced updates can hinder training.

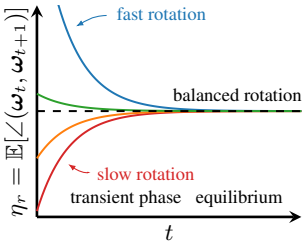

Figure 1: Conceptual figure of the angular updates of the weight vector $\boldsymbol{\omega}_t$ for different normalized neurons (each line color) over time t with a constant learning rate.

In this work we explore how weight decay can balance the *average angular update* $\eta_r = \mathbb{E}[\angle(\boldsymbol{\omega}_t, \boldsymbol{\omega}_{t+1})]$ of different weight vectors $\boldsymbol{\omega}$ used for dot products (e.g. the weights of one neuron, but not the bias). When weight decay is applied to such vectors, it can interact with normalization layers and gradient noise to cause the norm $\|\boldsymbol{\omega}\|$ to converge to a stable equilibrium value $\widehat{\|\boldsymbol{\omega}\|}$ via Spherical Motion Dynamics [35]. This process causes η_r to converge to a stable value $\widehat{\eta_r}$ as well after an *initial transient* phase (see Fig 1). For some optimizers and normalization setups this value is identical across all or most $\boldsymbol{\omega}$, resulting in *balanced rotation*. We extend prior research

*Equal contribution. Our code is available at https://github.com/epfml/rotational-optimizers and an extended version of this work is accessible at https://arxiv.org/abs/2305.17212.

Mathematics of Modern Machine Learning Workshop at NeurIPS 2023.

Table 1: Update and equilibrium norms for a parameter $\boldsymbol{\omega} \in \mathbb{R}^C$ for different optimizers. The RMS update size η_g is for all parameters, $\widehat{\eta_r}$ and $\widehat{\|\boldsymbol{\omega}\|}$ only apply to a scale-invariant $\boldsymbol{\omega}$ in equilibrium.

	RMS update size $\widehat{\eta_g}$	Expected angular update $\widehat{\eta_r}$	Equilibrium norm $\widehat{\|\boldsymbol{\omega}\|}$
SGDM	$\eta\sqrt{\frac{\mathbb{E}[\|\boldsymbol{g}\|^2]}{1-\alpha^2}}$	$\sqrt{\frac{2\eta\lambda}{1+\alpha}}$	$\sqrt[4]{\frac{\eta\mathbb{E}[\|\tilde{\boldsymbol{g}}\|^2]}{2\lambda\cdot(1-\alpha)}}$
AdamW	$\eta\sqrt{C\frac{1-\beta_1}{1+\beta_1}}$	$\sqrt{2\eta\lambda\frac{1-\beta_1}{1+\beta_1}}$	$\sqrt{\frac{\eta C}{2\lambda}}$

in this area (Appx B) to new optimizers like AdamW [21] and investigate the impact of balanced rotation when training deep neural networks, demonstrating its importance for the effectiveness of multiple methods.

2 Methods & Analysis

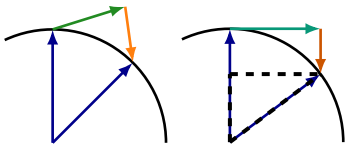

Figure 2: Equilibrium perspectives

The main goal of our analysis is to obtain expressions for $\widehat{\|\boldsymbol{\omega}\|}$, $\widehat{\eta_r}$ and $\widehat{\eta_g}$, the predicted root-expected-square size of a parameter update arising from the gradient (not the weight decay). We do this by analysing the geometry of a neural network undergoing a random walk. The left side of Fig 2 shows a conceptual view of equilibrium where the weight magnitude (blue) is preserved as the gradient component (green) and weight decay (or L_2 regularization, orange) component of an update balance out on average. To simplify the effects of momentum, we construct an alternative view (Fig 2 right) by considering the total weight change throughout training from the gradient and L_2 regularization at a given time step. This makes the components orthogonal in expectation, allowing us to apply the Pythargorean theorem to the dashed triangle and then solving for the equilibrium weight magnitude $\widehat{\|\boldsymbol{\omega}\|}$. Combined with an expression for the absolute size of the update $\widehat{\eta_g}$ for this weight magnitude, we obtain $\widehat{\eta_r}$. See Table 1 for predictions, Appx E for the setup, and F-I for specific optimizers.

Scale-invariance: In real optimization problems (i.e. not a random walk) the gradient can have an average component in the direction of the weights (a radial component). We find that this component can act like an additional weight decay, resulting in rotational behavior similar to the random walk but for an adjusted weight decay value (see Appx L for details). Some weights $\boldsymbol{\omega}$ are scale-invariant w.r.t. the loss function $\mathscr{L}(\boldsymbol{\omega}, \ldots)$, meaning that $\mathscr{L}(r\boldsymbol{\omega}, \ldots) = \mathscr{L}(\boldsymbol{\omega}, \ldots), \forall r > 0$, typically due to normalization (Appx D). The gradient $\nabla_{\boldsymbol{\omega}}\mathscr{L}(\boldsymbol{\omega}, \ldots)$ for a scale-invariant vector $\boldsymbol{\omega}$ is orthogonal to the vector. This eliminates the radial gradient component resulting in a behavior better aligned with the random walk. Different normalization configurations can result in scale-invariance on a different granularity. For example when Batch Normalization is placed immediately after a layer, the weight vector of each neuron (filter) becomes scale-invariant on their own as opposed to Layer Normalization which only makes the whole matrix scale-invariant. We believe the primary advantage of Weight Standardization [27] (a.k.a. Centered Weight Normalization [10]) on top of Layer Normalization [1] or Group Normalization [38] is ensuring scale-invariance on a finer granularity, balancing rotation across neurons.

AdamW vs Adam+L_2: We find that Adam with L_2-regularization (Adam+L_2) does not result in balanced rotation, even in optimally normalized setups, unlike every other optimizer we considered. This is because the $\widehat{\eta_r}$ has a complicated dependency on the average gradient magnitude, which can vary across layers and neurons. We believe this could be the main reason for the effectiveness of the decoupled weight decay in AdamW [21] compared to L_2-regularization. See Appx I for the analysis and empirical evidence in our experiments section.

Rotational Optimizer Variants Standard optimizers only achieve balanced rotation after the weight norms converge to equilibrium. This creates a transient phase at the start of training and when hyperparameters change (e.g. step-wise change in the learning rate), where rotation can be imbalanced (see e.g. Fig 1) and far away from the equilibrium value, creating a mismatch between the specified learning rate schedule and the resulting effective step sizes over time. We propose Rotational Variants (RVs) of these optimizers that explicitly control the average angular update, forcing it to match equilibrium throughout training. We do this by making the updates orthogonal to the weights and then controlling the average magnitude of the update relative to the weights according to the values

Table 2: Test set performance using three different random seeds for the baseline optimizer, AdamW, and its rotational variant (RV). We use the baseline hyperparameters directly for the zero-shot results. For the best shot results, minor tuning was applied. The final column shows results for Adam+L_2 directional updates with a balanced angular update speed $\widehat{\eta}_r$ based on AdamW across all neurons.

Dataset	Model	Batch Size	Metric (⇅)	AdamW Baseline	RV-AdamW Zero Shot	RV-AdamW Best Shot	Wrapped Adam+L_2
CIFAR-10	ResNet-20	128	Top-1 Acc. (↑)	92.2±0.11	92.3±0.25	N/A	92.3±0.18
CIFAR-10	ResNet-20	2048	Top-1 Acc. (↑)	91.5±0.22	91.2±0.21	91.9 ±0.29	91.8±0.15
CIFAR-10	DeiT tiny	64	Top-1 Acc. (↑)	95.9±0.07	96.3±0.25	N/A	96.3±0.17
Imagenet-1k	DeiT tiny	1024	Top-1 Acc. (↑)	72.1	71.5	72.3	N/A
IWSLT2014 de-en	Transformer-S	4096	Bleu (↑)	34.6±0.06	19.9±0.14	34.7±0.10	34.5±0.04
Wikitext	GPT-like	55	Perplexity (↓)	19.6±0.07	19.1±0.21	N/A	19.3±0.12

given in Table 1. This eliminates the transient phases completely and can balance the rotation of different layers and neurons without relying on normalization. By default we target the equilibrium dynamics of Weight Standardization [27]. We keep the weight magnitude constant and optionally introduce a learnable gain to compensate, which can matter for scale-sensitive weights and avoids numerical issues. The form of the rotational wrapper closely resembles that of relative optimizers like LARS [39] and others discussed in Appx B, revealing a connection to the equilibrium dynamics of standard optimizers. Further details and algorithm are given in Appx M.

3 Experiments

Experimental Setup: We conducted experiments on several well-known datasets and standard network architectures (details in Appendix O). When using our RVs, we apply rotational updates to all convolutional and linear layers. For transformers, these layers are not always fully scale-invariant; the weight norms matter. To address this, we introduced learnable gains as previously mentioned.

Constraining the Rotational Dynamics: We explore the impact of constraining the rotational dynamics by comparing the performance of AdamW and RV-AdamW across various network architectures and tasks. Table 2 lists results (mean±std) for three seeds (additional results for SGDM and Lion in Appendix N). The RVs achieve comparable performance to the original versions without any additional hyperparameter tuning (zero-shot) or light tuning (best-shot). This suggests that the simplified learning dynamics (e.g. no transient phase) with the RVs are sufficient. Although the RVs have many interesting properties and good performance, we view them primarily as a research tool.

Learning Rate vs Weight Decay: Different combinations of a learning rate η and weight decay λ with a constant product $\eta\lambda$ result in the same expected angular update $\widehat{\eta}$ in equilibrium but different RMS update sizes $\widehat{\eta}_g$ (see Table 1). This affects the rotational weights (especially when scale-invariant) differently than gains and biases (which have $\lambda=0$). Figure 3 (left) explores the impact of this. For small η, biases and gains update slowly since $\widehat{\eta}_g \propto \eta$, resulting in accuracy comparable to removing or freezing them (90.8%). Conversely, large η results in large updates to these parameters,

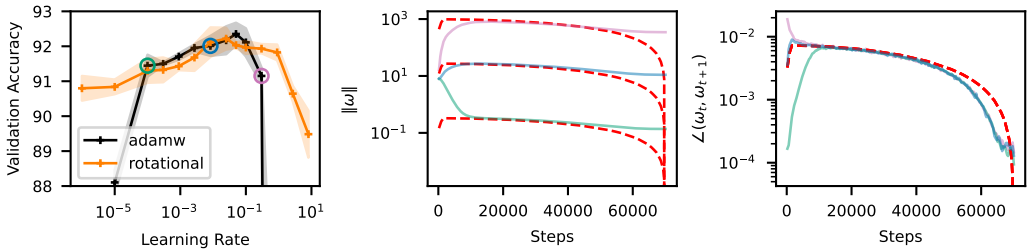

Figure 3: **Left:** Validation accuracy for ResNet-20 on CIFAR-10 for different learning rate, weight decay pairs with a constant product ($\eta\lambda = 5{\cdot}10^{-4}$) resulting in a specific $\widehat{\eta}_r$ (Table 1). **Right:** The weight norm $\|\boldsymbol{\omega}\|$ and angular update size η_r over time for three (η, λ) pairs corresponding to the colored circles on the left with predictions in dashed red. Note the difference in the initial/final phase.

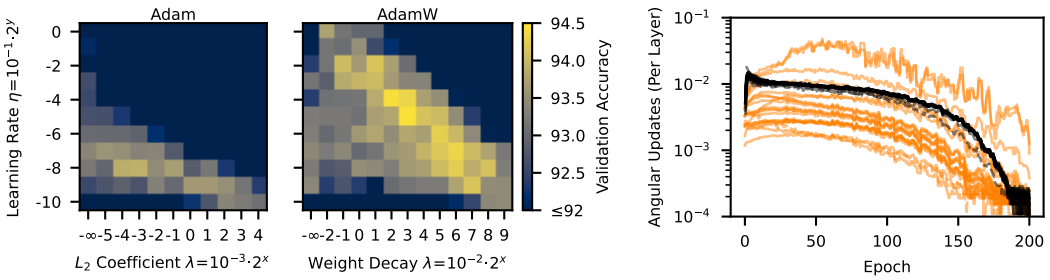

Figure 4: **Left:** A hyperparameter sweep for the learning rate and L_2 regularization / weight decay of Adam and AdamW on CIFAR-10 ResNet-18. Adam is unable to match the performance of AdamW. **Right:** The average angular per-step updates (radians) of each layer over the course of training for the best configuration of Adam (orange) and AdamW (black). In both cases the updates change with the learning rate schedule (cosine, no warmup) but for Adam they vary considerably across layers. The dashed line corresponds to the final fully-connected layer which is scale-sensitive.

potentially causing unstable training. We attribute the performance difference between AdamW and the RV to varying effective update size schedules (see Figure 3 middle and right).

Scheduling Effects: In Figure 3 (right), there are noticeable deviations between the measurements of $\|\boldsymbol{\omega}\|$ and η_r and the predictions listed in Table 1 for a cosine decay learning rate schedule and a five epoch warmup. In the initial phase we observe the transient phase, corresponding to the transition towards equilibrium, while the fall off in the end phase indicates that the learning rate is too small for the weights to decay fast enough to maintain equilibrium. The same effective step size schedule could be achieved by the RV with an adjusted learning rate schedule. See Appendix O for details.

Adam vs AdamW: Figure 4 (left) reproduces the performance gap between AdamW and Adam+L_2 observed by [21], around 0.5% on the validation set. The right panel shows that Adam+L_2 results in an unbalanced rotation unlike AdamW, confirming our observations from Section I.1. To determine whether this contributes to the performance gap, we create a special RV that combines the update direction $\Delta_g \boldsymbol{p}$ from Adam+L_2 with the $\widehat{\eta_r}$ of AdamW, ensuring balanced angular updates across all neurons and layers. For the experiment in Figure 4, the special RV performs identically to a standard RV-AdamW, outperforming Adam+L_2 by roughly 0.5%. In the final column of Table 2 we show that this holds for more settings, with the special RV and RV-AdamW performing similarly, roughly matching or outperforming AdamW in all cases.

Training Poorly Normalized Networks: Layer Normalization can make a whole layer scale-invariant but not individual neurons (unlike e.g. Batch Normalization). For standard optimizers, this can result in imbalanced rotation across neurons but the rotational optimizer variants ensure balanced rotation irrespective of the normalization. Figure 5 (left) shows that this results in improved performance across learning rates when training a layer-normalized ResNet-18 on CIFAR-100. We also observed a slight performance increase for the small GPT-like model in Table 1, potentially for a similar reason.

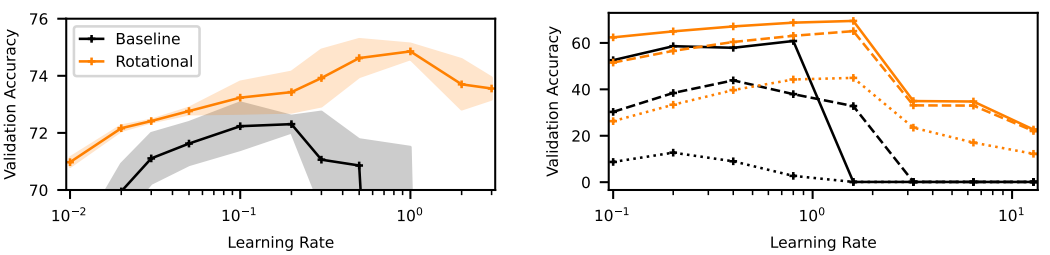

Figure 5: Comparison of the final validation accuracy of runs with different learning rates, using either SGDM (black) or RV-SGDM (orange). **Left:** Layer normalized ResNet-18 training on CIFAR-100, accuracy given as (mean±std of 5 runs). **Right:** 10 epoch training of ResNet-50 on ImageNet-1k without a learning rate warmup using large batch sizes (2k solid, 8k dashed, 32k dotted).

Need for Learning Rate Warmup: For standard optimizers the initial transient phase can result in imbalanced and overly fast rotation that does not match the learning rate schedule. We conjecture that the common practice of learning rate warmup is often beneficial in part by counteracting this effect. In Figure 5 (right) we train a ResNet-50 on ImageNet for 10 epochs using large batch sizes and different learning rates without warmup. The RV is more stable and achieves higher accuracy.

Imbalanced Rotation: In Appendix N we further quantify the impact of imbalanced rotation by artificially scaling η_r for a fraction of the neurons in each layer using a special RV. We find that even small variations in η_r can significantly affect performance.

4 Conclusion

We believe rotational dynamics can provide valuable insights into many phenomena in deep learning optimization. The learning rate and weight decay jointly determine two distinct effective step sizes used for different parameters and their schedule over time (along with the initialization and learning rate schedule). Balanced rotation seems to play a key role in the effectiveness of AdamW over Adam+L_2 and the benefit of Weight Standardization [27] or Batch Normalization [12] over e.g. Layer Normalization [1]. Learning rate warmup may aid training by stabilizing the chaotic transient phase where rotation can be imbalanced and/or overly fast. Although this may not be the sole reason a warmup is beneficial/needed in every case, we believe it is an important effect to be aware of.

5 Acknowledgements

We thank Maksym Andriushchenko for insightful conversations related to this study. We also appreciate Dongyang Fan and Nikita Doikov for their feedback on the manuscript which enhanced its quality and clarity.

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A Reproducibility

The code used for our experiments is available at `https://github.com/epfml/rotational-optimizers`. We provide more experimental details in Appendix O including hyperparameters not listed in the main body. The analysis for the other optimizers are also provided in the Appendix, SGDM in Appendix G, Lion in Appendix H and Adam+L_2 in Appendix I. Additional information about the analytical setting used in Section 2 and how it compares to practical settings can be found in Appendix K.

B Expanded Related Work

In this section we discuss more related works divided into four main categories.

B.1 Scale-invariance and Effective Learning Rates

Several works have investigated how the scale-invariance results in a certain "effective learning rate" based on the relative change that varies based on the norm of the weights, often in the form of $\eta/\|\boldsymbol{\omega}\|^2$. The works in this section do not describe how $\|\boldsymbol{\omega}\|$ can converge to an equilibrium value that results in a fixed relative or rotational learning rate. In Weight Normalization, Salimans and Kingma [31] point out how normalization can make parameters scale-invariant and that the gradient magnitude varies based on the weight magnitude. They describe how the gradient can "self-stabilize its norm", with larger gradients becoming smaller over time due to growth in the weight magnitude, but do not consider the effects of weight decay on this process. Zhang et al. [41] and Hoffer et al. [9] empirically find that the regularization effects of weight decay are primarily caused by increases in the effective learning rate due to decreased weight norms. Li and Arora [16] show that weight decay can be replaced by an exponentially increasing learning rate when optimizing scale-invariant weights with SGDM.

B.2 Equilibrium

These works also consider the fact that the weight norm converges to a specific value and the resulting effects on the relative update size. Van Laarhoven [34] points out the scale-invariance property of normalization and how it interacts with L2 regularization. They derive the $\eta/\|\boldsymbol{\omega}\|^2$ as the effective learning rate and also how there exists a fixed point where the weight norms are stable. Their work does not consider convergence of the weight magnitude as a separate process from the overall convergence of the loss and weights. In Online Normalization, Chiley et al. [6] show a simple derivation of the equilibrium condition in SGD and how it results in a relative update size that is identical across layers. The Spherical Motion Dynamics (SMD) [35] expands on prior work by proving the convergence of the weight norm and extending the analysis to include momentum. They also show plots of the weight norm over the course of training, providing empirical evidence for early convergence of the weight norm and also how it can fall out of equilibrium with sudden learning rate changes or when the learning rate becomes too small. They also consider the angular updates, empirically and analytically showing that they converge to an equilibrium value. Li et al. [17] analyze the convergence of SGD to equilibrium by modelling it as a stochastic differential equation, arriving at similar conclusion as the SMD paper (without momentum). This is expanded upon by Li et al. [19].

B.3 Projected Optimization

Some existing works are based on projections onto a sphere but do not scale the update to be proportional to the weight magnitude. Although similar to our Rotational Variants, this has a very different effect. Instead of rotating all scale-invariant groups at the same rate, they are kept at different rates based on their magnitude and resulting gradient norm where applicable. AdamP [8] orthogonalizes the update of the Adam and SGDM optimizers by removing the radial component of $\Delta_g\boldsymbol{\omega}$. The main reason for this is to avoid a rapid increase the weight norm during the initial phases of training. However, they use low amounts of weight decay if any which can also prevent the norms from growing as we have shown. Our Rotational Optimizer Variants keep the relative learning rate at a fixed level and do not suffer from the effect they report. Zhou et al. [42] propose keeping the weight magnitude constant, projecting it onto a sphere after every step and removing the

weight decay. Kodryan et al. [14] analyze training using projections onto the unit sphere after every optimizer update. They consider the union of all scale-invariant weights and do not normalize their step sizes so their effective learning rate per group can vary, giving dynamics that are generally quite different from SMD equilibrium. Roburin et al. [29] analyzes the spherical projection of the Adam optimization trajectory during standard training.

B.4 Relative Optimization

LARS [39] and LAMB [40] are variants of SGDM and AdamW that scale the update of a weight to be proportional to its norm (sometimes a clamped version of the weight norm). They apply this to linear and convolutional layer weights, keeping the original update for weights and biases. LARS and LAMB were proposed for large batch size training and found to work well there. Although they are not inspired by the Spherical Motion Dynamics, their form is quite similar to the Rotation Wrapper (Algorithm 1) with a few important distinctions. The default form of our Rotational Optimizer Variants (RVs) is applied filter-wise, centers the weights and allows the update magnitude to vary between steps while keeping the average relative update constant. The RV also doesn't apply weight decay while LARS and LAMB consider it a part of the update and take into account when scaling the relative update. Finally the RVs adjust the learning rate based on the SMD equilibrium value. This makes it more compatible with the underlying optimizer variants in terms of hyperparameters. One notable difference is the square root dependency on the relative updates in the SMD, while LARS and LAMB are directly proportional. This means that any learning rate schedule for these optimizers is more similar to applying a squared version of this schedule to standard optimizers or the RVs. This does not fully resolve the differences however, because changing the schedule also affects gains and biases where the update magnitude is directly proportional to the learning rate for all the optimizers and variants discussed here.

Nero [20] is another optimizer that applies relative updates that are directly proportional to the learning rate and weight magnitude. Like LARS and LAMB, Nero is not inspired by the SMD and to the best of our knowledge their relationship has not been pointed out before. Like the RVs, Nero is applied filter-wise and centers the weights. Overall, Nero is similar to the SGDM RV without momentum and the hyperparameter mapping, but also applies Adam like updates to the gains and biases, using a separate learning rate. By making the relative updates directly proportional to the learning rate, it has the same learning rate scheduling differences as LARS and LAMB mentioned above. Nero lacks momentum which is something that we observed can hurt large batch size training (exploratory experiments not shown).

Instead of controlling the average relative update size, Brock et al. [3] and Li et al. [18] clip the maximum relative update size instead. The Adaptive Gradient Clipping from Brock et al. [3] is applied on a per filter basis and is constant throughout training, i.e. does not vary with the learning rate or weight decay. The clipping introduced in Li et al. [18] scales with the learning rate and weight decay in a manner inspired by the equilibrium norm for SGD. They seem to apply this globally (i.e., not per neuron or layer).

C Preliminaries

C.1 Normalization and Scale-Invariance

We say that a weight vector $\boldsymbol{\omega}$ is *scale-invariant* with respect to the loss $\mathscr{L}(\boldsymbol{\omega}, \ldots)$, which generally depends on other network parameters too, if scaling it by a positive factor does not affect the loss, i.e. $\mathscr{L}(r\boldsymbol{\omega}, \ldots) = \mathscr{L}(\boldsymbol{\omega}, \ldots), \forall r > 0$. When placed correctly relative to $\boldsymbol{\omega}$, many normalization operations such as Batch Normalization [12], Layer Normalization [1] and more [11, 10, 38, 27] result in approximate scale-invariance for a given vector. Different forms of normalization can result in scale-invariance on different granularities, for example whole-layer for Layer Normalization and per-neuron for Batch Normalization when performed immediately following a given layer. Note that the concatenation of scale-invariant vectors is also scale-invariant. The gradient $\nabla_{\boldsymbol{\omega}}\mathscr{L}(\boldsymbol{\omega}, \ldots)$ for a scale-invariant vector $\boldsymbol{\omega}$ has two important properties:

$$\text{Gradient orthogonality:} \qquad \nabla_{\boldsymbol{\omega}}\mathscr{L}(\boldsymbol{\omega}, \ldots) \perp \boldsymbol{\omega} \tag{1}$$

$$\text{Inverse proportionality:} \qquad \|\nabla_{\boldsymbol{\omega}}\mathscr{L}(\boldsymbol{\omega}, \ldots)\| \propto \|\boldsymbol{\omega}\|^{-1} \tag{2}$$

When the weights are centered directly like in Weight Standardization [27], we also have $\nabla_{\boldsymbol{\omega}} \mathscr{L}(\boldsymbol{\omega}, \ldots) \perp \mathbf{1}$. See Appendix D for further discussion about normalization and a derivation of these properties, which have also been described before [16, 35].

C.2 Defining Measures of the Effective Update Size

We use $\boldsymbol{\omega}$ to denote a weight vector that can achieve rotational equilibrium and refer to such vectors as *rotational weights*. They are typically used in dot products e.g. the weights of fully connected and convolutional layers and are often scale-invariant (but not always, see later sections). We use $\boldsymbol{p}$ for an arbitrary parameter that is not necessarily rotational. When training with weight decay, we break an update $\boldsymbol{p}_t \to \boldsymbol{p}_{t+1}$ (and analogously $\boldsymbol{\omega}_t \to \boldsymbol{\omega}_{t+1}$) for each parameter into two components:

$$\boldsymbol{p}_{t+1} - \boldsymbol{p}_t = \Delta_g \boldsymbol{p}_t + \Delta_\lambda \boldsymbol{p}_t \tag{3}$$

where $\Delta_{\boldsymbol{g}} \boldsymbol{p}_t$ comes from the loss gradient, which we denote with $\boldsymbol{g} := \nabla_{\boldsymbol{p}} \mathscr{L}(\boldsymbol{p}, \ldots)$, and $\Delta_\lambda \boldsymbol{p}_t$ from the weight decay or L_2 regularization. Both of these terms can include averaging over time like in SGD with momentum, causing them to depend on previous parameter and gradient values.

When measuring the size of an update in this work, we focus on the gradient component $\Delta_{\boldsymbol{g}} \boldsymbol{p}_t$, considering the weight decay as a separate process. Note that some scale-sensitive parameters like biases and gains are often excluded from weight decay, see for example the PyTorch Image Models [36] or Hugging Face Transformers [37] libraries. We can measure the update size of parameters using $\|\Delta_g \boldsymbol{p}_t\|$ or on average with the *root-mean-square (RMS) update size*:

$$\eta_g := \sqrt{\mathbb{E}[\|\Delta_g \boldsymbol{p}\|^2]} \tag{4}$$

For rotational weight vectors, we consider the *expected angular update size* defined as:

$$\eta_r := \mathbb{E}[\angle(\boldsymbol{\omega}_t, \boldsymbol{\omega}_{t+1})] = \mathbb{E}\left[\arccos\left(\frac{\langle \boldsymbol{\omega}_t, \boldsymbol{\omega}_{t+1} \rangle}{\|\boldsymbol{\omega}_t\| \|\boldsymbol{\omega}_{t+1}\|}\right)\right] \tag{5}$$

where $\langle \cdot \rangle$ denotes an inner product. Note that for a scale-invariant $\boldsymbol{\omega}$, only the direction $\boldsymbol{\omega}/\|\boldsymbol{\omega}\|$ matters since the magnitude $\|\boldsymbol{\omega}\|$ does not. As a result, η_r is a more natural way to measure the effect of an update on $\boldsymbol{\omega}$, compared to other metrics such as η_g that are not scale-invariant. We can also often scale one layer by a constant and undo the effects of this by scaling another parameter or layer (see e.g. [23]). In such cases relative metrics like the angular update may still capture the effect of an update better than η_g, even if $\boldsymbol{\omega}$ is not strictly scale-invariant by itself.

In related literature there are multiple definitions of an "effective" learning rate that vary slightly between works [34, 6, 35]. We use η_r as a measure of an *effective update size* that is closely related to these quantities. However, we want to emphasize the difference between measuring the size of a single update compared to the change over a longer time interval. The total weight change over extended periods is affected by both the magnitude of individual steps (i.e. η_g or η_r) as well as the consistency in the update direction over the period which is affected by momentum and not captured by these metrics. The longer term change may be a better measure of an "effective" learning rate, but the step-wise metrics are easier to measure and control. For a given momentum coefficient they may be roughly proportional (e.g. in a random walk).

D Normalization and Scale-Invariance

Setup: We use Batch Normalization [12] as an example of a normalization operation. Let $\boldsymbol{x} = \boldsymbol{Z}\boldsymbol{p}$ for $\boldsymbol{x} \in \mathbb{R}^{B \times 1}$, $\boldsymbol{p} \in \mathbb{R}^{C \times 1}$ and $\boldsymbol{Z} \in \mathbb{R}^{B \times C}$ correspond to a single output feature of a linear layer (i.e. a neuron). We can write the batch normalization of this feature as:

$$\hat{\boldsymbol{x}} = N(\boldsymbol{x}) = \frac{\boldsymbol{x} - \mu}{\sqrt{\sigma^2 + \varepsilon}}, \qquad \mu = \frac{1}{B}\sum_{i=1}^{B} x_i, \qquad \sigma^2 = \frac{1}{B}\sum_{i=1}^{B}(x_i - \mu)^2 \tag{6}$$

where $\boldsymbol{x} = [x_1, \ldots, x_B]^\top \in \mathbb{R}^B$ is a vector and $\varepsilon \geq 0$ is a small hyperparameter added for numerical stability. Backpropagation accurately treats μ and σ as functions of $\boldsymbol{x}$. When ε is sufficiently small to be ignored, the output of the normalization is not affected by a positive scaling of the input:

$$N(r\boldsymbol{x}) = (r\boldsymbol{x} - r\mu)/\sqrt{r^2\sigma^2 + \varepsilon} \approx (\boldsymbol{x} - \mu)/\sqrt{\sigma^2 + \varepsilon} = N(\boldsymbol{x}), \qquad r > 0 \tag{7}$$

If the training loss $\mathscr{L}$ does not depend on $\boldsymbol{x}$ in other ways than through $N(\boldsymbol{x})$, this makes $\boldsymbol{x}$ scale-invariant with respect to the loss, i.e. $\mathscr{L}(r\boldsymbol{\omega}) = \mathscr{L}(\boldsymbol{\omega})$ for $r > 0$. Note that although we sometimes write $\mathscr{L}(\boldsymbol{\omega})$ for brevity the loss generally depends on other weights and inputs as well, $\boldsymbol{\omega}$ is generally only a portion of the parameters used in the network, and could for example be a particular row in the weight matrix of a fully connected layer. Some normalization operations like Centered Weight Normlization [10] a.k.a. Weight Standardization [27] are performed directly on the weights instead of activations. This also makes the weight scale-invariant and in case of the aforementioned methods also makes $\nabla_{\boldsymbol{\omega}}\mathscr{L}(\boldsymbol{\omega}) \perp \mathbf{1}$.

Properties: Scale-invariance results in the properties stated in Equations (1) and (2), repeated below:

$$\text{Gradient orthogonality:} \quad \nabla_{\boldsymbol{\omega}}\mathscr{L}(\boldsymbol{\omega}) \perp \boldsymbol{\omega} \tag{8}$$

$$\text{Inverse proportionality:} \quad \|\nabla_{\boldsymbol{\omega}}\mathscr{L}(\boldsymbol{\omega})\| \propto \|\boldsymbol{\omega}\|^{-1} \tag{9}$$

Intuition: The first property is a result of the loss surface being invariant along the direction of $\boldsymbol{\omega}$. Hence the directional derivative of $\mathscr{L}(\boldsymbol{\omega})$ in the direction of $\boldsymbol{\omega}$ is zero:

$$\langle \nabla_{\boldsymbol{\omega}}\mathscr{L}(\boldsymbol{\omega}), \frac{\boldsymbol{\omega}}{\|\boldsymbol{\omega}\|} \rangle = \lim_{h \to 0} \frac{\mathscr{L}(\boldsymbol{\omega} + h\boldsymbol{\omega}/\|\boldsymbol{\omega}\|) - \mathscr{L}(\boldsymbol{\omega})}{h} \tag{10}$$

$$= \lim_{h \to 0} \frac{\mathscr{L}(\boldsymbol{\omega}) - \mathscr{L}(\boldsymbol{\omega})}{h} \tag{11}$$

$$= \lim_{h \to 0} \frac{0}{h} = 0 \tag{12}$$

The second property is a result of the backpropagation through N, which scales the gradient by the factor used on the forward pass $1/\sqrt{\sigma^2 + \varepsilon} \approx \sigma^{-1}$ as if it were a constant, and the fact that $\sigma \propto \|\boldsymbol{\omega}\|$.

Backpropagation: The properties can also be shown using expressions for the backpropagation through the normalization layers. For completeness we include the learnable affine transformation that typically follows normalization operations:

$$\boldsymbol{y} = \gamma\hat{\boldsymbol{x}} + \beta \tag{13}$$

For the backpropagation we have:

$$\nabla_{\gamma}\mathscr{L}(\boldsymbol{p}) = \langle \hat{\boldsymbol{x}}, \nabla_{\boldsymbol{y}}\mathscr{L}(\boldsymbol{p}) \rangle \tag{14}$$

$$\nabla_{\beta}\mathscr{L}(\boldsymbol{p}) = \langle \mathbf{1}_B, \nabla_{\boldsymbol{y}}\mathscr{L}(\boldsymbol{p}) \rangle \tag{15}$$

$$\nabla_{\boldsymbol{x}}\mathscr{L}(\boldsymbol{p}) = \tfrac{\gamma}{\sqrt{\sigma^2+\varepsilon}} \cdot \left[\nabla_{\boldsymbol{y}}\mathscr{L}(\boldsymbol{p}) - \tfrac{1}{B}\langle \mathbf{1}_B, \nabla_{\boldsymbol{y}}\mathscr{L}(\boldsymbol{p}) \rangle \mathbf{1}_B - \tfrac{1}{B}\langle \hat{\boldsymbol{x}}, \nabla_{\boldsymbol{y}}\mathscr{L}(\boldsymbol{p}) \rangle \hat{\boldsymbol{x}} \right] \tag{16}$$

Assuming that ε is small gives:

$$\nabla_{\boldsymbol{x}}\mathscr{L}(\boldsymbol{p}) = \tfrac{\gamma}{\sigma} \left[\nabla_{\boldsymbol{y}}\mathscr{L}(\boldsymbol{p}) - \tfrac{1}{B}\langle \mathbf{1}_B, \nabla_{\boldsymbol{y}}\mathscr{L}(\boldsymbol{p}) \rangle \mathbf{1}_B - \tfrac{1}{B}\langle \hat{\boldsymbol{x}}, \nabla_{\boldsymbol{y}}\mathscr{L}(\boldsymbol{p}) \rangle \hat{\boldsymbol{x}} \right] \tag{17}$$

In this case we have:

$$\langle \nabla_{\boldsymbol{x}}\mathscr{L}(\boldsymbol{p}), \mathbf{1}_B \rangle = \tfrac{\gamma}{\sigma} \left[\langle \mathbf{1}_B, \nabla_{\boldsymbol{y}}\mathscr{L}(\boldsymbol{p}) \rangle - \tfrac{1}{B}\langle \mathbf{1}_B, \nabla_{\boldsymbol{y}}\mathscr{L}(\boldsymbol{p}) \rangle \underbrace{\langle \mathbf{1}_B, \mathbf{1}_B \rangle}_{=B} - \tfrac{1}{B}\langle \hat{\boldsymbol{x}}, \nabla_{\boldsymbol{y}}\mathscr{L}(\boldsymbol{p}) \rangle \underbrace{\langle \hat{\boldsymbol{x}}, \mathbf{1}_B \rangle}_{=0} \right]$$

$$= 0 \tag{18}$$

and similarly:

$$\langle \nabla_{\boldsymbol{x}}\mathscr{L}(\boldsymbol{p}), \hat{\boldsymbol{x}} \rangle = \tfrac{\gamma}{\sigma} \left[\langle \hat{\boldsymbol{x}}, \nabla_{\boldsymbol{y}}\mathscr{L}(\boldsymbol{p}) \rangle - \tfrac{1}{B}\langle \mathbf{1}_B, \nabla_{\boldsymbol{y}}\mathscr{L}(\boldsymbol{p}) \rangle \underbrace{\langle \mathbf{1}_B, \hat{\boldsymbol{x}} \rangle}_{=0} - \tfrac{1}{B}\langle \hat{\boldsymbol{x}}, \nabla_{\boldsymbol{y}}\mathscr{L}(\boldsymbol{p}) \rangle \underbrace{\langle \hat{\boldsymbol{x}}, \hat{\boldsymbol{x}} \rangle}_{=B} \right]$$

$$= 0 \tag{19}$$

which gives:

$$\langle \nabla_{\boldsymbol{x}}\mathscr{L}(\boldsymbol{p}), \boldsymbol{x} \rangle = \langle \nabla_{\boldsymbol{x}}\mathscr{L}(\boldsymbol{p}), \sigma\hat{\boldsymbol{x}} + \mu\mathbf{1}_B \rangle = 0 \tag{20}$$

This allows us to obtain the properties of the weight gradient:

$$\nabla_{\boldsymbol{p}} \mathscr{L}(\boldsymbol{p}) = \boldsymbol{Z}^\top \nabla_{\boldsymbol{x}} \mathscr{L}(\boldsymbol{p}) \tag{21}$$

First we note that:

$$\|\nabla_{\boldsymbol{p}} \mathscr{L}(\boldsymbol{p})\| \propto \|\nabla_{\boldsymbol{x}} \mathscr{L}(\boldsymbol{p})\| \propto \sigma^{-1} \propto \|\boldsymbol{p}\|^{-1} \tag{22}$$

where the second proportionality follows from (17) and the final one from (6). This gives the inverse proportionality listed in Equation (9).

We can also derive the gradient orthogonality in Equation (8) as follows:

$$\begin{aligned}
\langle \nabla_{\boldsymbol{p}} \mathscr{L}(\boldsymbol{p}), \boldsymbol{p} \rangle &= \langle \boldsymbol{Z}^\top \nabla_{\boldsymbol{x}} \mathscr{L}(\boldsymbol{p}), \boldsymbol{p} \rangle && (23)\\
&= \boldsymbol{p}^\top \boldsymbol{Z}^\top \nabla_{\boldsymbol{x}} \mathscr{L}(\boldsymbol{p}) && (24)\\
&= \boldsymbol{x}^\top \nabla_{\boldsymbol{x}} \mathscr{L}(\boldsymbol{p}) && (25)\\
&= \langle \nabla_{\boldsymbol{x}} \mathscr{L}(\boldsymbol{p}), \boldsymbol{x} \rangle && (26)\\
&= 0 && (27)
\end{aligned}$$

These properties can also be shown directly from the scale-invariance using calculus theorems as done in Wan et al. [35].

E Geometric Model For Equilibrium

The main goal of our analysis is to obtain simple expressions for the equilibrium magnitude $\widehat{\|\boldsymbol{\omega}\|}$ and the expected angular update in equilibrium, $\widehat{\eta_r}$ (See Appx C for background and definitions). To do this we analyze a simplified setting, the optimization of a random loss function, resulting in a *random walk* for the neural network parameters. In practice, this could be implemented by randomly sampling the inputs and the gradients for the network outputs (e.g. logits) at each step independently from a zero-mean normal distribution. Backpropagating the output gradients results in parameter gradients that are zero-mean and independent between steps, with a distribution that does not change over time due to the optimization progress on an underlying objective function. Appendix K gives further information and explores differences between a random walk and real neural network optimization and how they affect the predictions. In our experiments we find that the final predictions hold well for a variety of networks despite being derived for this simplified setting.

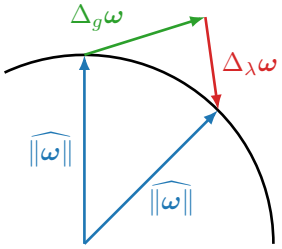

Figure 6: Weight norm equilibrium where a single expected optimizer step is split into components $\Delta_g \boldsymbol{\omega}$ due to the loss gradient and $\Delta_\lambda \boldsymbol{\omega}$ from weight decay.

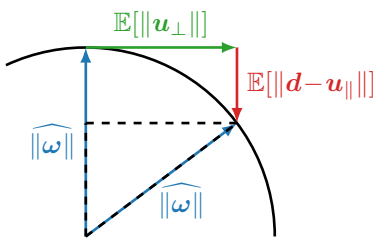

Figure 7: Weight norm equilibrium. The loss gradient causes an update $\boldsymbol{u}$ and the weight decay $\boldsymbol{d}$.

In this section we present a simple geometric derivation of the equilibrium norm $\widehat{\|\boldsymbol{\omega}\|}$ for different optimizers inspired in part by the analysis in Online Normalization [6]. Equilibrium is an abstract state where the effects of the gradient component of the update $\Delta_g \boldsymbol{\omega}$ and the weight decay component $\Delta_\lambda \boldsymbol{\omega}$ on the expected weight magnitude balance out on average. These components typically have different monotonic dependencies on the weight magnitude, with weight decay being proportional while the gradient component is either constant or inversely proportional, depending on the setting. As a result, the effects of these components can balance out in expectation for a particular magnitude, which we call the equilibrium norm $\widehat{\|\boldsymbol{\omega}\|}$. As shown in Figure 6, the geometry of this is not necessarily simple. Due to the averaging effects of momentum over time, $\Delta_g \boldsymbol{\omega}$ is not necessarily orthogonal to the weights even in cases where individual gradients are (e.g. for scale-invariant weights). In the same way, the weight decay (or L_2-regularization) component $\Delta_\lambda \boldsymbol{\omega}$ may not be perfectly anti-parallel to the weights with momentum.

To simplify the effects of momentum, we instead consider a different view of equilibrium shown in Figure 7. Here we consider the total weight change throughout training derived from the weight and gradient at a given time step, instead of the update that is applied in that iteration. We thus define $\boldsymbol{u}$ for time step t as the sum of the contributions of $\nabla_{\boldsymbol{\omega}}\mathscr{L}(\boldsymbol{\omega}_t, \ldots)$ to subsequent updates $\boldsymbol{\omega}_t \to \boldsymbol{\omega}_{t+1}$, $\boldsymbol{\omega}_{t+1} \to \boldsymbol{\omega}_{t+2}$, and so on. Analogously, the weight decay term $\boldsymbol{d}$ is defined as the total weight change due to the weight decay or L_2-regularization of the weights $\boldsymbol{\omega}_t$ at iteration t. Note that without momentum $\boldsymbol{u} = \Delta_g\boldsymbol{\omega}$, $\boldsymbol{d} = \Delta_\lambda\boldsymbol{\omega}$ and that if $\Delta_g\boldsymbol{\omega}$ and $\Delta_\lambda\boldsymbol{\omega}$ balance out on average, then so must $\boldsymbol{u}$ and $\boldsymbol{d}$.

In many cases $\boldsymbol{u}$ is orthogonal to the weights on average due to scale-invariance or randomness, but otherwise can we split it into orthogonal $\boldsymbol{u}_\perp$ and radial $\boldsymbol{u}_\parallel$ components. The weight decay term $\boldsymbol{d}$ is anti-parallel to the weights in every case we consider. If we can obtain an expression for $\|\boldsymbol{u}_\perp\|$ and $\|\boldsymbol{d} - \boldsymbol{u}_\parallel\|$, this allows us to apply the Pythagorean theorem to the dashed triangle in Figure 7:

$$(\widehat{\|\boldsymbol{\omega}\|} - \mathbb{E}[\|\boldsymbol{d} - \boldsymbol{u}_\parallel\|])^2 + \mathbb{E}[\|\boldsymbol{u}_\perp\|]^2 = \widehat{\|\boldsymbol{\omega}\|}^2 \tag{28}$$

We can then solve for $\widehat{\|\boldsymbol{\omega}\|}$ after accounting for the dependency of $\boldsymbol{u}$ and $\boldsymbol{d}$ on the weight norm.

Once we have an expression for $\widehat{\|\boldsymbol{\omega}\|}$, we can compute a prediction for the RMS update size $\widehat{\eta_g}$. Combining $\widehat{\eta_g}$ for the equilibrium magnitude $\widehat{\|\boldsymbol{\omega}\|}$ allows us to compute the expected relative update size $\widehat{\eta_g}/\widehat{\|\boldsymbol{\omega}\|}$ which closely approximates η_r in equilibrium. We do this for AdamW, SGDM and Lion [5] in Appendix F, G and H respectively. The results for each optimizer are summarized in Table 1.

F AdamW Equilibrium

The standard version of AdamW [21] can be written as:

$$\boldsymbol{m}_t = \beta_1\boldsymbol{m}_{t-1} + (1-\beta_1)\boldsymbol{g}_t \tag{29}$$

$$\boldsymbol{v}_t = \beta_2\boldsymbol{v}_{t-1} + (1-\beta_2)\boldsymbol{g}_t^2 \tag{30}$$

$$\boldsymbol{p}_t = \boldsymbol{p}_{t-1} - \eta \cdot \left(\frac{\boldsymbol{m}_t/(1-\beta_1^t)}{\sqrt{\boldsymbol{v}_t/(1-\beta_2^t)}+\varepsilon} + \lambda\boldsymbol{p}_{t-1}\right) \tag{31}$$

Where $\boldsymbol{p}_t \in \mathbb{R}^C$ is a parameter vector at time t, $\boldsymbol{g}_t = \nabla_{\boldsymbol{p}}\mathscr{L}(\boldsymbol{p}_t, \ldots)$ is the gradient, $\boldsymbol{m}$ is the first moment and $\boldsymbol{v}$ is the second moment. The learning rate ($\eta \geq 0$), weight decay ($\lambda \geq 0$), moment coefficients ($0 < \beta_1 < 1, 0 < \beta_2 < 1$) and $\varepsilon \geq 0$ are hyperparameters. For simplicity we assume that ε and the bias correction can be ignored, i.e. that ε, β_1^t and β_2^t are all effectively zero.

Equilibrium magnitude: For a rotational weight $\boldsymbol{\omega}$, we can write $\boldsymbol{u}$ and $\boldsymbol{d}$ from Section E as:

$$\boldsymbol{u} = -\eta\textstyle\sum_{k=t}^{\infty}\beta_1^{k-t}(1-\beta_1)\frac{\boldsymbol{g}_t}{\sqrt{\boldsymbol{v}_k}}, \qquad \boldsymbol{d} = -\eta\lambda\boldsymbol{\omega} \tag{32}$$

We note that due to symmetry, each coordinate of $\boldsymbol{u}$ has a zero-mean symmetric distribution in the random walk setup. Since $\boldsymbol{u}$ is independent from $\boldsymbol{\omega}$, this makes them orthogonal in expectation i.e. $\mathbb{E}[\langle\boldsymbol{u}, \boldsymbol{\omega}\rangle] = 0$. When the gradient distribution is not changing over time, it is also reasonable to assume that the variance of each coordinate remains constant resulting in $\forall t, k : \mathbb{E}[\|\boldsymbol{g}_t/\sqrt{\boldsymbol{v}_k}\|^2] = C$ (for $\boldsymbol{\omega} \in \mathbb{R}^C$) and by extension $\mathbb{E}[\|\boldsymbol{u}\|^2] = \eta^2 C$. Defining $\omega = \|\boldsymbol{\omega}\|$, $u = \|\boldsymbol{u}\|$, $u_\parallel = \langle\boldsymbol{\omega}, \boldsymbol{u}\rangle/\|\boldsymbol{\omega}\|$, $u_\perp^2 = u^2 - u_\parallel^2$ and $d = \|\boldsymbol{d}\|$ we can write a recurrence relation based on Equation (28):

$$\mathbb{E}[\omega_{i+1}^2] = \mathbb{E}[(\omega_i - d + u_\parallel)^2 + u_\perp^2] \tag{33}$$

$$= \mathbb{E}[\omega_i^2 - 2d\omega_i + 2u_\parallel\omega_i - 2du_\parallel + d^2 + u_\parallel^2 + (u^2 - u_\parallel^2)] \tag{34}$$

$$= \mathbb{E}[\omega_i^2](1 - 2\eta\lambda + \eta^2\lambda^2) + \eta^2 C \tag{35}$$

where we have used independence, $\mathbb{E}[u_\parallel] = 0$ and $\mathbb{E}[u] = \eta^2 C$. The solution is:

$$\mathbb{E}[\omega_i^2] = \mathbb{E}[\omega_0^2]a^i + \tfrac{\eta^2 C}{2\eta\lambda - \eta^2\lambda^2}(1 - a^i), \qquad a = (1 - 2\eta\lambda + \eta^2\lambda^2) \tag{36}$$

The recurrence relation is written in terms of $\boldsymbol{u}$ and $\boldsymbol{d}$ instead of $\Delta_g\boldsymbol{\omega}$ and $\Delta_\lambda\boldsymbol{\omega}$. This is thus only an approximation of how the real system converges to equilibrium over time, but still informative. It may be a good approximation if $\|\boldsymbol{\omega}\|$ changes slowly compared to how quickly $\boldsymbol{u}$ is applied (i.e. $\boldsymbol{m}$ changes) and $\boldsymbol{v}$ is updated. In either case, the limit gives us the equilibrium norm listed in Table 1:

$$\widehat{\|\boldsymbol{\omega}\|} = \sqrt{\frac{\eta C}{2\lambda - \eta\lambda^2}} \approx \sqrt{\frac{\eta C}{2\lambda}} \qquad \text{(for } \lambda\eta \ll 2\text{)} \tag{37}$$

Update size: We can estimate the RMS update size η_g of $\Delta_g \boldsymbol{p} = \frac{\boldsymbol{m}_t}{\sqrt{\boldsymbol{v}_t}}$ as follows:

$$\mathbb{E}[\|\Delta_g \boldsymbol{p}\|^2] = \mathbb{E}[\|\tfrac{1}{\sqrt{\boldsymbol{v}_t}}(1-\beta_1)\textstyle\sum_{k=0}^{t-1}\beta_1^{t-k}\boldsymbol{g}_{t-k}\|^2] \tag{38}$$

$$= (1-\beta_1)^2 \textstyle\sum_{k=0}^{t-1}\beta_1^{2t-2k}\mathbb{E}[\|\frac{\boldsymbol{g}_{t-k}}{\sqrt{\boldsymbol{v}_t}}\|^2] \tag{39}$$

$$\approx \eta^2 \tfrac{1-\beta_1}{1+\beta_1} C \tag{40}$$

where we have approximated the geometric sum with its limit $t \to \infty$, used the fact that for the random walk $\forall j \neq k : \mathbb{E}[\langle \boldsymbol{g}_j, \boldsymbol{g}_k \rangle] = 0$ as well as our previous assumption $\forall t, k : \mathbb{E}[\|\boldsymbol{g}_t / \sqrt{\boldsymbol{v}_k}\|^2] = C$. This gives us the prediction $\widehat{\eta_g} \approx \mathbb{E}[\sqrt{\|\Delta_g \boldsymbol{p}\|^2}]$ listed in Table 1. Approximating the equilibrium angular update size with the expected relative update size $\sqrt{\mathbb{E}[\|\Delta_g \boldsymbol{\omega}\|^2]}/\widehat{\|\boldsymbol{\omega}\|}$ gives the $\widehat{\eta_r}$ value. This approximation is good for small relative updates and a relatively small radial component in $\Delta_g \boldsymbol{\omega}$.

G SGDM Equilibrium

The standard version of SGD with Momentum can be written as:

$$\boldsymbol{m}_t = \alpha \boldsymbol{m}_{t-1} + \boldsymbol{g}_t + \lambda \boldsymbol{p}_{t-1} \tag{41}$$

$$\boldsymbol{p}_t = \boldsymbol{p}_{t-1} - \eta \cdot \boldsymbol{m}_t \tag{42}$$

Where $\boldsymbol{p}_t \in \mathbb{R}^C$ is a parameter vector at time t, $\boldsymbol{g}_t = \nabla_{\boldsymbol{p}} \mathscr{L}(\boldsymbol{p}_t)$ is the gradient, $\boldsymbol{m}$ is the first moment and $\boldsymbol{v}$ is the second moment. The learning rate ($\eta \geq 0$), weight decay ($\lambda \geq 0$), momentum coefficient ($0 < \alpha < 1$) are hyperparameters.

We compute the total weight change due to $\boldsymbol{g}_t$, i.e. $\boldsymbol{u}$ in Equation (28) as:

$$\boldsymbol{u} = -\eta \textstyle\sum_{k=t}^{\infty} \alpha^{k-t} \boldsymbol{g}_t = \frac{-\eta}{1-\alpha} \boldsymbol{g}_t \tag{43}$$

Analogously, the total weight change due to $\boldsymbol{w}_t$, i.e. $\boldsymbol{d}$ in Equation (28) is:

$$\boldsymbol{d} = -\eta \textstyle\sum_{k=t}^{\infty} \alpha^{k-t} \lambda \boldsymbol{p}_t = \frac{-\eta\lambda}{1-\alpha} \boldsymbol{p}_t \tag{44}$$

Combining (43) and (44), this allows us to solve (28) for a scale-invariant weight vector $\boldsymbol{\omega}$. Here we assume scale-invariance since it slightly changes the resulting expression due to the dependency of $\|\boldsymbol{u}\|$ on $\|\boldsymbol{\omega}\|$. It also simplifies the math a bit, with $\boldsymbol{u} \perp \boldsymbol{\omega}$, not just in expectation. We get:

$$\widehat{\|\boldsymbol{\omega}\|}^2 = \mathbb{E}\left[(\widehat{\|\boldsymbol{\omega}\|} - \|\boldsymbol{d}\|)^2 + \|\boldsymbol{u}\|^2\right] \tag{45}$$

$$= \mathbb{E}\left[\left(\widehat{\|\boldsymbol{\omega}\|} - \tfrac{\eta\lambda}{1-\alpha}\widehat{\|\boldsymbol{\omega}\|}\right)^2 + \left(\tfrac{\eta}{1-\alpha}\tfrac{\|\tilde{\boldsymbol{g}}\|}{\widehat{\|\boldsymbol{\omega}\|}}\right)^2\right] \tag{46}$$

Where we define $\tilde{\boldsymbol{g}}_t = \boldsymbol{g}_t \|\boldsymbol{\omega}_t\|$ using $\|\boldsymbol{g}_t\| \propto \|\boldsymbol{\omega}_t\|^{-1}$ due to the inverse proportionality of the gradient magnitude, see Equation (2) or (9). We can interpret $\tilde{\boldsymbol{g}}_t$ as the gradient for weights of unit norm $\|\boldsymbol{\omega}_t\| = 1$.

Solving for $\widehat{\|\boldsymbol{\omega}\|}$ and assuming that $\eta\lambda \ll 2 \cdot (1-\alpha)$ gives:

$$\widehat{\|\boldsymbol{\omega}\|} = \sqrt[4]{\frac{\eta\mathbb{E}[\|\tilde{\boldsymbol{g}}\|^2]}{2\lambda\cdot(1-\alpha) - \eta\lambda^2}} \approx \sqrt[4]{\frac{\eta\mathbb{E}[\|\tilde{\boldsymbol{g}}\|^2]}{2\lambda\cdot(1-\alpha)}} \tag{47}$$

To obtain the absolute size of an update, we further assume that $\mathbb{E}[\|\boldsymbol{g}_t\|^2]$ can be approximated as a constant $\mathbb{E}[\|\boldsymbol{g}\|^2]$ when computing the size of $\boldsymbol{m}_t$, and that successive gradients are roughly orthogonal giving $\boldsymbol{m}_{t-1} \perp \boldsymbol{g}_t$ in expectation. For the random walk setting, the first is reasonable when the norm is stable e.g. around equilibrium and the second always holds. The average square size of an update is then:

$$\mathbb{E}[\|\Delta_g \boldsymbol{p}\|^2] = \eta^2 \mathbb{E}\left[\|\alpha \boldsymbol{m}_{t-1} + \boldsymbol{g}_t\|^2\right] \tag{48}$$

$$= \eta^2 \mathbb{E}\left[\|\alpha \boldsymbol{m}_{t-1}\|^2\right] + \mathbb{E}\left[\|\boldsymbol{g}_t\|^2\right] \tag{49}$$

$$= \eta^2 \textstyle\sum_{k=0}^{t} \mathbb{E}[\left(\alpha^{t-k}\|\boldsymbol{g}_k\|\right)^2] \tag{50}$$

$$\approx \eta^2 \tfrac{\mathbb{E}[\|\boldsymbol{g}\|^2]}{1-\alpha^2} \tag{51}$$

where (49) comes from the orthogonality, (50) by recursively writing out $\boldsymbol{m}$ in terms of $\boldsymbol{g}$, and (51) from assuming that t is high enough to approximate the sum of the geometric series as an infinite sum.

Simplifying, we get the $\eta_g = \sqrt{\mathbb{E}[\|\Delta_g \boldsymbol{p}\|^2]}$ and $\eta_r = \sqrt{\mathbb{E}[\|\Delta_g \boldsymbol{\omega}\|^2]}/\widehat{\|\boldsymbol{\omega}\|}$ in Table 1. We note that the derived rates are consistent with the ones derived in the Spherical Motion Dynamics [35] and Online Normalization [6] (when $\alpha = 0$).

H Lion Equilibrium

The standard version of Lion [5] can be written as:

$$\boldsymbol{v}_t = \operatorname{sign}(\beta_1 \boldsymbol{m}_{t-1} + (1-\beta_1)\boldsymbol{g}_t) \tag{52}$$

$$\boldsymbol{p}_t = \boldsymbol{p}_{t-1} - \eta \cdot \Big(\boldsymbol{v}_t + \lambda \boldsymbol{p}_{t-1}\Big) \tag{53}$$

$$\boldsymbol{m}_t = \beta_2 \boldsymbol{m}_{t-1} + (1-\beta_2)\boldsymbol{g}_t \tag{54}$$

Where $\boldsymbol{p}_t \in \mathbb{R}^C$ is a parameter vector at time t, $\boldsymbol{g}_t = \nabla_{\boldsymbol{p}} \mathscr{L}(\boldsymbol{p}_t)$ is the gradient, $\boldsymbol{m}$ is the first moment and $\boldsymbol{v}$ is the update velocity. The learning rate ($\eta \geq 0$), weight decay ($\lambda \geq 0$), moment coefficients ($0 < \beta_1 < 1, 0 < \beta_2 < 1$).

In our analysis we look at the arguments of the sign function which we define as:

$$\boldsymbol{n}_t := \beta_1 \boldsymbol{m}_{t-1} + (1-\beta_1)\boldsymbol{g}_t, \qquad \boldsymbol{v}_t = \operatorname{sign}(\boldsymbol{n}_t) \tag{55}$$

To obtain an estimate of the magnitude $\|\boldsymbol{n}_t\|$, we assume that the gradient magnitude $\mathbb{E}[\|\boldsymbol{g}_t\|^2]$ can be approximated as a constant $\mathbb{E}[\|\boldsymbol{g}\|^2]$ and that successive gradients are roughly orthogonal giving $\boldsymbol{m}_{t-1} \perp \boldsymbol{g}_t$ in expectation. For the random walk setting, the first is reasonable when the norm is stable e.g. around equilibrium and the second always holds. This gives:

$$\mathbb{E}[\|\boldsymbol{n}_t\|^2] = \mathbb{E}\left[\|\beta_1 \boldsymbol{m}_{t-1} + (1-\beta_1)\boldsymbol{g}_t\|^2\right] \tag{56}$$

$$= \beta_1^2 \mathbb{E}\left[\|\beta_2 \boldsymbol{m}_{t-1} + (1-\beta_2)\boldsymbol{g}_{t-1}\|^2\right] + (1-\beta_1)^2 \mathbb{E}\left[\|\boldsymbol{g}_t\|^2\right] \tag{57}$$

$$= \beta_1^2 (1-\beta_2)^2 \sum_{k=0}^{k=t-1} \beta_2^{t-1-k} \mathbb{E}[\|\boldsymbol{g}_k\|^2] + (1-\beta_1)^2 \mathbb{E}\left[\|\boldsymbol{g}_t\|^2\right] \tag{58}$$

$$\approx \beta_1^2 (1-\beta_2)^2 \sum_{k=0}^{\infty} \beta_2^k \mathbb{E}[\|\boldsymbol{g}\|^2] + (1-\beta_1)^2 \mathbb{E}\left[\|\boldsymbol{g}\|^2\right] \tag{59}$$

$$= \left((1-\beta_1)^2 + \beta_1^2 \frac{1-\beta_2}{1+\beta_2}\right) \mathbb{E}[\|\boldsymbol{g}\|^2] \tag{60}$$

where we have used the gradient orthogonality and constant magnitude and approximated the geometric sum as extending to infinity.

To compute the gradient contribution $\|\boldsymbol{u}\|$ in Equation (28), we first need to model how the sign non-linearity affects the magnitude and direction of the update. We note that for $\boldsymbol{p}_t \in \mathbb{R}^C$:

$$\|\boldsymbol{v}_t\| = \sqrt{C} \tag{61}$$

so the sign function has an average scaling effect:

$$\frac{\|\boldsymbol{v}_t\|}{\|\boldsymbol{n}_t\|} = \sqrt{\frac{C}{\left((1-\beta_1)^2 + \beta_1^2 \frac{1-\beta_2}{1+\beta_2}\right) \mathbb{E}[\|\boldsymbol{g}\|^2]}} \tag{62}$$

The sign function will also rotate $\boldsymbol{n}_t$ resulting in two components, one parallel to $\boldsymbol{n}_t$ and the other orthogonal. We will assume that the orthogonal one cancels out on average without significantly affecting equilibrium and focus on the parallel component. This component depends on the average angle between $\boldsymbol{n}_t$ and $\operatorname{sign}(\boldsymbol{n}_t)$ which is determined by the distribution and correlation between the elements. In the random walk setting, we can assume the components of $\boldsymbol{n}_t = [n_1, \ldots, n_C]$ are normally distributed with mean zero. However, the expression for the average angle is still complicated unless the components are independent and identically distributed (i.i.d.) so we make

this assumption for this step with $n_k \sim \mathcal{N}(0, \sigma^2)$ i.i.d. for all k. Then we can use the known expected absolute value for a centered normal distribution to get:

$$\mathbb{E}[\langle \boldsymbol{n}_t, \operatorname{sign}(\boldsymbol{n}_t)\rangle] = C \cdot \mathbb{E}[|n_k|] = C \cdot \sqrt{\frac{2\sigma^2}{\pi}} = \sqrt{\frac{2}{\pi}} \cdot \|\boldsymbol{n}_t\| \cdot \|\operatorname{sign}(\boldsymbol{n}_t)\| \tag{63}$$

Note that the angle is still bounded regardless of the distribution but will result in a different factor in the range that $\|\boldsymbol{n}\|_1/(\sqrt{C}\|\boldsymbol{n}\|_2)$ can take, i.e. $[C^{-\frac{1}{2}}, 1]$ instead of $\sqrt{2/\pi}$.

Based on the preceding analysis we will model the sign function for the computation of $\|\boldsymbol{u}\|$ as:

$$\operatorname{sign}(\boldsymbol{n}_t) \approx \sqrt{\frac{2}{\pi}} \frac{\|\boldsymbol{v}_t\|}{\|\boldsymbol{n}_t\|} \boldsymbol{n}_t = \sqrt{\frac{2C}{\mathbb{E}[\|\boldsymbol{g}\|^2] \cdot \pi \cdot \left((1-\beta_1)^2 + \beta_1^2 \frac{1-\beta_2}{1+\beta_2}\right)}} \boldsymbol{n}_t \tag{64}$$

which gives:

$$\mathbb{E}[\|\boldsymbol{u}\|^2] = \frac{2\eta C}{\mathbb{E}[\|\boldsymbol{g}\|^2] \cdot \pi \cdot \left((1-\beta_1)^2 + \beta_1^2 \frac{1-\beta_2}{1+\beta_2}\right)} \mathbb{E}\left[\left\|(1-\beta_1)\boldsymbol{g} + \beta_1(1-\beta_2)\sum_{k=0}^{\infty} \beta_2^k \boldsymbol{g}\right\|^2\right] \tag{65}$$

$$= \frac{2\eta C}{\pi \cdot \left((1-\beta_1)^2 + \beta_1^2 \frac{1-\beta_2}{1+\beta_2}\right)} \tag{66}$$

Combined with $\boldsymbol{d} = -\eta\lambda\boldsymbol{\omega}_{t-1}$, this allows us to solve (28) for a scale-invariant weight vector $\boldsymbol{\omega}$:

$$\widehat{\|\boldsymbol{\omega}\|}^2 = \mathbb{E}\left[(\widehat{\|\boldsymbol{\omega}\|} - \|\boldsymbol{d}\|)^2 + \|\boldsymbol{u}\|^2\right] \tag{67}$$

$$= (\widehat{\|\boldsymbol{\omega}\|} - \eta\lambda\widehat{\|\boldsymbol{\omega}\|})^2 + \mathbb{E}[\|\boldsymbol{u}\|^2] \tag{68}$$

$$= (1 - 2\eta\lambda + \eta^2\lambda^2)\widehat{\|\boldsymbol{\omega}\|}^2 + \frac{2\eta C}{\pi}\left((1-\beta_1)^2 + \beta_1^2\frac{1-\beta_2}{1+\beta_2}\right)^{-1} \tag{69}$$

Solving for $\widehat{\|\boldsymbol{\omega}\|}$ and assuming $\eta\lambda \ll 1$ gives:

$$\widehat{\|\boldsymbol{\omega}\|} = \sqrt{\tfrac{2\eta C}{\pi\cdot(2\lambda - \eta\lambda^2)}}\left((1-\beta_1)^2 + \beta_1^2\tfrac{1-\beta_2}{1+\beta_2}\right)^{-\frac{1}{2}} \approx \sqrt{\tfrac{\eta C}{\pi\lambda}}\left((1-\beta_1)^2 + \beta_1^2\tfrac{1-\beta_2}{1+\beta_2}\right)^{-\frac{1}{2}} \tag{70}$$

Combined with $\|\boldsymbol{v}_t\| = \sqrt{C}$ for $\boldsymbol{\omega}, \boldsymbol{p} \in \mathbb{R}^C$ we get:

$$\widehat{\eta_g} = \sqrt{\mathbb{E}[\|\Delta_g \boldsymbol{p}\|^2]} = \eta\sqrt{C} \tag{71}$$

$$\widehat{\eta_r} = \sqrt{\mathbb{E}[\|\Delta_g \boldsymbol{\omega}\|^2]}/\widehat{\|\boldsymbol{\omega}\|} = \sqrt{\pi\eta\lambda} \cdot \left((1-\beta_1)^2 + \beta_1^2\frac{1-\beta_2}{1+\beta_2}\right)^{\frac{1}{2}} \tag{72}$$

I Adam+L_2

I.1 Decoupled Weight Decay vs L_2-Regularization in Adam

Loshchilov and Hutter [21] proposed the use of decoupled weight decay instead of L_2-regularization in Adam [13]. In their experiments they find that Adam with decoupled weight decay (i.e. AdamW, see Equation 31) outperforms the L_2-regularized form (i.e. Adam+L_2, Equation 75) across a wide range of settings. Since then AdamW has been widely adopted, but as far as we know the reason for its effectiveness over Adam+L_2 is not well understood.

Figure 8 compares the dynamics of Adam+L_2 and AdamW for a random walk in a simple system described in Appendix J. We observe that the weight magnitude and angular updates converge to a stable value in expectation for both optimizers. However, these values are identical across different scale-invariant weight vectors with AdamW but not Adam+L_2 In Appendix I.2 we analyze the geometric model for Adam+L_2, revealing a complicated dependency on the gradient magnitude, resulting in imbalanced rotation. Note that for SGDM the equilibrium norm also depends on the gradient norm but not the rotation (Table 1). We believe the balanced vs imbalanced equilibrium rotation is a key difference between Adam+L_2 and AdamW which may explain why decoupled weight decay is more effective for Adam-like methods. We explore this further in our experiments.

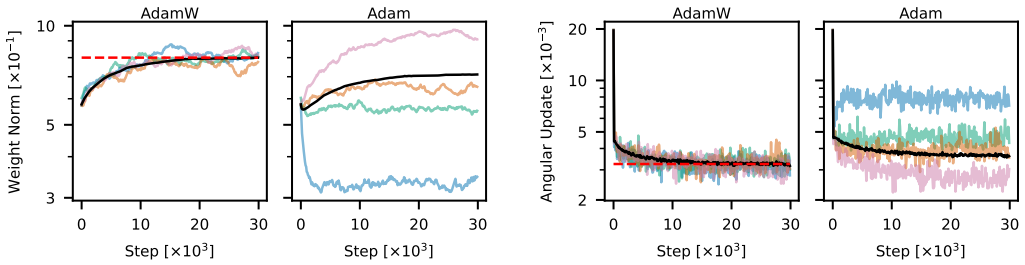

Figure 8: Comparing the equilibrium behavior of AdamW and Adam+L_2 in a simplified random setup for weight norm (left) and angular updates (right). Colors (pink, orange, green, blue) represent scale-invariant weight vectors with various gradient norms. Black denotes layer-wide averages, and the red dashed lines represents predictions from Table 1.

I.2 Equilibrium Analysis for Adam+L_2

In this section we apply the geometric model from Section E to Adam [13] with L_2 regularization to gain some insight into how the rotational equilibrium differs from that of Adam with decoupled weight decay (AdamW, see Section F). We will write the Adam update as follows:

$$\boldsymbol{m}_t = \beta_1 \boldsymbol{m}_{t-1} + (1-\beta_1)(\boldsymbol{g}_t + \lambda \boldsymbol{p}_{t-1}) \tag{73}$$

$$\boldsymbol{v}_t = \beta_2 \boldsymbol{v}_{t-1} + (1-\beta_2)(\boldsymbol{g}_t + \lambda \boldsymbol{p}_{t-1})^2 \tag{74}$$

$$\boldsymbol{p}_t = \boldsymbol{p}_{t-1} - \eta \cdot \left(\frac{\boldsymbol{m}_t / (1-\beta_1^t)}{\sqrt{\boldsymbol{v}_t / (1-\beta_2^t)} + \varepsilon} \right) \tag{75}$$

Similar to AdamW, $\boldsymbol{p}_t \in \mathbb{R}^C$ is a parameter vector at time t, $\boldsymbol{g}_t = \nabla_{\boldsymbol{p}} \mathscr{L}(\boldsymbol{p}_t)$ is the gradient, and all operations (e.g. division, squaring) are performed elementwise. In Adam $\boldsymbol{m}$ and $\boldsymbol{v}$ are the first and second moment of the gradient including an L_2 regularization term. This differs from AdamW (see Equation 31) where the L_2 regularization (or weight decay) does not affect $\boldsymbol{m}$ and $\boldsymbol{v}$. The learning rate ($\eta \geq 0$), L_2 regularization coefficient ($\lambda \geq 0$), moment coefficients ($0 < \beta_1 < 1, 0 < \beta_2 < 1$) and $\varepsilon \geq 0$ are hyperparameters similar to AdamW.

The rotational dynamics of Adam are more complicated than those of AdamW and we will only explore a further simplified setting using strong approximations. Like for the other optimizers, we will assume the rotational equilibrium is a steady state with constant hyperparameters, a constant expected weight norm $\mathbb{E}[\|\boldsymbol{\omega}_t\|] = \widehat{\|\boldsymbol{\omega}\|}$ and gradient norm $\mathbb{E}[\|\boldsymbol{g}_t\|^2] = \mathbb{E}[\|\boldsymbol{g}\|^2]$ over time. Furthermore, we will assume that the second moment is constant over time and elements, and that it can be expressed simply in terms of the gradient and parameter norms i.e.:

$$\boldsymbol{v}_t = \mathbf{1} \sqrt{\frac{\mathbb{E}[\|\boldsymbol{g}\|^2] + \lambda^2 \mathbb{E}[\|\boldsymbol{p}\|^2]}{C}} =: \mathbf{1} v \tag{76}$$

This last approximation is quite rough and is only likely to hold well in special cases. Finally we will assume that ε and the bias correction can be ignored, i.e. that ε, β_1^t and β_2^t are all effectively zero.

The total weight change due to the gradient $\boldsymbol{g}_t$, i.e. $\boldsymbol{u}$ in Equation (28) is given by:

$$\boldsymbol{u} = -\eta \textstyle\sum_{k=t}^{\infty} \beta_1^{k-t}(1-\beta_1) \frac{\boldsymbol{g}_t}{\boldsymbol{v}} = -\eta \frac{\boldsymbol{g}_t}{v} \tag{77}$$

Similarly the total change due to the weight decay is:

$$\boldsymbol{d} = -\eta \sum_{k=t}^{\infty} \beta_1^{k-t}(1-\beta_1) \frac{\lambda \boldsymbol{p}_{t-1}}{\boldsymbol{v}} = -\eta \frac{\lambda \boldsymbol{p}_{t-1}}{v} \tag{78}$$

Here we assume scale-invariance since it slightly changes the resulting expression due to the dependency of the gradient on $\|\boldsymbol{\omega}\|$. For a scale-invariant weight vector $\boldsymbol{p} = \boldsymbol{\omega}$, we can then write

Equation 28 for the equilibrium norm as:

$$\widehat{\|\boldsymbol{\omega}\|}^2 = \mathbb{E}\left[(\widehat{\|\boldsymbol{\omega}\|} - \|\boldsymbol{d}\|)^2 + \|\boldsymbol{u}\|^2\right] \tag{79}$$

$$= (1 - \tfrac{\eta\lambda}{v})^2(\widehat{\|\boldsymbol{\omega}\|}^2 + \eta^2 \tfrac{\mathbb{E}[\|\boldsymbol{g}\|^2]}{v^2} \tag{80}$$

$$= (1 - \tfrac{2\eta\lambda}{v} + \tfrac{\eta^2\lambda^2}{v^2})\widehat{\|\boldsymbol{\omega}\|}^2 + \eta^2\left(C - \tfrac{\lambda^2}{v^2}\widehat{\|\boldsymbol{\omega}\|}^2\right) \tag{81}$$

$$= (1 - \tfrac{2\eta\lambda}{v})\widehat{\|\boldsymbol{\omega}\|}^2 + \eta^2 C\widehat{\|\boldsymbol{\omega}\|}^2 \tag{82}$$

$$= \left(1 - \sqrt{\tfrac{4\eta^2\lambda^2 C}{\mathbb{E}[\|\tilde{\boldsymbol{g}}\|^2]\widehat{\|\boldsymbol{\omega}\|}^{-2} + \lambda^2\widehat{\|\boldsymbol{\omega}\|}^2}}\right)\widehat{\|\boldsymbol{\omega}\|}^2 + \eta^2 C \tag{83}$$

where we have used Equation 76 twice and defined $\|\tilde{\boldsymbol{g}}\| = \|\boldsymbol{p}\|\|\boldsymbol{g}\|$ as the gradient norm corresponding to a unit weight norm based on Equation 2.

Rearranging and simplifying, this gives us a third order polynomial in $\widehat{\|\boldsymbol{\omega}\|}^2$:

$$4\lambda^2\widehat{\|\boldsymbol{\omega}\|}^6 - C\eta^2\lambda^2\widehat{\|\boldsymbol{\omega}\|}^4 - C\eta^2\mathbb{E}[\|\tilde{\boldsymbol{g}}\|^2] = 0 \tag{84}$$

Although a closed form solution exists, it is unwieldy and we were unable to simplify it much in the general case. For known coefficients, it is easy to find the real root with numerical methods. Note that the equilibrium norm depends on $\mathbb{E}[\|\tilde{\boldsymbol{g}}\|^2]$ in a complicated manner, unlike for AdamW.

To obtain the absolute size of an update for $\boldsymbol{\omega}$ in equilibrium, we again assume that successive gradients are roughly orthogonal. Similar to AdamW, we can then write the average square size of an update as:

$$\mathbb{E}[\|\Delta_g\boldsymbol{\omega}\|^2] = \eta^2(1-\beta_1)^2\textstyle\sum_{k=0}^{t}\beta_1^{2t-2k}\mathbb{E}\left[\left\|\frac{\boldsymbol{g}_k}{\boldsymbol{v}}\right\|^2\right] \tag{85}$$

$$\approx \eta^2\tfrac{(1-\beta_1)^2}{1-\beta_1^2}\left(C - \tfrac{\lambda^2}{v^2}\widehat{\|\boldsymbol{\omega}\|}^2\right) \tag{86}$$

$$\approx \eta^2 C\tfrac{1-\beta_1}{1+\beta_1}\left(1 - \tfrac{\lambda^2\widehat{\|\boldsymbol{\omega}\|}^4}{\mathbb{E}[\|\tilde{\boldsymbol{g}}\|^2] + \lambda^2\widehat{\|\boldsymbol{\omega}\|}^4}\right) \tag{87}$$

where we used (76) twice and assumed that t was high enough to approximate the sum with its limit.

In general the dependency of $\mathbb{E}[\|\Delta_g\boldsymbol{\omega}\|^2]$ and $\widehat{\|\boldsymbol{\omega}\|}$ on $\mathbb{E}[\|\tilde{\boldsymbol{g}}\|^2]$ does not cancel out in the relative update size $\eta_r = \sqrt{\mathbb{E}[\|\Delta_g\boldsymbol{\omega}\|^2]}/\widehat{\|\boldsymbol{\omega}\|}$. The equilibrium η_r values are therefore not necessarily equal, even for scale invariant parameters. This behavior differs from all the other optimizers we studied in this work i.e. AdamW, SGDM and Lion. We believe the imbalanced rotational update sizes could be the main reason for the poor performance of L_2 regularization compared to decoupled weight decay for Adam-like methods. We explore this further in our experiments (Section 3).

In the simple system described in Appendix J, we find that Equation 84 and Equation 87 are able to predict the equilibrium norm and angular update size (via η_r) reasonably well. This is especially true when the gammas are initialized with ones instead of randomly, resulting in more uniform elements in $\boldsymbol{v}$. For random gammas or very small loss scales the predictions get worse.

J Simple System for Random Walks

Definition: We define the simple system as:

$$f(\boldsymbol{X}) = \boldsymbol{\gamma}_{\text{out}} \odot N\big(\boldsymbol{W}(\boldsymbol{\gamma}_{\text{in}} \odot \boldsymbol{X})\big) \tag{88}$$

where $\boldsymbol{X} \in \mathbb{R}^{C\times B}$, $\boldsymbol{W} \in \mathbb{R}^{K\times C}$, $\boldsymbol{\gamma}_{\text{in}} \in \mathbb{R}^{C\times 1}$, $\boldsymbol{\gamma}_{\text{out}} \in \mathbb{R}^{K\times 1}$ and N is a batch normalization function (see Equation 6) applied to each feature independently. The only learnable parameters are the weights $\boldsymbol{W}$, the gammas $\boldsymbol{\gamma}_{\text{in}}$ and $\boldsymbol{\gamma}_{\text{out}}$ are kept constant. We initialize the weights using the default initialization for a linear layer in PyTorch [26] i.e. each element is sampled independently and uniformly from the interval $[-\frac{1}{\sqrt{C}}, \frac{1}{\sqrt{C}}]$. The gammas are initialized with elements independent and identically distributed (i.i.d.) following a standard normal distribution. The inputs are also sampled i.i.d. from a standard normal distribution at each iteration. The gradients of $f(\boldsymbol{X})$, which are used to compute other gradients via the chain-rule or backpropagation, are sampled i.i.d. from a normal distribution with standard deviation $\frac{1}{KB}$ where the B simulates the typical averaging of the loss over a batch and the K gives a scale more similar to the derivatives of softmax-cross-entropy (the

difference of two vectors with an L_1-norm of 1 each). We can also scale the initial gradients further with a loss scale to obtain different gradient norms (especially important for Adam).

Rationale: We use this system to study a random walk in a neural network as described in Section 2, which serves as a simplified model of a real optimization problem. The gammas give different variances for each input and output channel, causing the second gradient moment in Adam/AdamW to vary between elements of $\boldsymbol{W}$ like they may in real neural network training due to the rest of the network. The normalization ensures that the system is scale-invariant for each row of $\boldsymbol{W}$. The randomly sampled inputs and initial gradients ensure that everything is orthogonal in expectation. Compared to a real neural network training, the dynamics of this system are simplified with no loss converging over time and steady input / gradient distributions. Other complicated effects such as dead ReLUs do also not happen in this system. This makes this simple system a good setting to study the equilibrium dynamics in a controlled manner.

Details of Figure 8: Here we use $B = 32, C = K = 128$. We use the default hyperparameters for Adam and AdamW from PyTorch 2.0 [26], except for the L_2-regularization coefficient of Adam which is zero by default but we use $\lambda = 10^{-3}$. For reference the other values are: learning rate $\eta = 10^{-3}$, first moment coefficient $\beta_1 = 0.9$, second moment coefficient $\beta_2 = 0.999$, epsilon $\varepsilon = 10^{-8}$, weight decay $\lambda = 10^{-1}$ (AdamW only). We do not use an additional loss scale for to scale the gradient norms. The experiments run for 30k steps, the plots are downsampled by a factor of 100x without averaging.

K Differences between Real Networks and a Random Walk

The random walk model we use in our analysis lets us make a variety of simplifications that would not strictly hold for real neural network optimization problems. For the real networks we therefore effectively make several approximations that can affect the accuracy of the predictions. In this section, we begin by discussing the random walk model. Subsequently, we present measurements to evaluate the accuracy of the approximations and outline the differences for real neural networks.

Random Walk Setup: The random walk setup models training dynamics of optimizing a random loss function. We can think of this as randomly sample the inputs and gradients for the network outputs at each step independently from a zero-mean normal distribution. The parameter gradients are computed via backpropagation, i.e. by using the chain rule starting from the randomly sampled output gradients. Due to the linearity of backpropagation, the gradient of each parameter coordinate will be a weighted sum of the output gradients. As a result these gradients remain zero-mean and are independent between steps, but they will accurately capture effects such as gradient orthogonality (Equation 1) and inverse proportionality (Equation 2). Additionally, because we are randomly sampling the inputs and gradients for the network outputs from the same distribution over time, the distribution of these gradients does not change over time. This is in contrast to standard optimization with a non-random loss, where we do not expect gradient independence between steps and the distribution may change over time in a non-trivial manner. We can model stochastic optimization as the sum of a noise component and an underlying true objective function (e.g. the true distribution loss). Although the noise component can easily depend on the progress on the underlying objective function, we can view the random walk as an approximation of this noise component. When the noise component dominates, which may be the case for stochastic gradient descent with sufficiently small mini-batches, the equilibrium dynamics may closely approximate that of the random walk.

Normalized Setup: We present measurements how closely the random walk model approximates the training dynamics of an original ResNet-20 trained on CIFAR-10 with a constant learning rate of $\eta = 0.01$ and a weight decay of $\lambda = 0.01$ in Figure 9. This standard ResNet has its convolutional layers followed by Batch Normalization, ensuring that the network is well-normalized. Consequently, we expect the convolutional weights to be scale-invariant.

Consistent with the expectation for this network, the angle between the gradient $\boldsymbol{g}_t$ and the weights $\boldsymbol{\omega}_{t-1}$ is close to zero. This is evident from the first row. The second row suggests that $\forall j \neq k$: $\mathbb{E}[\langle \boldsymbol{g}_j, \boldsymbol{g}_k \rangle] = 0$, which in average holds in random walk scenarios, also roughly holds here. We use $\boldsymbol{m}_{t-1} \perp \boldsymbol{g}_t$ as a measurement for this. It gives us information about the orthogonality of $\boldsymbol{g}_t$ and the previous update directions.

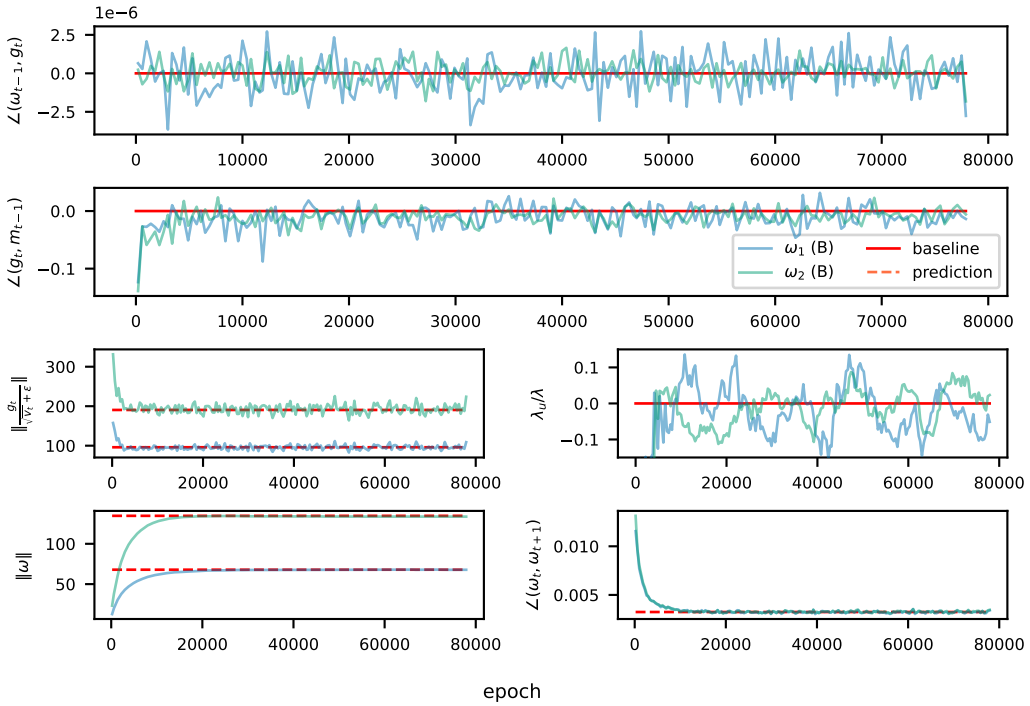

Figure 9: Measurement how closely the random walk model approximates the training dynamics of two convolutional filters of a ResNet-20 trained on CIFAR-10 with a constant learning rate.

In the third row, we assess the simplifications related to the scaled gradient $\tilde{\boldsymbol{u}}_t = \boldsymbol{g}_t/(\sqrt{\boldsymbol{v}} + \varepsilon)$, an approximation of $\boldsymbol{u}$ with constant $\boldsymbol{v}$. The left panel depicts how $\mathbb{E}[\|\tilde{\boldsymbol{u}}_t\|]$ evolves over time. Our observations indicate that it closely aligns with our approximation $\sqrt{C}$ in this setup.

We further measure the weight decay component of the scaled gradient λ_u by projecting it on the weight vector $\boldsymbol{\omega}$, $\lambda_u = \langle \boldsymbol{\omega}, \tilde{\boldsymbol{u}}_t \rangle / \|\boldsymbol{\omega}\|^2$. We take this approach to relate the weight decay component of the scaled gradient with the weight decay denoted as λ. The right panel of the third row illustrates this measurement. Notably, the gradient's weight decay component is relatively small, staying roughly within 10% of the weight decay.

Finally, in the fourth row, we compare the observed weight norms $\|\boldsymbol{\omega}_i\|$ and angular updates η_r with our predictions from Table 1. We find that the predictions closely match the measurements after the initial transient phase in this setup.

Poorly Normalized Setup: In this section, we evaluate how closely the random walk model approximates the training dynamics of a GPT-like model trained on Wikitext with learning rate $\eta = 0.0005$ and weight decay $\lambda = 0.5$ in Figure 10. The architecture does not incorporate Layer Normalization immediately after the linear layers. As a result, we do not expect the weight vectors $\boldsymbol{\omega}$ to be fully scale-invariant. The measurements in the first row supports this. For the angle between the gradient $\boldsymbol{g}_t$ and the weights $\boldsymbol{\omega}_{t-1}$ we measure a small bias in average opposed to the measurements of the normalized setup in Figure 9.

It is therefore not surprising that the scaled gradient $\tilde{\boldsymbol{u}}_t$, projected on the weight vector $\boldsymbol{\omega}$ has a more significant contribution in this setup, as evident in the right panel of the third row. As a consequence of the additional negative weight decay of the scaled gradient component—reducing the effective weight decay—our equilibrium norm prediction tends to underestimate the measured weight norm $\|\boldsymbol{\omega}\|$ and over-estimate the expected angular update η_r. By defining the error as a scaling factor of λ

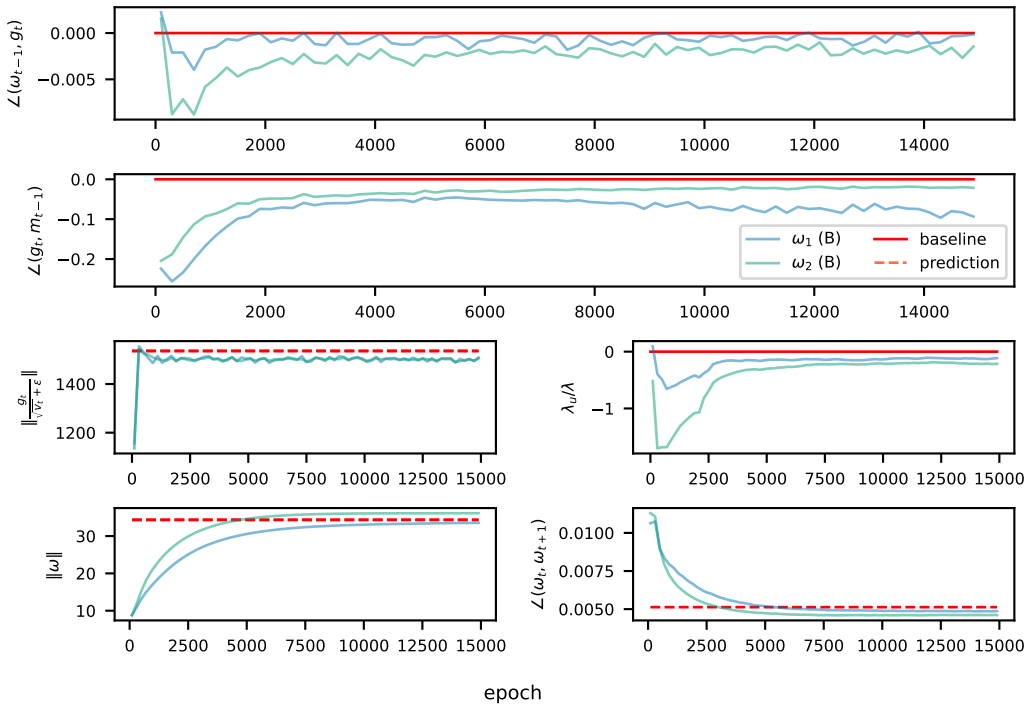

Figure 10: Measurement how closely the random walk model approximates the training dynamics of two linear layers for a GPT-like model trained on wikitext.

(represented as $\lambda_{err} = \frac{\lambda_e}{\lambda}$), we observe the following impact on our prediction

$$\eta_r \approx \sqrt{2\eta(\lambda \cdot \lambda_{err})\frac{1-\beta_1}{1+\beta_1}} = \widehat{\eta}_r \cdot \sqrt{\lambda_{err}} \tag{89}$$

$$\|\boldsymbol{\omega}\| \approx \sqrt{\frac{\eta C}{2(\lambda \cdot \lambda_{err}}}) = \widehat{\|\boldsymbol{\omega}\|} \cdot \frac{1}{\sqrt{\lambda_{err}}} \tag{90}$$

At the same time we can see from the left panel in row three, that our prediction over-estimates the scaled gradient norm. This means we tend to over-estimate the equilibrium norm with our prediction, but not the expected angular update $\widehat{\eta}_r$. For $\widehat{\eta}_r$ the scaled gradient norm cancels out. For the equilibrium norm and error estimate C_{err}, $\|\tilde{\boldsymbol{g}}\| = C \cdot C_{err}$, we have:

$$\|\boldsymbol{\omega}\| \approx \sqrt{\frac{\eta(C \cdot C_{err})}{2\lambda}} = \widehat{\|\boldsymbol{\omega}\|} \cdot \sqrt{C_{err}} \tag{91}$$

Interestingly, we observe in the last row that for $\boldsymbol{\omega}_1$ these effects on the equilibrium weight norm seem to cancel out and our prediction $\widehat{\|\boldsymbol{\omega}\|}$ holds roughly for $\|\boldsymbol{\omega}_1\|$. For $\boldsymbol{\omega}_2$ we do under-estimate the equilibrium norm but note that the error is only a few percent.

Finally, the second row suggests that the approximation $\forall j \neq k : \mathbb{E}[\langle \boldsymbol{g}_j, \boldsymbol{g}_k \rangle] = 0$, measured by $\boldsymbol{m}_{t-1} \perp \boldsymbol{g}_t$ does not strictly hold in this case. In fact it seems that consequent updates point slightly in opposite directions. This means that we expect additional negative terms in Equation (39) and thus to over-estimate the approximated RMS update size η_g in Equation (40).

Even though, we notice that the random walk model is only an approximation, our predictions hold fairly well, as evident from the last row.

L Rotational Dynamics of Scale-Sensitive Parameters

Prior work has primarily focused on the dynamics of scale-invariant weights. Note that any weight vector can be made scale-invariant by simply applying normalization to it, for example in the form of Weight Standardization [27]. For a random walk the gradient component is always orthogonal in expectation, but for real tasks scale-sensitive weights sometimes have an average radial gradient component $\mathbb{E}[\boldsymbol{u}_{\|}] \neq 0$ (see Appx E). Here we explore how this affects the rotational dynamics of these weights (for SGDM). We find that a radial component acts like an additional weight decay $\lambda_u = -\mathbb{E}[\langle \boldsymbol{u}, \boldsymbol{\omega} \rangle]/(\eta \|\boldsymbol{\omega}\|^2)$ that can be combined with λ to give a new "effective" weight decay $\lambda_e = \lambda + \lambda_u$, resulting in dynamics similar to scale-invariant weights with the adjusted value.

Most neural network architectures have some scale-sensitive parameters. This commonly includes gains and biases as well as a final fully connected layer that is typically not followed by normalization. In networks without normalization, with infrequent normalization, or poorly placed normalization, most weight vectors can be scale-sensitive. The original, un-normalized, VGG [32] architecture is a good example of this, it consists of a series of convolutional layers with ReLUs and occasional pooling layers between them and series of fully connected layers towards the end. In this section we use it to investigate the rotational dynamics of scale-sensitive weights.

First we would like to note that the magnitude of scale-sensitive weights can also be largely arbitrary. Although they can't be scaled directly without affecting the loss, we can often scale two of them without affecting the network output. Consider two successive layers with a ReLU between them:

$$f(\boldsymbol{X}, \boldsymbol{W}_1, \boldsymbol{W}_2, \boldsymbol{b}_1, \boldsymbol{b}_2) = \text{ReLU}(\boldsymbol{X}\boldsymbol{W}_1 + \boldsymbol{b}_1)\boldsymbol{W}_2 + \boldsymbol{b}_2 \tag{92}$$

where $\boldsymbol{W}_1, \boldsymbol{W}_2 \in \mathbb{R}^{C \times C}$ are weight matrices, $\boldsymbol{b}_1, \boldsymbol{b}_2 \in \mathbb{R}^{1 \times C}$ are vectors, $\boldsymbol{X} \in \mathbb{R}^{B \times C}$ are inputs and we broadcast the operations. Note that the ReLU is positively homogeneous, so for a positive scalar $r > 0$ we have:

$$f(\boldsymbol{X}, r\boldsymbol{W}_1, r^{-1}\boldsymbol{W}_2, r\boldsymbol{b}_1, \boldsymbol{b}_2) = \text{ReLU}(\boldsymbol{X}\boldsymbol{W}_1 r + \boldsymbol{b}_1 r)\boldsymbol{W}_2 r^{-1} + \boldsymbol{b}_2 = f(\boldsymbol{X}, \boldsymbol{W}_1, \boldsymbol{W}_2, \boldsymbol{b}_1, \boldsymbol{b}_2) \tag{93}$$

Assuming the weights are scaled in-place (i.e. we don't modify the computation graph, only the weight values), this type of rescaling operation scales the relative update of $\boldsymbol{W}_1$ by r^{-2} and $\boldsymbol{W}_2$ by r^2 when optimizing using SGD. This can significantly affect the learning dynamics as studied in e.g. Path-SGD [23].

For a scale-sensitive weight $\boldsymbol{\omega}$, the gradient orthogonality (1) and inverse scaling (2) do not necessarily hold. The inverse scaling holds in terms of rescaling operations like the ones mentioned above if they are applicable. Generally, the gradient has some radial component in the direction of the weight. The expected magnitude of this component depends on the average angle between the gradient and the weight as well as the expected gradient magnitude itself. If we separate the gradient into radial and perpendicular components and view the radial component as a modification of the weights decay, we have a very similar setup to the one we analyzed for scale-invariant weights. If a stable equilibrium exists, this could give rise to rotational dynamics which may vary from weight to weight based on the "effective weight decay" for each one.

We explore this with VGG-13 training on CIFAR-10 using SGDM. We compare two versions, a standard unnormalized one and a variant where weight standardization is applied to every convolutional and fully connected layer. For each one, we measure the angular updates, the weight norms and the relative radial gradient magnitude:

$$\lambda_u = \mathbb{E}[\langle \boldsymbol{\omega}, \nabla_{\boldsymbol{\omega}} \mathscr{L} \rangle / \|\boldsymbol{\omega}\|^2] = E[\cos(\angle(\boldsymbol{\omega}, \nabla_{\boldsymbol{\omega}} \mathscr{L})) \cdot \|\nabla_{\boldsymbol{\omega}} \mathscr{L}\| / \|\boldsymbol{\omega}\|] \tag{94}$$

Note that we have written this in the case of no momentum by using $-\eta \nabla_{\boldsymbol{\omega}} \mathscr{L}$ instead of $\boldsymbol{u}$, but for the standard implementation of SGDM the momentum magnifies both this version of λ_u and the standard "weight decay" (L_2 regularization) term the same way so they are comparable. The λ_u term can therefore be viewed as modifying the weight decay, the effective weight decay parameter is $\lambda_e = \lambda + \lambda_u$ and accounts for the entire radial portion of a weight update. We replace the standard λ with λ_e when showing predicted values for the unnormalized network.

The results can be seen in Figure 11 for two weights. For the first one on the left, λ_u is relatively small compared to $\lambda = 5 \cdot 10^{-4}$ and the weight behaves similarly in both setups, showing "standard" Spherical Motion Dynamics in the unnormalized setup. The equilibrium predictions match well early in training after the initial transition phase but the weight falls out of equilibrium towards the end

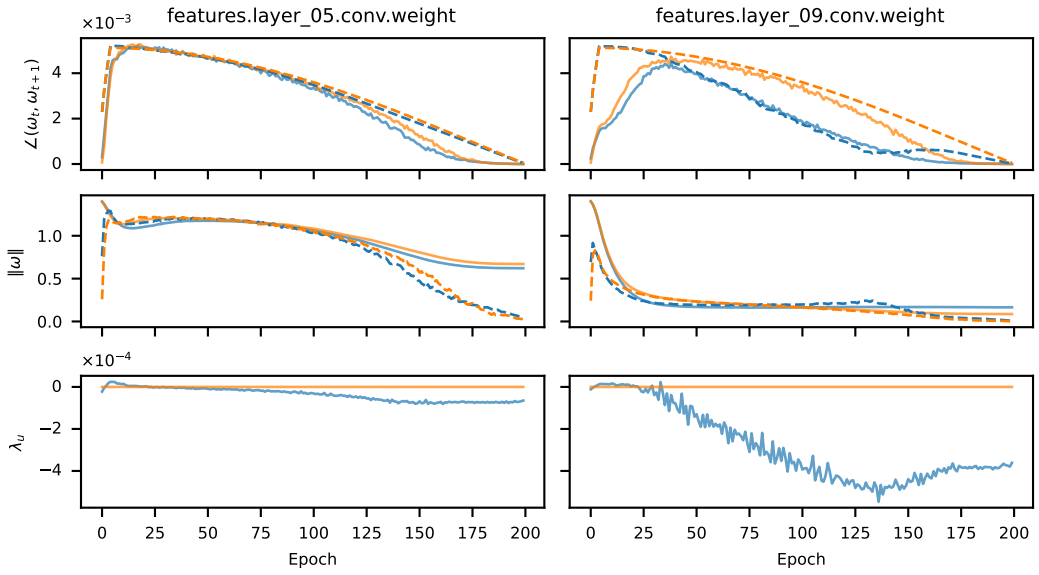

Figure 11: Measured (solid) and predicted equilibrium values (dashed) when training unnormalized (blue) and weight standardized (orange) variants of VGG-13 on CIFAR-10. The blue predictions account for the modified "effective" weight decay caused by the radial component of the gradient.

when it can't decay fast enough to keep up with the equilibrium weight magnitude. For the second weight shown on the right, λ_u is large causing a significant difference between the scale-invariant and scale-sensitive setups. The modified equilibrium predictions using λ_e capture the behavior well in the middle phase of training, after the initial transition before the weight falls out of equilibrium towards the end. We note that in the unnormalized setup λ_u changes over the course of training, starting out around 0 corresponding to an orthogonal gradient and growing larger in the later phases. This is likely due to the cross-entropy loss used, which is minimized with large (infinite) output magnitudes once the network has learned to accurately classify (overfit) the training data.

Our results for VGG13 suggest that scale-sensitive weights can also have rotational dynamics in real neural networks. The dynamics are less regular than in the normalized setup, with weights rotating at different speeds depending on the size of the radial gradient component. The weight magnitude can also not vary freely like for scale-invariant weights, where we can trade off the weight decay and learning rate without affecting the dynamics much (once equilibrium is achieved). Using large amounts of weight decay in unnormalized networks can bring the weight norms out of balance, resulting in issues like vanishing gradients or activations. In unnormalized networks the magnitude of one weight matrix also affects the gradient magnitude of all others layers, further complicating the effect of weight decay. Our rotational optimizer variants constrain the dynamics to match the equilibrium dynamics of weight standardized networks throughout training, eliminating some of these effects.

M Rotational Variants of Optimizers (RVs)

In other sections we have discussed rotational equilibrium in standard optimizers like AdamW. Training with these optimizers requires the weights to transition towards equilibrium at the start of training and when hyperparameters such as the learning rate change. This creates transient phases where rotation can be imbalanced (see e.g. Figure 1) and far away from the equilibrium value, creating a mismatch between the specified learning rate schedule and the resulting effective step sizes over time. Standard optimizers also rely on proper normalization to avoid radial gradient components that can cause imbalanced rotation in equilibrium, see Section L. In this section we create an experimental tool to study these phenomena. Algorithm 1 shows how we can create rotational variants of existing optimizers by explicitly controlling the average angular update size, forcing it to match equilibrium throughout training. This eliminates the transient phases completely and can balance the rotation of different layers and neurons without relying on normalization. By default

Algorithm 1 Our proposed Rotational Wrapper for training with constrained rotational dynamics.

Require: Inner optimizer F, decay factor $0 \leq \beta < 1$, $\varepsilon \geq 0$ for numerical stability, iteration count T, set Ω

1: **for** $\boldsymbol{p}$ in Ω **:** ▷ *For weights we choose to treat as rotational*
2: $\quad \nu_{\boldsymbol{p}} \leftarrow 0$ ▷ *Initialize the update RMS tracker*
3: $\quad n_{\boldsymbol{p}} \leftarrow \|\boldsymbol{p}\|$ ▷ *Save the initial magnitude*
4: **for** $t \in \{1, ..., T\}$ **:**
5: $\quad$ Perform backpropagation, obtain gradients for all parameters
6: $\quad$ **for all** $\boldsymbol{p}$ **:** ▷ *For each parameter*
7: $\quad\quad \Delta_g \boldsymbol{p}, \Delta_\lambda \boldsymbol{p} \leftarrow \text{F.get\_update}(\boldsymbol{p}, \nabla_{\boldsymbol{p}} \mathscr{L}(\boldsymbol{p}, \ldots))$ ▷ *Get update components (Equation (3))*
8: $\quad\quad$ **if** $\boldsymbol{p} \in \Omega$ **:** ▷ *If $\boldsymbol{p}$ should be treated as rotational*
9: $\quad\quad\quad \Delta_g \boldsymbol{p} \leftarrow \Delta_g \boldsymbol{p} / \eta$ ▷ *Remove the effect of the learning rate η used in F*
10: $\quad\quad\quad \Delta_g \boldsymbol{p} \leftarrow \Delta_g \boldsymbol{p} - \frac{\langle \Delta_g \boldsymbol{p}, \boldsymbol{p} \rangle}{\|\boldsymbol{p}\|^2} \boldsymbol{p}$ ▷ *Remove the projection onto $\boldsymbol{p}$, making $\Delta_g \boldsymbol{p} \perp \boldsymbol{p}$*
11: $\quad\quad\quad \nu_{\boldsymbol{p}} \leftarrow \beta \cdot \nu_{\boldsymbol{p}} + (1 - \beta) \cdot \|\Delta_g \boldsymbol{p}\|^2$ ▷ *Update RMS tracker*
12: $\quad\quad\quad \boldsymbol{p} \leftarrow \boldsymbol{p} + \widehat{\eta_r} \cdot n_{\boldsymbol{p}} \cdot \frac{\Delta_g \boldsymbol{p}}{\sqrt{\nu_{\boldsymbol{p}} / (1 - \beta^t)} + \varepsilon}$ ▷ *Rotate $\boldsymbol{p}$ by $\widehat{\eta_r}$ from Table 1 on average*
13: $\quad\quad\quad \boldsymbol{p} \leftarrow n_{\boldsymbol{p}} \cdot \frac{\boldsymbol{p} - \bar{\boldsymbol{p}}}{\|\boldsymbol{p} - \bar{\boldsymbol{p}}\|}$ ▷ *Center and normalize $\boldsymbol{p}$ to the initial magnitude*
14: $\quad\quad$ **else:** ▷ *Treat $\boldsymbol{p}$ as non-rotational*
15: $\quad\quad\quad \boldsymbol{p} \leftarrow \boldsymbol{p} + \Delta_g \boldsymbol{p} + \Delta_\lambda \boldsymbol{p}$ ▷ *Perform standard update*

we target the equilibrium dynamics of Weight Standardization [27]. We keep the weight magnitude constant and optionally introduce a learnable gain to compensate, which can matter for scale-sensitive weights and avoids numerical issues. Note that the form of the rotational wrapper closely resembles that of relative optimizers like LARS [39] and others discussed in Appendix B, revealing a connection to the equilibrium dynamics of standard optimizers.

In the remainder of this section, we provide further details on the algorithmic design choices used in our rotation optimizer wrapper, as shown in Algorithm 1. Note that the method can act as a wrapper around any given existing optimizer F with a known $\widehat{\eta_r}$. In cases where the true value is unknown or undesirable, we can also specify some different value of our choice.

Rotational and Non-Rotational Updates: We use Ω to specify weights we apply rotational updates to, so a parameter $\boldsymbol{p}$ is treated differently based on whether $\boldsymbol{p} \in \Omega$ or not. Note that we can choose the scale of the weights in Ω as well, e.g. each filter can be considered on its own or as a part of a larger group such as the whole layer. The rotational wrapper leaves the update of non-rotational parameters unchanged. Rotational parameters are rotated by $\widehat{\eta_r}$ on average and their magnitude is kept constant. In practice, we may choose to treat some scale-sensitive weights as rotational, constraining their magnitude. Since their magnitude can actually matter for efficient learning, we optionally introduce a learnable gain to allow the network to learn the right magnitude for these weights. This gain can be absorbed into the weights for inference.

Keeping the weight magnitude constant: Alternatively, we could vary the weight magnitude according to our derived value for the equilibrium norm. However, with a learning rate schedule this value can become arbitrarily small causing numerical issues. For scale-invariant weights the magnitude doesn't matter so we simply keep it constant. This has the added benefit of removing the inverse scaling effect of the weight norm on the gradient magnitude (2), potentially making it a more meaningful metric.

Controlling the rotation instead of the relative update: The rotation of a scale-invariant weight $\boldsymbol{\omega}$ is generally caused by both $\Delta_g \boldsymbol{\omega}$ and $\Delta_\lambda \boldsymbol{\omega}$ as can be seen in Figure 6 and 7. In equilibrium, the sum of these components is roughly orthogonal to the weight vector. We want to avoid having to apply the weight decay and our constrained magnitude is generally not equal to the equilibrium magnitude. We therefore project $\Delta_g \boldsymbol{p}$ to be orthogonal to $\boldsymbol{p}$ and control the average size of this projected version of $\Delta_g \boldsymbol{p}$ instead of the original $\boldsymbol{p}$. This lets us explicitly control the angular update, regardless of any radial component in $\Delta_g \boldsymbol{p}$ that the weight decay would eliminate on average. If we apply rotational updates to scale-sensitive weights, performing Line 11 after Line 10 prevents any radial component in the gradient from affecting the rotational speed.

Centering the weights: Different normalization setups can result in slightly different SMD properties. Layer Normalization typically makes an entire weight matrix scale-invariant whereas Batch Normalization makes individual filters (i.e., rows or columns) independent. The default form of the rotational wrapper corresponds to the rotational equilibrium dynamics obtained with Weight Standardization [27] also known as Centered Weight Normalization [10], where each filter is scale

and shift invariant. We remove the mean $\bar{\boldsymbol{p}} = \frac{1}{C}\sum_{i=1}^{C} p_i$ of $\boldsymbol{p} = [p_1, \ldots, p_C]$ since it is irrelevant in this setup. This removal was also found to be beneficial in NF-Nets [2, 3].

Hyperparameters: The algorithm requires an ε value for numerical stability but otherwise only adds one hyperparameter, a decay factor β similar to those in Adam. It determines the rate at which we update our estimate of the average update magnitude (Line 11). This in turn controls how much we let the rotation vary between steps. We could potentially derive an analytical value for β based on the convergence speed towards equilibrium. For example β should perhaps be roughly equal to $\sqrt{a}$ from Equation (36) for AdamW, when trying to match the dynamics exactly. However, this rate may not be optimal and generally depends on the learning rate (which may be scheduled). We use a default of $\beta = 0.99$ which should keep the expected angular update close to the equilibrium value over time, while still allowing some variation from step to step. There is likely batch size dependence in the optimal value of β, with larger batches potentially benefiting from smaller values since balancing the average rotation within in each step could be sufficient in these cases. An Adam-like bias correction is applied to the average update magnitude when it is used (Line 12).

Resource Requirements: We need to keep track of two scalars $\nu_{\boldsymbol{p}}$ and $n_{\boldsymbol{p}}$ for each rotational parameter. Since $\boldsymbol{p}$ is generally a vector, such as a row in a weight matrix, the memory requirement is negligible compared to quantities like momentum that store a scalar for every element of $\boldsymbol{p}$. The computational requirements in terms of floating-point operations are also relatively small, linear in the number of network (scalar) parameters like standard optimizers. However, the rotational variants are not applied fully elementwise, making efficient (parallel) implementations slightly harder.

N Additional Experiments

Imbalanced Rotation: To further quantify the impact of imbalanced rotation we experiment with artificially scaling η_r for a fraction of the neurons in each layer. Figure 12 shows the result of varying the fraction (middle) and scale for half the neurons (right), after tuning the learning rate for each configuration. We observe that even small variations in η_r can significantly affect performance.

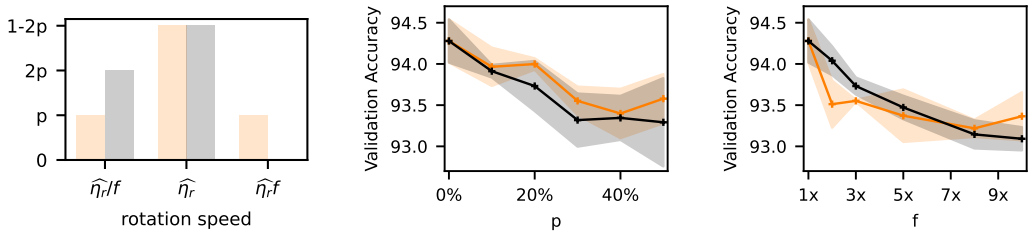

Figure 12: **Left:** Two artificially imbalanced angular update size distributions (black/orange). A portion p of the neurons is rotating f times slower and/or faster than rest using a modified RV for ResNet-18 training on CIFAR-10. **Middle:** Varying p for a fixed $f = 10$. The performance with $p = 50\%$ is comparible to a network half the width (93.5%). **Right:** Varying $f \in [1, 10]$ for $p = 50\%$.

Learning Rate vs Weight Decay for a Transformer model: In this section, we replicate the experiment previously described using a GPT-like model trained on Wikitext. The outcomes are illustrated in Figure 13. Unlike the ResNet-20 model trained on CIFAR-10, for transformers, linear layers are not fully scale-invariant; the weight norms matters. Thus, as mentioned before, we introduce a learnable gain for these layers when training with our RVs. Varying the learning rate affects the updates to biases and gains in the RV. Thus we expect the performance of this network to be more sensitive to changes in the learning rate. Again, we believe the noticeable performance difference between AdamW and RV is primarily attributed to differences in their effective step size schedules.

Constraining the Rotational Dynamics: In this section, we examine the performance of our RV of SGDM and Lion. Table 3 confirms the results we have seen with AdamW for both SGDM and Lion. Without any additional hyperparameter tuning (zero-shot) or light tuning (best-shot), our RVs closely match the performance of the original variants. This supports the idea that we can simplify deep neural network (DNN) training by standardizing the update size to η_r and skipping the initial phase.

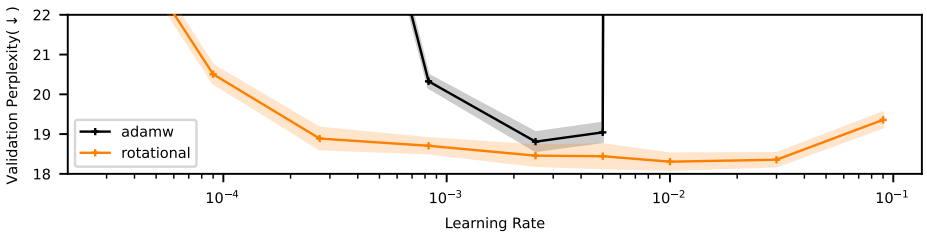

Figure 13: Validation perplexity for GPT-like model on Wikitext for different learning rate, weight decay pairs with a constant product ($\eta\lambda = 2.5{\cdot}10^{-3}$) resulting in a specific $\widehat{\eta_r}$ (Table 1).

Table 3: Test set performance for baseline optimizers SGDM and Lion and their RVs. Zero shot results for RVs use the baseline hyparparameters, the best shot is lightly tuned.

Dataset	**Model**	**Optimizer**	**Batch Size**	**Metric (⇅)**	**Baseline**	**Zero Shot**	**Best Shot**
CIFAR-10	ResNet-20	SGD	128	Top-1 Acc. (↑)	92.7 ±0.10	92.6 ±0.18	N/A
CIFAR-10	ResNet-20	SGD	2048	Top-1 Acc. (↑)	92.0 ±0.14	91.9 ±0.11	92.1 ±0.42
CIFAR-10	ResNet-20	Lion	128	Top-1 Acc. (↑)	92.1 ±0.12	91.7 ±0.13	N/A
CIFAR-10	ResNet-20	Lion	2048	Top-1 Acc. (↑)	91.8 ±0.22	91.5 ±0.12	91.6 ±0.30
Imagenet-1k	ResNet-50	SGD	256	Top-1 Acc. (↑)	77.4	77.3	N/A

Best-Shot Sweep: In practice, we have observed, that the difference between the initial learning rate, when the baseline optimizer has not yet transitioned to equilibrium, can provide a good starting point for best shot hyper-parameter tuning. Figure 14 depicts the tracked angular updates for the first 40 epochs of the DeiT tiny experiment on Imagenet-1k, as reported in Table 4. We can see that at beginning of training the learning rate of the baseline is approximately 2 times higher than for the updates of the RV. Using Table 2 and 4, we can confirm that a 2 times higher effective relative update allows the RV to match performance of the original optimizer.

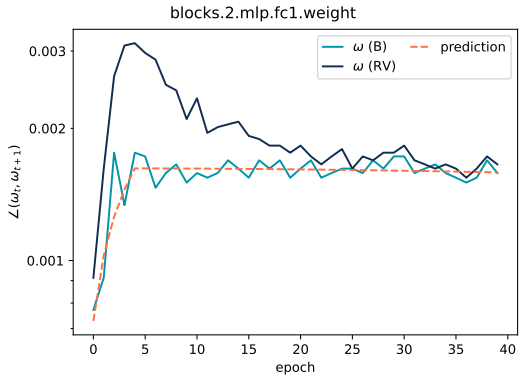

Figure 14: η_r measured for the first 40 epochs of a DeiT tiny trained on Imagenet-1k. Before AdamW transitions to equilibrium we observe a difference of roughly two times higher learning rate compared to its rotational variant.

Hyperparameter Sensitivity: The RVs introduce one hyperparameter, i.e., the decay rate β. Further, we can decide whether to enable (y) or disable (n) centering of the weights (zero-mean) and whether to enable scale-invariance on tensor (t) or filter (c) level. In this section we study the sensitivity of these choices in two different setups. We train a ResNet-18 on a random train split of CIFAR-10 with the RV of SGDM and a GPT-like model on Wikitext with the RV of AdamW. The results are shown in Figure 15. They indicate that the performance of the RVs remains relatively stable when the hyperparameters are varied. In our Adam experiments, using a special RV that combines the update direction $\Delta_g \boldsymbol{p}$ from Adam+L_2 with the $\widehat{\eta_r}$ of AdamW, we have observed that varying β noticeably affects the effective step size schedule and thus required slight tuning to increase performance.

Dataset	Model	Zero Mean	Invariance
CIFAR-10	ResNet-18	94.6±0.19 (**yc**) 94.9±0.10 (**nc**)	94.6±0.19 (**yc**) 94.6±0.47 (**yt**)
Wikitext	GPT-like	18.5 ±0.42 (**yc**) 18.5 ± 0.42 (**nc**)	18.5±0.42 (**yc**) 18.4±0.15 (**yt**)

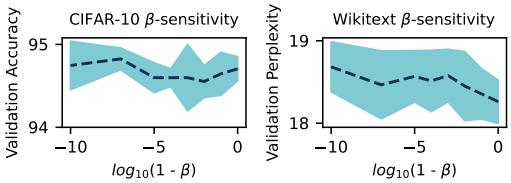

Figure 15: Experimental results for hyperparameter sensitivity. We report perplexity (↓) on Wikitext validation dataset and top-1 Acc. (↑) on a random validation split on CIFAR-10. In the default set up weight centering (y) and per filter (c) scale-invariance is enabled.

O Experimental Details

We perform our experiments on several popular datasets, i.e., CIFAR-10/100 [15] and Imagenet-1k [30] for image classification, IWSLT2014 [4] for German-English translation, and Wikitext [22] for language modelling. Our code utilizes the TIMM library [36] for vision tasks and FairSeq [24] for translation.

We train commonly used architectures, namely ResNet-20, ResNet-18, ResNet-50 [7], DeiT tiny [33], a small transformer, and a small GPT-like network [28] from scratch.

General Setup: For all experiments trained with SGDM we use a momentum of 0.9, for experiments trained with AdamW we used $\beta_1 = 0.9$ and $\beta_2 = 0.999$ and for experiments trained with our RVs we used $\beta = 0.99$, unless otherwise stated. For Lion we used $\beta_1 = 0.9$ and $\beta_2 = 0.999$ exclusively. The experiments on Imagenet-1k have been run on a single V100-SXM2-32GB GPU. All other experiments are run on a single NVIDIA A100-SXM4-40GB GPU.

Note that we used default architectures for the baseline experiments, but used a learnable gain for DeiT trained on Imagenet-1k, Transformer-S trained on IWSLT2014 de-en and the GPT-like architecture on Wikitext. As mentioned in Section M this was necessary because for transformers the layers are not fully scale-invariant and the norm of the weights matters in practice.

Here we list additional details referenced in the **details**-column in Table 4 and 5:

D-1 We pre-process the data by normalizing it with mean $(0.49140.48220.4465)$ and std $(0.20230.19940.2010)$. For training we used simple data augmentation from He et al. [7].

D-2 We use the standard data augmentation from He et al. [7] for Imagenet.

D-3 Analogously to Touvron et al. [33] we apply strong data augmentation. We use color jitter brightness up to 0.3, auto-augmentation, random erase with probability 0.25, drop path with probability 0.1, mixup with probability 0.8 and cutmix with probability 1.0. Additionally, we use label smoothing of 0.1.

Note that we used image up-scaling when DeiT tiny is trained on CIFAR-10 because DeiT tiny is designed for Imagenet images.

D-4 We use standard FairSeq library [24] with dropout probability 0.3.

Note that additionally to using weight standardization with learnable gain, we set the weight decay to 0 for scale-sensitive weights. This is done by default for the vision tasks in TIMM library [36], but not by FairSeq library [24] we used for this task. We observed no difference in performance for the baseline model, yet this adjustment allowed to tune the effective learning rate for the scale-invariant weights, without affecting the learning rate of the scale-sensitive weights significantly.

D-5 For this experiment we use the llm-baseline library library [25]. For the GPT-like architecture, we use vocabulary size of 50304, sequence length of 512, embedding size of 768. The model features 12 repeated blocks, each comprising a self-attention block followed by a two-layer MLP block with hidden dimension 3072. This results in a total of 124 million parameters. For the drop out layers we use a probability of 0.2.

Table 4: Experimental set up (include training set and test set definition).

Dataset	Model	Optimizer	Batch Size	zero shot	best shot	Wrapped Adam+$L2$	details	lr schedule	warmup	epochs (e)/ iteration (it)	train duration	precision
CIFAR-10	ResNet-20	SGD	128	wd=1e−4 lr=0.5	N/A	N/A	(D-1)	cosine	lr=1e−6 5 epochs	200 (e)	35min	float32
CIFAR-10	ResNet-20	SGD	2048	wd=1e−4 lr=4.8	wd=1e−4 lr=16	N/A	(D-1)	cosine	lr=1e−6 5 epochs	200 (e)	35min	float32
CIFAR-10	ResNet-20	AdamW	128	wd=1e−2 lr=5e−2	N/A	$\beta = 0.9$	(D-1)	cosine	lr=1e−6 5 epochs	200 (e)	35min	float32
CIFAR-10	ResNet-20	AdamW	2048	wd=1e−2 lr=1.6e−1	wd=1e−2 lr=0.8	$\beta = 0.0$	(D-1)	cosine	lr=1e−6 5 epochs	200 (e)	35min	float32
CIFAR-10	ResNet-20	Lion	128	wd=1.0 lr=5e−4	N/A	N/A	(D-1)	cosine	lr=1e−6 5 epochs	200 (e)	35min	float32
CIFAR-10	ResNet-20	Lion	2048	wd=1.0 lr=1.6e−2	wd=1.0 lr=8e−2	N/A	(D-1)	cosine	lr=1e−6 5 epochs	200 (e)	35min	float32
Imagenet-1k	ResNet-50	SGD	256	wd=1e−4 lr=1e−1	N/A	N/A	(D-2)	cosine	lr=1e−6 5 epochs	90 (e)	30h	float16
Imagenet-1k	DeiT tiny	AdamW	1024	wd=5e−2 lr=5e−4	wd=1e−1 lr=5e−4	N/A	(D-3)	cosine	lr=1e−6 5 epochs	300 (e)	70h	float16
CIFAR-10	DeiT tiny	AdamW	64	wd=5e−2 lr=5e−4	N/A	$\beta = 0.99$	(D-3)	cosine	lr=1e−6 5 epochs	600 (e)	16h	float16
IWSLT2014 de-en	Transformer-S	AdamW	4096	wd=1e−4 lr=5e−4 $\beta_2 = 0.98$	wd=2e−1 lr=5e−4 $\beta_2 = 0.98$	$\beta = 0.99$	(D-4)	cosine	4000 (it)	22021 (it)	50min	float16
Wikitext	GPT-like	AdamW	55	wd=0.5 lr=5e−3 $\beta_2 = 0.95$	N/A	$\beta = 0.9$	(D-5)	cosine div_f=1e2 final_div_f=1e4	2e−2 (%)	15000 (it)	3h	bfloat16

Constraining the Rotational Dynamics: The experimental details for the experiments reported in Table 2 and 3 can be found in Table 4. Note that for the Wrapped Adam+L_2 experiments we used the best shot settings. For these experiments, we have observed that varying β noticeably affects the effective step size schedule. Thus we used minimal hyperparameter tuning for β. The parameter configuration is shown in the respective column of Table 4.

Learning Rate vs Weight Decay: For the ResNet-20 experiment on CIFAR-10 and language model task on Wikitext with a GPT-like model we use the best shot (zero shot) setting reported in Table 4 as default. We then sweep over the learning rate keeping $\eta\lambda = 5{\cdot}10^{-4}$, $\eta\lambda = 2.5{\cdot}10^{-3}$ respectively, constant.

Scheduling Effects: In the experiment described in the preceding paragraph, we monitored $\|\boldsymbol{\omega}\|$, $\angle(\boldsymbol{\omega}_t, \boldsymbol{\omega}_{t+1})$ during training for one of the runs with the chosen learning rates: $1{\cdot}10^{-4}$, $8.3{\cdot}10^{-3}$, and $3{\cdot}10^{-1}$.

Adam vs AdamW: For the sweep we train a ResNet-18 on a 90/10 train/val split from the original train set. We use a step-wise cosine schedule and train for 200 epochs without warmup. In this sweep we try to reproduce the results in Figure 2 from [21], albeit with a slightly different network and training for 200 epochs instead of 100. The best configuration for Adam was $\eta = 7.813 \cdot 10^{-4}$, $\lambda = 1.250 \cdot 10^{-4}$ resulting in a validation set accuracy of 93.919. The best configuration for AdamW was $\eta = 1.25 \cdot 10^{-2}$, $\lambda = 8.0 \cdot 10^{-2}$ with a validation accuracy of 94.319. On the test set we run each configuration over 5 different seeds, using the AdamW hyperparameters for both RV-AdamW and the special wrapped Adam+L_2 RV. The results were Adam+L_2 94.08 ± 0.16, AdamW 94.74 ± 0.14, RV-AdamW 94.57 ± 0.12 and the special wrapped Adam+L_2 94.55 ± 0.07. Note that the performance of the two RVs is almost identical and higher than standard Adam+L_2, but slightly lower than AdamW which is not surprising for a zero-shot transfer of a well tuned baseline and likely due to scheduling effects.

Imbalanced Rotation: We trained a ResNet-18 on a 90/10 train/val split from the original train set with weight decay $\lambda = 0.01$ and varying learning rates $\eta \in \{2.7{\cdot}10^{-3}, 8.3{\cdot}10^{-3}, 2.5{\cdot}10^{-2}, 5{\cdot}10^{-2}, 1{\cdot}10^{-1}, 3{\cdot}10^{-1}, 0.9{\cdot}10^{-1}, 4.5\}$). For each rotation speed scaling f and portion p we report the best performance of this sweep. All other settings are equivalent to the settings reported in Table 4.

Training Poorly Normalized Networks: We train a ResNet-18 with layer normalization on CIFAR-100 using the same augmentation, learning rate schedule and base hyperparameters as for the ResNet-20 on CIFAR-10 experiments, unless otherwise noted below. We train on a random subset containing 90% of the train set and use the remaining 10% for validation which we report. The inputs are normalized for mean `(0.5071, 0.4867, 0.4408)` and std `(0.2675, 0.2565, 0.2761)`. We use a weight decay of $5 \cdot 10^{-4}$ and a batch size of 256. The layer normalization is implemented with a Group Normalization [38] using a single group.

Need for Learning Rate Warmup: These experiments follow the same base setup as we report in Table 4. We train for a total of 10 epochs using a cosine decay schedule (applied stepwise) and no warmup. We use local accumulation on top of batches of size 256 to emulate larger batch sizes.

Hyperparameter Sensitivity: The experimental details for the experiments reported in Figure 15 can be found in Table 5.

Table 5: Experimental details for hyperparameter senstivity study.

Dataset	Model	Optimizer	training dataset	validation dataset	hyper-parameters	details	lr schedule	warm up	epochs (e)/ iterations (it)	train duration	precision
CIFAR-10	ResNet-18	SGD	90% Train	10% Train	wd=1e−4 lr=0.5	(D-1)	cosine	lr=1e−6 5 epochs	200 (e)	35min	float32
Wikitext	GPT-like	AdamW	Train	Validation	wd=0.5 lr=4e−3 $\beta_2 = 0.95$	(D-5)	cosine div_f=1e2 final_div_f=1e4	2e−2 (%)	15000 (it)	3h	bfloat16

P Extended Discussion & Conclusion

We believe rotational dynamics can provide valuable insights into many phenomena in deep learning optimization. Our analysis shows how the average angular update size η_r in equilibrium and the RMS gradient size η_g have a different dependency on the learning rate η and weight decay λ (Table 1). In the common setting where gains and biases are excluded from weight decay, this results in a scaling of their "effective" update size, η_g, compared to that of the rotational weights, η_r. In equilibrium, η and λ therefore jointly determine two effective update sizes used for different types of parameters. Curiously, $\eta_r \propto \sqrt{\eta}$ while $\eta_g \propto \eta$, raising questions about the impact of this when varying the learning rate η with a schedule or scaling the batch size. We do not explore this here but note that this is not the case in most relative optimizers, such as Nero [20] for example.

In our analysis and experiments we found that the equilibrium in Adam with L_2 regularization significantly differs from that of AdamW and other optimizers, as the expected angular update in equilibrium depends on the gradient magnitude. This results in unbalanced rotation across vectors that have different gradient norms. Forcing Adam+L_2 to have balanced rotation through our rotational wrapper seems to eliminate most of the performance degradation compared to AdamW.

Normalization can make weights scale-invariant, ensuring the gradients are orthogonal to the weights. A consistent gradient component parallel to a weight vector results in a faster or slower rotation in equilibrium (Section L), potentially resulting in unbalanced rotation across different vectors. Different types of normalization result in different granularity of scale-invariance e.g. layer level with Layer Normalization and neuron level with Batch Normalization and Weight Standardization. With more coarse normalization operations this can result in unbalanced equilibrium rotation on the neuron level. We observe that the use of Rotational Optimizer Variants can improve the performance of layer normalized ResNets and also the GPT-like model, likely by enforcing balanced rotation. This could also help explain how Weight Standardization helps aid optimization when applied on top of Layer Normalization or Group Normalization, i.e. by ensuring scale-invariance on a finer level.

In standard optimizers different weight vectors must converge to equilibrium. Depending on the initialization weight magnitude, learning rate, weight decay and gradient norm (for SGDM only), this can result in faster or slower angular updates in the initial transient phase. This causes a mismatch between the specified learning rate schedule and the resulting angular step sizes over time. A learning rate warmup may potentially be needed in part to counteract fast initial rotation due to this effect. Using the rotational optimizer variants (RVs) eliminates this effect, ensuring angular updates follow the specified learning rate schedule. This may also explain why other relative optimizers like LARS seem to have a reduced need for learning rate warmup. Although this may not be the sole reason a warmup is needed in every case, we believe it is an important effect to be aware of.

In this project we have explored the importance of rotational dynamics in deep learning optimization aiming for a high level understanding. We hope our insights can be of use to practitioners when debugging or tuning neural network training, and that they may provide a new perspective for theorists. We see a lot of potential for future work in this area in terms of more formal and rigorous theory as well as practical application to e.g. optimizer and scheduler design.