

Not All Dimensions Are Equal: Outlier-Aware Attention Linearization with Grouped Quadratic Kernels

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Linearizing softmax attention offers a route to efficient Transformer inference, but preserving pretrained attention remains challenging. We identify an important source of approximation error overlooked by standard linearization methods: a small number of outlier dimensions contribute disproportionately to query–key dot products and account for a large fraction of the approximation error. Motivated by this heterogeneity, we propose Outlier-Aware Attention Linearization (OAL), a dimension-aware method for constructing quadratic attention approximations. Starting from a second-order Taylor approximation, we introduce a Hadamard-product formulation that makes individual query–key contributions explicit. This formulation relaxes globally shared coefficients to support dimension-specific linear terms and cross-dimensional quadratic interactions while preserving separable evaluation. To retain this flexibility with fewer parameters, OAL groups dimensions with similar contribution scales to share coefficients, while allowing distinct treatment of the few outliers. The resulting positive quadratic kernel is calibrated against pretrained attention outputs and admits linear-time evaluation with respect to sequence length for fixed head dimensions. On Qwen2.5-0.5B, OAL reduces frozen-attention output relative MSE by approximately 8.9% compared with DiJiang. Replacing middle-layer attention modules and jointly fine-tuning the kernel coefficients with LoRA further lowers WikiText-103 perplexity from 31.71 to 27.50 compared with a learned global quadratic, while improving accuracy on PIQA, HellaSwag, and ARC-E. These results demonstrate the value of outlier-aware dimension grouping for both attention approximation and pretrained-model conversion. The code is available at [GitHub](#).

1. Introduction

Softmax attention is central to modern Transformers, but computing all pairwise token interactions incurs quadratic computation as context length grows (Vaswani et al., 2017). Linear attention offers an alternative: separable query and key features permit recurrent evaluation with linear sequence-length complexity for a fixed feature dimension (Katharopoulos et al., 2020). Approximating softmax attention in pretrained Transformers with linear attention therefore offers a promising route to improving inference efficiency while reusing pretrained representations. Realizing this potential, however, requires approximations that closely preserve the original attention behavior, particularly when only limited adaptation is available to compensate for approximation errors.

Existing methods approximate the softmax kernel using feature maps that separate query and key computations, or low-order polynomials that admit efficient evaluation (Choromanski et al., 2021; Nguyen et al., 2026; Chen et al., 2024b; Nauen et al., 2024). To better preserve pretrained attention, these approximations can also be learned or adapted to match the original model’s attention behavior (Chowdhury et al., 2022; Zhang et al., 2025; Lan et al., 2025). Despite these advances,

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approximation errors are still to be desired under the limited-adaptation setting. To better understand the remaining approximation error, we examine how it relates to the contributions of individual query–key dimensions.

Our analysis reveals that a small subset of query–key dimensions contributes disproportionately to attention logits (Section 3). Diagnostic experiments further show that approximating these dimensions alone can induce substantial attention-output error. This suggests that the approximation should have sufficient flexibility to accommodate differences in dimension-wise contributions. The representative approaches discussed above generally formulate their approximations through vector-level feature mappings or functions of the aggregated query–key score. Their parameterization, however, is not explicitly organized around the unequal contributions of individual query–key dimensions. Our findings motivate making these contributions explicit and using their statistics to guide the approximation, so that outlier dimensions can be treated differently from the remaining dimensions.

Therefore, we propose *Outlier-Aware Attention Linearization* (OAL), which explicitly adapts the approximation to dimension-wise heterogeneity in pretrained attention. We use a quadratic kernel as a tractable starting point for introducing dimension-specific parameters while retaining separable evaluation. To enable dimension-wise treatment, we introduce a Hadamard-product formulation that makes individual query–key contributions explicit. This formulation generalizes the globally shared coefficients in a quadratic approximation to dimension-specific linear coefficients and coefficients for cross-dimensional quadratic interactions. It thereby allows different dimensions to be approximated differently while preserving the separable structure required for linear-time evaluation.

A fully dimension-specific parameterization, however, introduces many coefficients. The empirical distribution of dimension-wise contributions suggests that much of this flexibility can be shared. This observation motivates grouping dimensions with similar contributions and sharing coefficients within groups and across corresponding group pairs. The resulting parameterization accommodates outlier dimensions while keeping the number of learned coefficients small. With the groups fixed, we calibrate the resulting positive quadratic kernel against pretrained softmax attention outputs using frozen queries, keys, and values.

On frozen Qwen2.5-0.5B attention states, OAL achieves the lowest error among the evaluated methods, reducing attention-output relative MSE by approximately 8.9% compared with DiJiang (Chen et al., 2024b). Under a common middle-layer attention replacement protocol with matched LoRA adaptation (Hu et al., 2022), OAL outperforms the evaluated converted methods on WikiText-103 perplexity and all three downstream tasks on both Qwen2.5-0.5B and Llama-3.2-1B (Llama Team, 2024). Relative to the learned global quadratic, it lowers Qwen2.5-0.5B perplexity from 31.71 to 27.50, although a gap to the LoRA-adapted softmax reference remains on both backbones. OAL also requires only $O(N)$ attention computation for a sequence of length N and fixed head dimensions, without explicitly constructing the $N \times N$ attention matrix.

Our contributions are threefold:

- We characterize dimension-wise heterogeneity in pretrained query–key interactions and its association with attention approximation error, motivating contribution-aware kernel design.
- We develop a dimension-aware quadratic formulation with separable evaluation, and introduce contribution-based grouping to retain this flexibility with a compact parameterization.
- We demonstrate improved frozen-attention fidelity and adapted model quality under a matched conversion protocol, and examine the role of group-aware flexibility through a diagnostic budget study.

2. Preliminaries

Softmax Attention For a single head, let $Q, K \in \mathbb{R}^{N \times d_h}$ and $V \in \mathbb{R}^{N \times d_v}$ denote the query, key, and value matrices, with sequence length N , query–key dimension d_h , and value dimension d_v . We write q_i, k_i, v_i for their respective token vectors at position i . Scaled dot-product attention (Vaswani et al., 2017) is

$$\text{Attn}(Q, K, V) = \text{Softmax}\left(\frac{QK^\top}{\sqrt{d_h}}\right)V,$$

with row-wise softmax. Direct evaluation costs $O(N^2(d_h + d_v))$ operations; explicit storage of the score matrix requires $O(N^2)$ memory.

Linear Attention A common class of linear attention methods replaces the pairwise attention kernel with a separable feature representation (Katharopoulos et al., 2020; Choromanski et al., 2021). Specifically, the softmax kernel can be approximated as $\exp(q^\top k / \sqrt{d_h}) \approx \kappa(q, k) = \phi(q)^\top \phi(k)$, where $\phi(\cdot) \in \mathbb{R}^m$ denotes a feature map. Substituting this factorization into normalized causal attention yields

$$\hat{y}_i = \frac{\left(\sum_{j \leq i} v_j \phi(k_j)^\top\right) \phi(q_i)}{\phi(q_i)^\top \left(\sum_{j \leq i} \phi(k_j)\right)}. \quad (1)$$

Here, $\hat{y}_i$ denotes the approximate attention output at position i . The key–value and key statistics in parentheses can be accumulated incrementally over the causal prefix. This reassociation avoids explicitly constructing the $N \times N$ attention matrix and gives linear scaling with sequence length for a fixed feature dimension and memory size (Gu and Dao, 2023; Sun et al., 2023; Peng et al., 2023; Chen et al., 2024a). Different linearization methods therefore mainly differ in the choice or construction of the feature map $\phi(\cdot)$ and the resulting approximation to the softmax kernel.

Softmax Kernel Approximation Separable softmax approximations can be obtained by constructing feature maps directly (Katharopoulos et al., 2020; Choromanski et al., 2021; Peng et al., 2021) or by expanding a low-order polynomial in the attention score into separable terms. We focus on the latter. Polynomial kernels based on low-order Taylor approximations have been explored in efficient attention and language modeling (Arora et al., 2024; Nauen et al., 2024). For a scalar attention score s , the second-order Taylor approximation is $\exp(s) \approx 1 + s + \frac{1}{2}s^2$. Allowing its coefficients to vary gives the global quadratic $p(s) = a + bs + cs^2$, where a , b , and c control the constant, linear, and quadratic terms. This is the starting point for our dimension-aware kernel.

3. Dimension-Wise Heterogeneity in Pretrained Attention

Motivated by activation outliers in pretrained Transformers (Kovaleva et al., 2021; Dettmers et al., 2022; Sun et al., 2024; Li et al., 2026b,a), where a small number of activation or representation dimensions can attain unusually large magnitudes or exert disproportionate influence on model behavior, we examine whether query–key contributions are similarly concentrated in a few dimensions and whether this concentration is associated with attention-approximation error.

3.1. Query–Key Outliers and Dimension-Wise Contributions

Figure 1 shows the post-RoPE query and key magnitudes for a representative Qwen2.5-0.5B head. Since either magnitude alone does not directly measure a dimension’s contribution to attention logits,

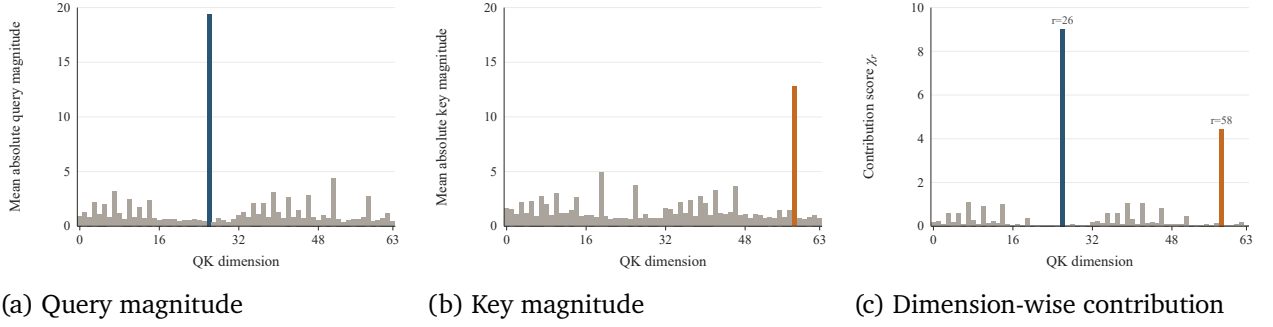


Figure 1 | **Dimension-wise heterogeneity in a representative Qwen2.5-0.5B attention head.** (a) Mean absolute query magnitude across QK dimensions. (b) Mean absolute key magnitude across the corresponding KV head. (c) Contribution score χ_r defined in Eq. (2). Dimensions $r = 26$ and $r = 58$ are highlighted across panels.

we analyze their coordinate-wise products for each layer–head pair, suppressing layer and head indices below. For calibration input p , let N_p denote its sequence length, and let $q_{p,i}, k_{p,j} \in \mathbb{R}^{d_h}$ denote the post-RoPE query and key vectors at positions i and j . Their r -th coordinates are denoted by $q_{p,i,r}$ and $k_{p,j,r}$. We define the causal pair set as $\Omega_p = \{(i, j) : 1 \leq j \leq i \leq N_p\}$.

We define the contribution of dimension r on input p and its equally weighted average over the P calibration inputs as

$$\chi_{p,r} = \frac{1}{|\Omega_p|} \sum_{(i,j) \in \Omega_p} \left| \frac{q_{p,i,r} k_{p,j,r}}{\sqrt{d_h}} \right|, \quad \chi_r = \frac{1}{P} \sum_{p=1}^P \chi_{p,r}. \quad (2)$$

The contribution profile in Figure 1 is concentrated on a few dimensions. The highlighted dimensions, $r = 26$ and $r = 58$, have high contribution scores and correspond to prominent coordinates in the query or key representations. A similarly concentrated dimension-wise contribution profile is observed in GPT-2 Medium; additional cross-model evidence is provided in Appendix A.

3.2. Approximation Error Concentration on High-Contribution Dimensions

We next test whether approximating only the high-contribution dimensions retains a substantial fraction of the full-head output error. For each evaluated attention head, we select a small subset S of high-contribution dimensions according to the pretrained contribution scores. We apply each approximation only to S while keeping the remaining dimensions exact. Let $\mathcal{E}_S$ denote the resulting attention-output MSE and $\mathcal{E}_{\text{full}}$ the error when the same method is applied to the full head. We report the retained-error ratio $R_S = \mathcal{E}_S / \mathcal{E}_{\text{full}}$. For the Gaussian controls, pretrained Q/K representations are replaced with i.i.d. $\mathcal{N}(0, 1)$ samples and the same contribution-based grouping and intervention procedures are applied. Further metric and control details are provided in Appendix A.

Approximating only about 5.5% of the dimensions retains a substantial fraction of the full-head attention-output MSE across all evaluated methods, with retained-error ratios consistently above the corresponding Gaussian controls (Table 1). In a separate random-coordinate control with 1,000 cardinality-matched subsets sampled from the pretrained bulk dimensions, none of the null replicates exceeded the observed retained-error ratio for any method, yielding empirical $p < 0.001$ in all cases. These retained-error ratios suggest an association between high contribution scores and attention-output approximation error. Appendix A provides additional control analyses and stability analyses across held-out inputs and sequence lengths for the representative Qwen2.5-0.5B head. These observations motivate the contribution-aware parameter sharing in Section 4, which allows

Table 1 | **Approximation error concentration on high-contribution dimensions.** The selected subsets account for approximately 5.5% of the evaluated head dimensions in aggregate. We report the retained-error ratio R_S (%) for the high-contribution subsets and for Gaussian controls obtained by replacing pretrained Q/K with i.i.d. $\mathcal{N}(0, 1)$ samples while keeping the remaining protocol unchanged. Higher values indicate stronger error concentration.

Method	High-contrib. error (%)	Gaussian control (%)
Performer	37.8	4.7
SpectraFormer	58.1	18.1
DiJiang	64.8	19.4
TaylorShift	63.9	13.6

dimensions with different contribution scales to use different approximation coefficients.

4. Outlier-Aware Attention Linearization

We introduce dimension-specific linear and cross-dimensional quadratic coefficients, with a Hadamard–diagonal (HD) formulation enabling exact kernel evaluation in linear time with respect to sequence length for fixed head dimensions. Contribution-based grouping then shares coefficients across dimensions and dimension pairs; enforcing positive kernel weights yields OAL.

4.1. Dimension-Aware Linearization

For a query–key pair $q, k \in \mathbb{R}^{d_h}$, let $x_r = q_r k_r / \sqrt{d_h}$ denote the contribution of dimension r , so that the attention score is $s = \sum_r x_r$. A global quadratic approximation to e^s can be written as

$$p(s) = a + bs + cs^2 = a + \sum_{r=1}^{d_h} b x_r + \sum_{r=1}^{d_h} \sum_{t=1}^{d_h} c x_r x_t.$$

All linear contributions therefore share b , and all quadratic interactions share c . TaylorShift uses this scalar polynomial structure with coefficients prescribed by the Taylor expansion (Nauen et al., 2024). To accommodate the dimension-wise heterogeneity observed in Section 3, we relax this sharing and allow separate coefficients for each dimension and dimension pair:

$$\kappa_{\text{dim}}(q, k) = a + \sum_{r=1}^{d_h} b_r x_r + \sum_{r=1}^{d_h} \sum_{t=1}^{d_h} C_{rt} x_r x_t, \quad C = C^\top. \quad (3)$$

Here, r and t index the d_h query–key dimensions; b_r weights the contribution of dimension r , C_{rr} weights its squared term, and the off-diagonal entries of C weight cross-dimensional interactions. Setting $b_r = b$ and $C_{rt} = c$ recovers the global quadratic. Degree two is the lowest degree that retains both squared terms and cross-dimensional interactions, using $O(d_h^2)$ features.

Hadamard-diagonal formulation Evaluating the quadratic kernel explicitly over all query–key pairs still incurs quadratic cost in sequence length. To avoid this pairwise computation, we use the following Hadamard–diagonal (HD) identity for the shared-coefficient quadratic term. For value channel $V_{:, \ell}$, let $D_\ell \in \mathbb{R}^{N \times N}$ be the diagonal matrix with entries $(D_\ell)_{jj} = V_{j, \ell}$. Then

$$[(QK^\top) \odot (QK^\top)] V_{:, \ell} = \text{diag}(Q[K^\top D_\ell K] Q^\top). \quad (4)$$

Here, $\odot$ denotes the Hadamard product, and diag extracts the diagonal entries of a matrix as a vector. By first accumulating $K^\top D_\ell K$ and then evaluating only the diagonal entries, the quadratic term can be computed in $O(Nd_h^2)$ time per value channel without forming an $N \times N$ matrix. For fixed head dimensions, this cost is linear in sequence length.

The (r, t) entry of $K^\top D_\ell K$ is $\sum_j V_{j,\ell} k_{j,r} k_{j,t}$, which sums value-weighted key products over token positions j . For a query at position i , restricting this sum to $j \leq i$ gives causal prefix statistics. The attention denominator uses the same kernel weights as the numerator, but replaces each value with 1. Thus, augmenting the values as $u_j = [v_j; 1]$ allows both to be accumulated together. Including the constant and linear kernel terms gives the states

$$H_{0,i} = \sum_{j \leq i} u_j, \quad H_{1,i,r} = \sum_{j \leq i} k_{j,r} u_j, \quad H_{2,i,rt} = \sum_{j \leq i} k_{j,r} k_{j,t} u_j. \quad (5)$$

Each $H_{2,i,rt}$ is a $(d_v + 1)$ -dimensional vector.

The query combines these statistics using the coefficients a , b_r , and C_{rt} from Eq. (3):

$$\begin{aligned} \tilde{u}_i &= aH_{0,i} + \frac{1}{\sqrt{d_h}} \sum_r b_r q_{i,r} H_{1,i,r} + \frac{1}{d_h} \sum_{r,t} C_{rt} q_{i,r} q_{i,t} H_{2,i,rt}, \\ \hat{y}_i &= \tilde{u}_{i,1:d_v} / \tilde{u}_{i,d_v+1}. \end{aligned} \quad (6)$$

This recurrence evaluates the specified quadratic kernel exactly without forming a token-by-token matrix (Appendix D). The last channel supplies the normalization denominator, whose positivity follows from the OAL parameterization in Section 4.3.

A fully dimension-specific quadratic has $O(d_h^2)$ coefficients. Contribution-based sharing reduces this parameter count while retaining the same states and readout.

4.2. Outlier-Aware Dimension Grouping

The contribution profile in Figure 1 suggests a compact sharing structure: low-contribution dimensions can share parameters, while the high-contribution tail may benefit from finer parameterization. Figure 2(a) illustrates a partition for the representative head in Section 3.1. For a partition $\{\mathcal{G}_g\}_{g=1}^G$ of the dimensions, recall that $x_r = q_r k_r / \sqrt{d_h}$ as the contribution of dimension r and $s_g = \sum_{r \in \mathcal{G}_g} x_r$ as the total contribution of group g . Thus, $s = \sum_{g=1}^G s_g$. Sharing b_r within each group and C_{rt} within each group pair reduces the number of learned coefficients to $O(G^2)$ while preserving the HD evaluation.

For each layer and query head, we sort dimensions by their calibration contribution scores χ_r and identify large gaps in the high-contribution tail that remain stable under prompt-level bootstrap resampling, following resampling-based stability assessment (Suzuki and Shimodaira, 2006; Hennig, 2007). These boundaries determine the group count, with at least two groups. Groups remain fixed during coefficient fitting and model adaptation. The procedure yields 2–5 groups per head for Qwen2.5-0.5B and 2–4 for Llama-3.2-1B; full details are provided in Appendix B.

As a separate diagnostic, Figure 2(b) illustrates the trade-off between attention-output MSE and parameter count across fixed group budgets for the representative head. For this head, a small number of groups captures most of the error reduction, with diminishing gains as the budget increases.

4.3. The OAL Algorithm

OAL uses a positive quadratic kernel whose coefficients are shared according to contribution-based dimension groups. We specify this parameterization below and combine it with the HD formulation to summarize the complete method in Algorithm 1.

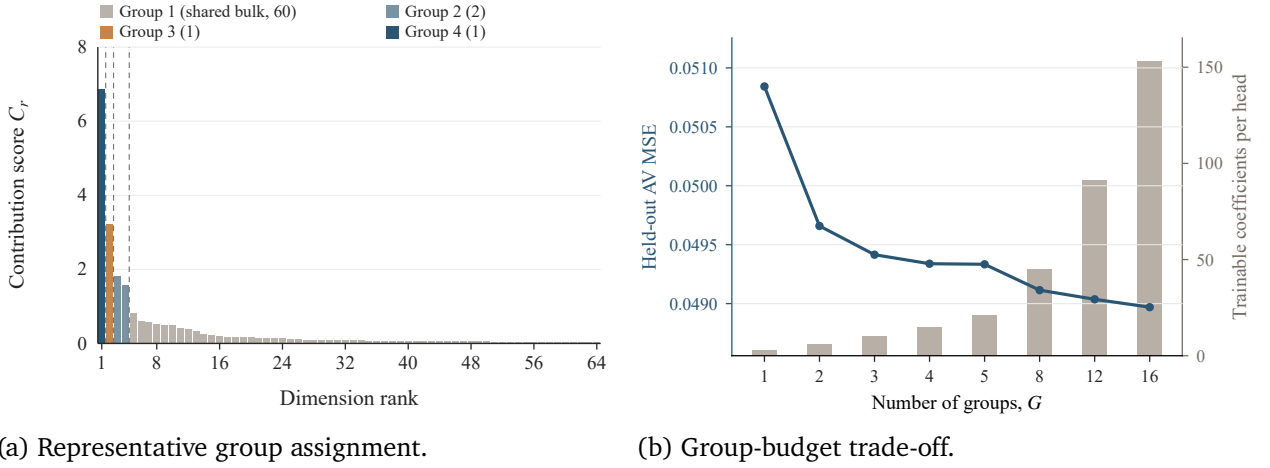


Figure 2 | **Contribution-aware dimension grouping.** (a) Sorted contribution scores χ_r and the resulting OAL groups for the representative Qwen2.5-0.5B Layer 4, Head 0. (b) Attention-output MSE and trainable coefficients per head under fixed group budgets for the same head. All budgets use the same data splits and coefficient-fitting protocol.

Positive group-aware kernel Let $g(r)$ denote the group containing dimension r . To enforce positive kernel weights while sharing coefficients within groups and group pairs, we introduce a learnable lower-triangular matrix $L \in \mathbb{R}^{(G+1) \times (G+1)}$ for each layer and query head and set $M = LL^\top$. The coefficients are $a = \epsilon + M_{00}$, $b_r = 2M_{0,g(r)}$, and $C_{rt} = M_{g(r),g(t)}$, with $\epsilon > 0$. With $\bar{\mathbf{s}} = [1, s_1, \dots, s_G]^\top$, this parameterization gives

$$\kappa_{\text{OAL}}(q, k) = \epsilon + \bar{\mathbf{s}}^\top LL^\top \bar{\mathbf{s}}, \quad \epsilon > 0, \quad (7)$$

Since $LL^\top \succeq 0$, we have $\kappa_{\text{OAL}}(q, k) \geq \epsilon > 0$, ensuring a positive normalization denominator.

With the groups fixed, we learn L by minimizing the squared error between normalized OAL and softmax attention outputs on an offline calibration set, using frozen pretrained queries, keys, and values. Calibration details are provided in Appendix B. The calibrated factors initialize subsequent joint adaptation with LoRA, during which the groups remain fixed.

Causal evaluation The coefficient mapping above expresses the grouped kernel in the dimension-wise form of Eq. (3), so it can be evaluated using the HD states and readout in Eqs. (5) and (6). This computation is exact for the specified quadratic kernel; approximation arises only from replacing the softmax kernel. Algorithm 1 summarizes the offline calibration and causal evaluation stages. The equivalence proof is provided in Appendix D.

4.4. Complexity Analysis

Direct pairwise evaluation costs $O(N^2(d_h + G^2 + d_v))$, whereas the causal evaluation stage of Algorithm 1 costs $O(Nd_h^2(d_v + 1))$. The HD states require $O(d_h^2(d_v + 1))$ space, excluding inputs and outputs. For fixed head dimensions, computation is linear in sequence length and state space is independent of it. Grouping reduces the learned parameter count to $O(G^2)$, while the implementation retains dimension-pair states. Detailed counts appear in Appendix D.

Figure 3 reports causal attention forward latency and peak GPU memory under the setup in Appendix E. OAL exhibits approximately linear runtime growth at longer sequences. With the same HD block-GEMM backend, OAL ($G = 5$) achieves latency and peak memory comparable to the single-group

Algorithm 1 Outlier-Aware Attention Linearization (OAL)**Require:** Calibration data and pretrained attention states; evaluation inputs Q, K, V ; $\epsilon > 0$

- 1: **Offline calibration (per layer and query head)**
- 2: Determine group map g from contribution statistics (Section 4.2)
- 3: Fit L by output matching with g and pretrained states fixed (Eq. (20))
- 4: **Causal evaluation (fixed g, L)**
- 5: Compute a, b_r, C_{rt} from L, g, ϵ ; initialize $H_0, H_1, H_2 = 0$
- 6: **for** $i = 1, \dots, N$ **do**
- 7: $u_i \leftarrow [v_i; 1]$; $H_0 \leftarrow H_0 + u_i$
- 8: $H_{1,r} \leftarrow H_{1,r} + k_{i,r}u_i$ for all r
- 9: $H_{2,rt} \leftarrow H_{2,rt} + k_{i,r}k_{i,t}u_i$ for all r, t
- 10: Compute $\hat{y}_i$ using Eq. (6)
- 11: **end for**
- 12: **return** $\hat{Y} = [\hat{y}_1, \dots, \hat{y}_N]^\top$

Table 2 | **Frozen-attention approximation fidelity on Qwen2.5-0.5B.** Results cover all query heads in Layers 4, 12, and 20 at sequence length 4096. Lower is better for all metrics.

Method	Before W_o	Before W_o	After W_o	After W_o
	MSE ↓	RelMSE ↓	MSE ↓	RelMSE ↓
Performer	0.103	0.569	0.038	0.710
SpectraFormer	0.106	0.585	0.041	0.772
DiJiang	0.086	0.474	0.033	0.631
TaylorShift	0.088	0.490	0.034	0.649
OAL	0.078	0.432	0.031	0.595

Learned Global baseline ($G = 1$), suggesting that the additional coefficient flexibility introduces little operator overhead in this configuration. Cross-method comparisons reflect the evaluated implementations, which differ in numerical precision and optimization level. FlashAttention remains faster and more memory-efficient throughout the tested range.

5. Experiments

We first evaluate OAL’s approximation fidelity on frozen pretrained attention states. We then examine whether OAL supports better language-modeling and downstream performance after attention replacement under a common, limited LoRA adaptation budget.

5.1. Frozen-Attention Approximation Fidelity

To isolate approximation error before model adaptation, we compare OAL and existing attention linearization methods on identical frozen queries, keys, and values from Qwen2.5-0.5B Base, with queries and keys taken after RoPE. The evaluation covers all query heads in Layers 4, 12, and 20 at sequence length $N = 4096$.

Approximation error is measured by MSE and relative MSE (RelMSE) against the corresponding softmax outputs, both before and after the output projection W_o . Exact metric definitions and aggregation procedures are provided in Appendix C.

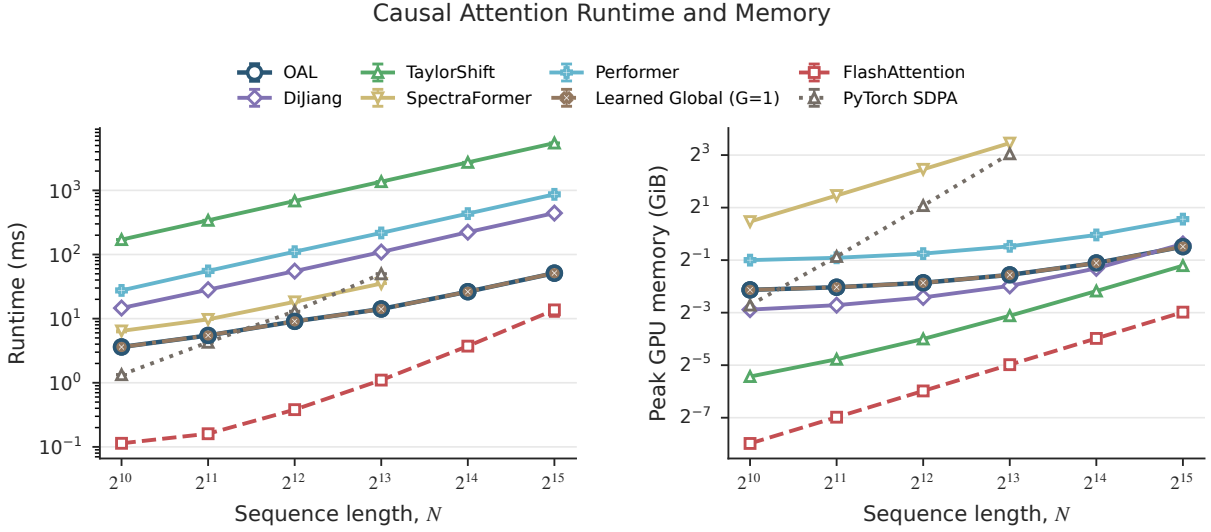


Figure 3 | **Causal attention forward runtime and memory.** Forward-only measurements use an A100-SXM4-40GB with BF16 inputs. OAL ($G = 5$) and Learned Global ($G = 1$) share the HD block-GEMM backend. Memory denotes total peak PyTorch-allocated tensor memory. SpectraFormer and SDPA (math backend) are capped at $N = 8192$ by a pre-run memory guard. Implementation and measurement details appear in Appendix E.

OAL achieves the lowest MSE and RelMSE both before and after the output projection W_o (Table 2). DiJiang is the strongest evaluated baseline under all four metrics.

5.2. End-to-End Conversion and Downstream Performance

The conversion experiments use Qwen2.5-0.5B Base and Llama-3.2-1B Base. We compare OAL with a softmax reference, a learned global quadratic, Performer, SpectraFormer, DiJiang, and TaylorShift. The learned global quadratic is the single-group ($G = 1$) instance of OAL and shares its calibration and adaptation protocol, providing a controlled comparison of global and contribution-based coefficient sharing. Following the hybrid conversion setting explored in LoLCATs (Zhang et al., 2025), we retain softmax attention in the initial and final layers. Our layer-wise activation analysis motivates converting the intervening layers, where large activations persist across depth (Appendix F). All converted methods use the same replacement range within each backbone.

All methods, including the softmax reference, use the same LoRA configurations, training data, and update budgets within each backbone. Adaptation uses next-token prediction on WikiText-103. For OAL, the calibrated kernel coefficients are jointly optimized with the LoRA parameters, while the dimension groups remain fixed. We adapt Performer, SpectraFormer, DiJiang, and TaylorShift from their official source code; modifications are detailed in Appendix F.1, and shared conversion and adaptation settings in Appendix F. Table 3 reports language-modeling and downstream results for both backbones. Evaluation protocols are detailed in Appendix F.3.

OAL achieves the lowest WikiText-103 perplexity and highest normalized accuracy on all three downstream tasks among the evaluated converted models on both backbones (Table 3). The LoRA-adapted softmax reference remains stronger on every metric, indicating a substantial conversion gap. Relative to the learned global quadratic ($G = 1$), OAL reduces perplexity from 31.71 to 27.50 on

Table 3 | Model-level quality after LoRA adaptation on two backbones. We report WikiText-103 test PPL (lower is better) and zero-shot normalized accuracy (% , higher is better). Bold highlights the leading results among converted methods within each backbone.

Method	Qwen2.5-0.5B				Llama-3.2-1B			
	PPL	PIQA	HellaSwag	ARC-E	PPL	PIQA	HellaSwag	ARC-E
Softmax (LoRA)	10.89	70.18	51.38	62.63	8.19	74.92	62.84	59.43
Performer	37.03	57.78	30.37	38.93	50.39	57.07	27.42	32.58
SpectraFormer	34.78	59.19	31.77	41.84	45.09	59.36	30.41	36.49
DiJiang	36.05	59.90	31.47	41.41	56.27	58.92	28.75	36.32
TaylorShift	39.50	59.90	31.18	40.78	62.67	57.40	28.95	35.77
Learned global	31.71	61.15	32.43	41.41	36.03	57.40	35.02	33.46
OAL	27.50	64.69	37.20	50.38	32.99	63.66	35.42	41.67

Qwen2.5-0.5B and from 36.03 to 32.99 on Llama-3.2-1B. These gains support the effectiveness of contribution-based grouping over globally shared coefficients under the evaluated protocol.

6. Related Work

Kernel Linearization and Pretrained-Model Conversion Kernel methods such as Performer, SpectraFormer, and DiJiang approximate softmax attention with feature representations that permit linear-time evaluation (Choromanski et al., 2021; Nguyen et al., 2026; Chen et al., 2024b). T2R converts pretrained Transformers through attention replacement and finetuning (Kasai et al., 2021). Hedgehog learns MLP feature maps by matching softmax attention weights (Zhang et al., 2024), while LoLCATs combines attention transfer with low-rank adaptation (Zhang et al., 2025). OAL shares the goal of preserving pretrained behavior, but organizes kernel coefficients by dimension-wise contribution, providing a structured alternative to learned MLP feature maps. Our comparison scope is detailed in Appendix F.1.

Polynomial Attention In language models trained from scratch, Based uses a second-order Taylor kernel (Arora et al., 2024), while ReBased modifies the quadratic kernel with query/key normalization and learnable affine transformations (Aksenov et al., 2024). PolySketchFormer approximates higher-degree polynomial attention through sketching (Kacham et al., 2024). OAL targets pretrained softmax recovery under limited adaptation. It retains original coordinate products and learns coefficients for linear terms and quadratic interactions, with sharing determined by contribution statistics.

Activation and Attention Outliers Massive activations and attention sinks reveal different forms of nonuniformity in pretrained Transformers: unusually large hidden activations and persistent attention to particular positions, respectively (Sun et al., 2024; Xiao et al., 2024; Li et al., 2026a,b). Systematic Outliers examines how alignment in specific query and key dimensions produces large attention weights (An et al., 2025). Building on these observations, we examine how concentrated query-key contributions relate to attention-linearization error, using this association to guide coefficient sharing.

7. Discussion

We introduce Outlier-Aware Attention Linearization (OAL), a contribution-aware method for linearizing pretrained softmax attention. Our analysis shows that approximation error is concentrated on a small subset of high-contribution query-key dimensions, motivating non-uniform parameter sharing. OAL captures this structure with a compact grouped quadratic kernel and supports exact recurrent evaluation with linear sequence-length complexity. It improves frozen-attention fidelity on Qwen2.5-0.5B and model-level performance after LoRA adaptation on both Qwen2.5-0.5B and Llama-3.2-1B. Our experiments are limited to 0.5B–1B backbones, so performance at larger scales remains unverified. Future work will evaluate OAL on larger and longer-context models and explore more compact implementations of the quadratic recurrent state.

AI Use Statement

We used generative AI tools to aid and polish writing, including language editing and improving the clarity and organization of the manuscript. We also used generative AI tools for retrieval and discovery, such as finding related work and assisting with literature search. All AI-assisted content was reviewed and verified by the authors, who take full responsibility for the final manuscript.

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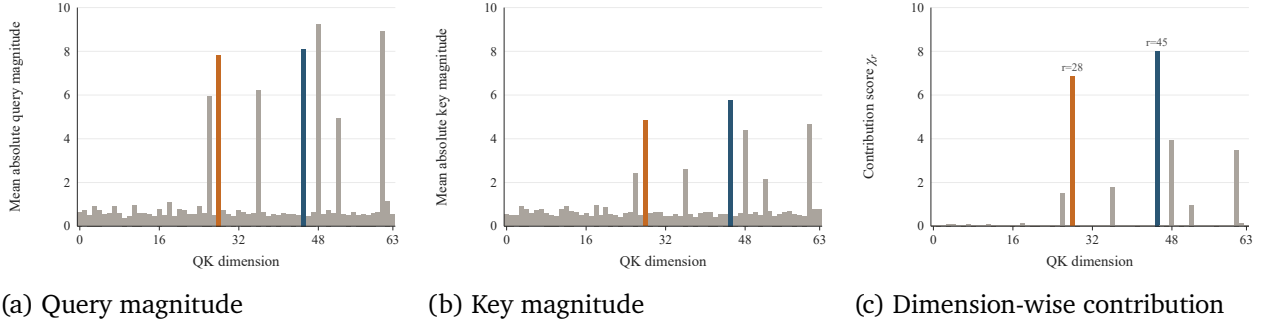


Figure 4 | **Dimension-wise heterogeneity in a representative GPT-2 Medium attention head.** GPT-2 exhibits the same qualitative relationship between nonuniform query/key representations and concentrated query–key contributions as observed in Qwen2.5-0.5B, although the dominant contribution is distributed across a broader set of dimensions.

A. Additional Analysis of Query–Key Heterogeneity

This section provides additional evidence for the dimension-wise heterogeneity discussed in Section 3.1. We first examine the same phenomenon in GPT-2 Medium and under isotropic Gaussian controls, and then evaluate the stability of the identified high-contribution structure across inputs and sequence lengths.

A.1. Cross-Model and Control Evidence

GPT-2 Medium We repeat the dimension-wise analysis on GPT-2 Medium to examine whether the heterogeneity observed in Qwen2.5-0.5B is model-specific. As shown in Figure 4, GPT-2 also exhibits nonuniform query/key magnitudes and a strongly concentrated contribution profile, although the dominant contribution is distributed across a broader set of dimensions.

Gaussian controls The Gaussian-control column in Table 1 uses a synthetic counterpart of the pretrained representations. Specifically, we replace the pretrained post-RoPE query and key representations with i.i.d. standard Gaussian samples drawn from $\mathcal{N}(0, 1)$, while keeping the remaining experimental protocol unchanged. We compute the same contribution statistics, apply the same grouping procedure to identify the corresponding high-contribution dimensions, and use the same approximation, intervention, and aggregation procedures as in the pretrained experiment. Thus, the Gaussian control changes only the distribution of the query and key representations.

The statistical significance reported in Section 3.2 is based on a separate matched-control randomization test on the original pretrained representations, described in Appendix A.2; it is not computed from the Gaussian controls.

As an additional diagnostic, for the representative Qwen2.5-0.5B attention head analyzed in Figure 5, we compare the top-5 contribution concentration of the pretrained representations with isotropic Gaussian controls of the same dimensionality. The Gaussian controls use the same head dimension and the same contribution statistic χ_r , but replace the pretrained query–key representations with i.i.d. standard Gaussian samples drawn from $\mathcal{N}(0, 1)$.

The Gaussian controls indicate that the observed concentration is not explained solely by dimensionality or by selecting particular coordinate indices.

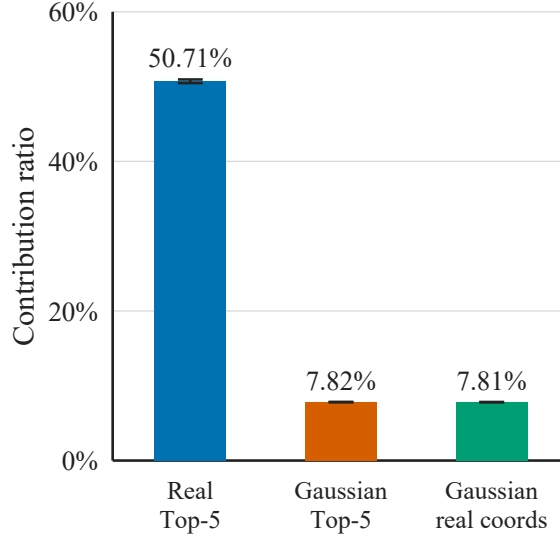


Figure 5 | **Top-5 contribution concentration.** The pretrained Qwen2.5-0.5B representation exhibits substantially stronger concentration than either isotropic Gaussian control.

A.2. High-Contribution-Dimension Intervention

We further specify the intervention used in Table 1. For each evaluated query head h , let $\mathcal{G}_{h,1}$ denote the bulk group produced by the fixed OAL grouping procedure, and let G_h denote the total number of groups. We define the high-contribution subset as

$$\mathcal{S}_h = \bigcup_{g=2}^{G_h} \mathcal{G}_{h,g},$$

that is, the union of all non-bulk groups for that head. The subset is determined independently for each layer–query-head pair using only the calibration contribution statistics.

Across the 42 evaluated query heads, the selected subsets contain 147 dimensions in total. Since $d_h = 64$, their aggregate fraction is

$$\frac{\sum_h |\mathcal{S}_h|}{42d_h} = \frac{147}{42 \times 64} \approx 5.47\%,$$

which gives the approximately 5.5% reported in Table 1.

Let $\bar{\mathcal{S}}_h$ denote the complement of $\mathcal{S}_h$. For a causal query–key pair (i, j) , the exact softmax kernel factorizes as

$$\kappa_{ij}^{\text{exact}} = \exp\left(\frac{q_i^\top k_j}{\sqrt{d_h}}\right) = \kappa_{ij, \mathcal{S}_h}^{\text{exact}} \kappa_{ij, \bar{\mathcal{S}}_h}^{\text{exact}}, \quad (8)$$

where

$$\kappa_{ij, \bar{\mathcal{S}}_h}^{\text{exact}} = \exp\left(\frac{\sum_{r \notin \mathcal{S}_h} q_{i,r} k_{j,r}}{\sqrt{d_h}}\right). \quad (9)$$

To measure the error associated with approximating only $\mathcal{S}_h$, we mask the query and key coordinates outside $\mathcal{S}_h$ and apply the corresponding baseline approximation to the masked vectors, retaining the kernel’s input preprocessing. The approximate factor for the selected dimensions and the exact factor

for their complement are computed separately. Denoting the resulting sub-kernel by $\widehat{\kappa}_{ij, S_h}^{(m)}$ for method m , we form the full-head hybrid kernel

$$\widetilde{\kappa}_{ij}^{(m, S_h)} = \widehat{\kappa}_{ij, S_h}^{(m)} \kappa_{ij, \bar{S}_h}^{\text{exact}}. \quad (10)$$

Thus, dimensions outside S_h remain exact at the unnormalized-kernel level. The hybrid kernel is then causally masked, normalized over the key positions, and applied to the pretrained values:

$$\widetilde{y}_i^{(m, S_h)} = \frac{\sum_{j \leq i} \widetilde{\kappa}_{ij}^{(m, S_h)} v_j}{\sum_{j \leq i} \widetilde{\kappa}_{ij}^{(m, S_h)}}. \quad (11)$$

The tilde distinguishes this hybrid output from a full-head approximation. The same intervention is used for all evaluated approximation methods; the methods differ only in the construction of $\widehat{\kappa}_{ij, S_h}^{(m)}$.

Error metric The intervention is evaluated at the attention-output level rather than on kernel values or attention probabilities. Let Y^{softmax} denote the exact causal softmax attention output. For an evaluated input and layer, we define

$$\mathcal{E}_S^{(m)} = \text{MSE}\left(\widetilde{Y}^{(m, S_h)}, Y^{\text{softmax}}\right), \quad (12)$$

$$\mathcal{E}_{\text{full}}^{(m)} = \text{MSE}\left(\widehat{Y}^{(m, \text{full})}, Y^{\text{softmax}}\right), \quad (13)$$

where the MSE is averaged over the evaluated query heads, query positions, and output-head dimensions. We report the retained-error ratio

$$R_S^{(m)} = \frac{\mathcal{E}_S^{(m)}}{\mathcal{E}_{\text{full}}^{(m)}}. \quad (14)$$

A larger value therefore indicates that restricting the approximation to the selected dimensions preserves a larger fraction of the full-method attention-output error.

Evaluation protocol We evaluate Qwen2.5-0.5B on zero-based Layers 4, 12, and 20, using all query heads in each layer. Evaluation uses 16 disjoint 4096-token WikiText-2 sequences, with 32 causal query positions sampled per sequence and layer. For stochastic approximation methods, we use five random seeds, $\{0, 1, 2, 3, 4\}$. The selected-subset and full-method MSEs are first averaged over these seeds before forming the retained-error ratio. TaylorShift is deterministic and is evaluated once. The resulting ratios are then macro-averaged across the evaluated layers and input sequences.

Matched-control randomization test To assess whether the retained-error ratio of the high-contribution subset can be explained by selecting an equally sized subset of ordinary dimensions, we construct $B = 1000$ matched control subsets. For each control, the same intervention and aggregation procedure is used to obtain a retained-error ratio R_b^{control} . We perform a one-sided Monte Carlo randomization test and report

$$p = \frac{1 + \sum_{b=1}^B \mathbf{1}[R_b^{\text{control}} \geq R_S]}{B + 1}. \quad (15)$$

The +1 correction avoids reporting a zero Monte Carlo p -value; with $B = 1000$, the minimum reportable value is $1/1001 \approx 0.000999$.

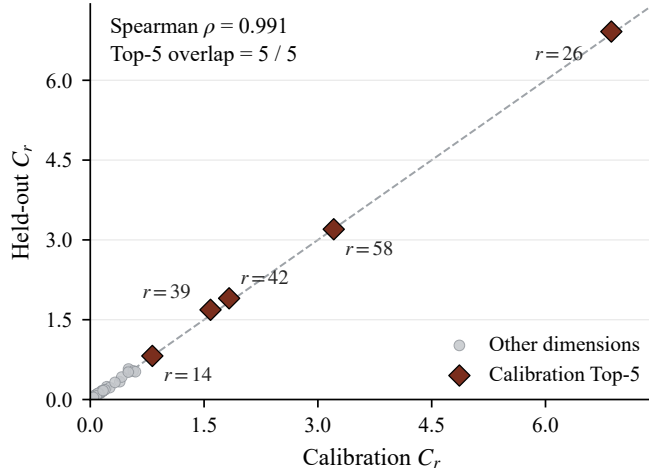


Figure 6 | **Stability of dimension-wise contributions across inputs.** Contribution scores computed from calibration and held-out inputs achieve Spearman correlation $\rho = 0.991$, with a top-5 overlap of 5/5.

Table 4 | **Stability across sequence lengths.** Normalized contribution $\chi_r / \sum_t \chi_t$ (%) of the five highest-contribution dimensions identified at sequence length 4096 for the representative attention head.

Length	$r = 26$	$r = 58$	$r = 19$	$r = 39$	$r = 7$
128	33.34	13.64	2.12	3.30	3.56
512	32.96	15.70	1.27	3.66	3.86
1024	32.19	15.91	1.27	3.78	3.86
2048	30.93	16.18	2.10	3.75	3.80
4096	28.75	17.53	4.11	3.71	3.66

A.3. Stability of Dimension-Wise Contributions

Across inputs We first examine whether the contribution profile depends strongly on the calibration inputs. Figure 6 compares the dimension-wise contribution scores computed from calibration and held-out inputs for the representative Qwen2.5-0.5B attention head. The two profiles are highly consistent, with Spearman correlation $\rho = 0.991$ and identical top-5 dimensions.

Across sequence lengths We further evaluate the same representative head using prefixes of a fixed sequence at different lengths. Table 4 reports the normalized contributions of the five dimensions identified at sequence length 4096.

The two dominant dimensions, $r = 26$ and $r = 58$, remain the largest contributors across all evaluated sequence lengths and jointly account for approximately 46–49% of the total contribution. Together with the held-out-input analysis, these results indicate that the dominant dimension-wise structure is stable under changes in both input content and sequence length.

B. Grouping and Coefficient-Calibration Details

Dimension grouping For a fixed layer and query head, we compute the dimension-wise contribution scores χ_r from the calibration set as defined in Eq. (2). The queries and keys are the post-RoPE representations used by the attention module, causal query-key pairs are used throughout, and query

heads follow the model’s native GQA mapping to key–value heads. Different prompts are aggregated independently and then averaged with equal weight.

We sort the d_h dimensions by increasing contribution,

$$\chi_{\pi_1} \leq \chi_{\pi_2} \leq \dots \leq \chi_{\pi_{d_h}},$$

using a stable ordering, and search for group boundaries only in the high-contribution tail. Given a tail fraction α , let

$$m = \lceil \alpha d_h \rceil.$$

Candidate cuts are restricted to

$$k \in \{\max(1, d_h - m), \dots, d_h - 1\}.$$

A cut at position k separates

$$\mathcal{L}_k = \{\pi_1, \dots, \pi_k\}, \quad \mathcal{U}_k = \{\pi_{k+1}, \dots, \pi_{d_h}\}.$$

Thus, the cut immediately preceding the top- m tail is included, while the endpoint $k = d_h$ is excluded so that both sides of every candidate cut are nonempty.

For each candidate cut, we first measure the adjacent contribution gap using

$$\rho_k = \frac{\chi_{\pi_{k+1}}}{\max(\chi_{\pi_k}, \varepsilon)}, \quad (16)$$

where ε is a small numerical constant used only to avoid division by zero. A large ρ_k indicates a sharp increase in contribution across the candidate boundary.

To test whether this separation is stable with respect to the calibration inputs, we perform B prompt-level bootstrap replicates. For replicate b , let

$$I_b = (i_1^{(b)}, \dots, i_P^{(b)})$$

denote the length- P sequence of prompt indices sampled independently with replacement from $\{1, \dots, P\}$. Repeated indices are retained with their sampling multiplicities. We then recompute the contribution scores as

$$\chi_r^{(b)} = \frac{1}{P} \sum_{t=1}^P \chi_{i_t^{(b)}, r}.$$

Importantly, the dimension ordering π obtained from the full calibration set is kept fixed across bootstrap replicates. For candidate cut k , we measure the separation between the two sides as

$$R_k^{(b)} = \frac{\min_{r \in \mathcal{U}_k} \chi_r^{(b)}}{\max(\max_{r \in \mathcal{L}_k} \chi_r^{(b)}, \varepsilon)}. \quad (17)$$

Its bootstrap support is then

$$S_k = \frac{1}{B} \sum_{b=1}^B \mathbf{1} \left[R_k^{(b)} \geq \tau \right], \quad (18)$$

which measures the fraction of bootstrap samples for which the two sides remain separated by at least a factor of τ .

With minimum bootstrap support γ , a candidate boundary is retained as a *natural cut* when

$$\rho_k \geq \tau \quad \text{and} \quad S_k \geq \gamma.$$

All natural cuts satisfying these conditions are retained; we do not apply an additional minimum group size, boundary-spacing constraint, or maximum group-count cap.

To prevent the grouped parameterization from collapsing to the global single-group case, we require at least $G_{\min}$ groups per head. If the natural cuts yield fewer than $G_{\min}$ groups, unused candidate cuts are added in descending lexicographic order of (S_k, ρ_k) until the minimum group count is reached. Bootstrap support is therefore prioritized over the full-sample gap magnitude; exact ties retain the original candidate-cut order. These *fallback cuts* are restricted to the same high-contribution tail but need not satisfy the natural-cut thresholds.

Let the final retained cuts be $k_1 < \dots < k_{G-1}$, with $k_0 = 0$ and $k_G = d_h$. The resulting groups are

$$\mathcal{G}_g = \{\pi_{k_{g-1}+1}, \dots, \pi_{k_g}\}, \quad g = 1, \dots, G. \quad (19)$$

Thus, the first group typically forms a large low-contribution shared bulk, while subsequent groups refine the high-contribution tail. Grouping is performed independently for every layer and query head, and the resulting assignments are mapped back to the original dimension indices. The groups are determined before kernel-coefficient fitting and remain fixed during subsequent calibration, model adaptation, and evaluation.

Grouping settings Both backbones use the same fixed set of 40 short calibration prompts and $B = 1000$ prompt-level bootstrap replicates. For the formal experiments, Qwen2.5-0.5B uses $(\alpha, \gamma) = (0.25, 0.9)$, while Llama-3.2-1B uses $(\alpha, \gamma) = (0.50, 0.5)$. Both use $\tau = 1.5$ and $G_{\min} = 2$, with no explicit maximum group-count constraint.

Coefficient calibration For a fixed layer and query head, let C be the coefficient-fitting sequences and $\hat{y}_i(L; x)$ the normalized causal OAL output for $x \in C$. With groups and pretrained queries, keys, and values fixed, we fit the lower-triangular L to the corresponding softmax outputs:

$$L^* = \arg \min_L \frac{1}{|C|} \sum_{x \in C} \sum_i \left\| \hat{y}_i(L; x) - y_i^{\text{softmax}}(x) \right\|_2^2. \quad (20)$$

Matching the normalized output directly accounts for both the kernel weights and their normalization denominator. The calibrated factors L^* initialize subsequent model-level adaptation, during which the dimension groups remain fixed while L is jointly optimized with the LoRA parameters under the causal language-modeling objective.

Coefficient-calibration settings Coefficient fitting uses WikiText-2 raw v1 training text, concatenated and divided into non-overlapping sequences of length 4096. We use blocks 16–23 as eight training sequences and blocks 24–25 as two validation sequences. This coefficient-fitting data is separate from the short-prompt calibration data used to determine the dimension groups.

We optimize the OAL factors with Adam at learning rate 0.02, without a learning-rate scheduler, for up to 1000 updates per initialization. Parameters are optimized in FP32 with gradient norm clipped at 10. Each update samples one training sequence and 32 causal query positions. Validation is performed every 100 updates, with early stopping after 400 updates without improvement.

We use three initialization candidates for OAL: an embedding of the learned global quadratic into the grouped parameterization, a second-order Taylor initialization, and a Gaussian-perturbed learned-global initialization. Within each initialization, the lowest-validation-MSE checkpoint is retained independently for each query head; the final coefficients are then selected per head across the three initialization candidates.

Layers are calibrated separately, with independent coefficients for each query head. The learned global quadratic uses the same training and validation tensors, output-matching objective, optimizer, learning rate, query-position sampling protocol, and checkpoint selection rule as OAL. Both methods use three initialization candidates with up to 1000 updates per initialization; they differ only in their kernel parameterization and corresponding initializations.

Frozen-attention evaluation uses blocks 0–15 of the same token stream, giving 16 additional sequences of length 4096. These blocks are disjoint from both coefficient training and validation and are not used for coefficient optimization or checkpoint selection.

Group-budget diagnostic Figure 2(b) uses the same representative Qwen2.5-0.5B query head (zero-based Layer 4, Head 0) for all group budgets. Dimensions are sorted by the same calibration contribution scores χ_r , and for a prescribed budget G we select the $G - 1$ largest adjacent gaps in the sorted profile as group boundaries. Thus, increasing G only further subdivides the same contribution ordering.

All budgets use the same coefficient-fitting and evaluation data: eight 4096-token WikiText-2 blocks for training, two for validation, and 16 disjoint blocks for held-out evaluation. Coefficients are fitted with the same Adam-based optimization sweep and selected by validation attention-output MSE; held-out data are never used for model selection. The plotted error is pooled held-out attention-output MSE over all tokens and output dimensions. Therefore, the points in Figure 2(b) are directly comparable across group budgets, with the group partition and corresponding parameter count as the only systematic changes.

C. Frozen-Attention Error Metrics

For each held-out sequence, let Y_b denote the exact softmax attention output and $\widehat{Y}_b$ the corresponding approximation, with all evaluated layers and heads collected into the same tensor. We report

$$\text{MSE} = \frac{1}{B} \sum_{b=1}^B \frac{\|\widehat{Y}_b - Y_b\|_F^2}{n_b}, \quad (21)$$

and

$$\text{RelMSE} = \frac{1}{B} \sum_{b=1}^B \frac{\|\widehat{Y}_b - Y_b\|_F^2}{\|Y_b\|_F^2}, \quad (22)$$

where B is the number of held-out sequences and n_b is the number of scalar output elements in sequence b . Thus, RelMSE normalizes the squared approximation error by the energy of the exact softmax output for each sequence before averaging across sequences.

For the metrics before W_o , Y_b and $\widehat{Y}_b$ contain the per-head attention outputs before the output projection. For the post- W_o metrics, the heads are first concatenated in their original order and passed through the frozen output projection W_o ; the same MSE and RelMSE definitions are then applied to the projected outputs. The post- W_o RelMSE is therefore normalized by the energy of the exact softmax output after W_o , rather than by the pre-projection attention output.

D. Global Reference and HD Realization

Positive global reference The learned global reference is a quadratic in the total score $s = q^\top k / \sqrt{d_h}$. A symmetric positive semidefinite matrix $M_{\text{global}} = [m_{uv}]_{u,v=0}^1 \in \mathbb{R}^{2 \times 2}$ parameterizes it as

$$\begin{aligned} \kappa_{\text{global}}(q, k) &= \epsilon + \begin{bmatrix} 1 & s \end{bmatrix} M_{\text{global}} \begin{bmatrix} 1 & s \end{bmatrix}^\top \\ &= (\epsilon + m_{00}) + 2m_{01}s + m_{11}s^2, \quad M_{\text{global}} \succeq 0. \end{aligned} \quad (23)$$

Thus this is the global polynomial $a + bs + cs^2$ with $a = \epsilon + m_{00}$, $b = 2m_{01}$, and $c = m_{11}$; the matrix constraint and $\epsilon > 0$ ensure strictly positive weights. It is contained in OAL: setting $M_{00} = m_{00}$, $M_{0g} = M_{g0} = m_{01}$, and $M_{gh} = m_{11}$ for all groups recovers the same polynomial because $s = \sum_g s_g$. This grouped matrix is positive semidefinite, since its quadratic form on any vector t equals the quadratic form of M_{global} on $[t_0, \sum_g t_g]^\top$, and hence admits $M = LL^\top$.

Evaluation equivalence For a single query, $(q_i^\top k_j)^2 = q_i^\top k_j k_j^\top q_i$ gives

$$\sum_j (q_i^\top k_j)^2 V_{j,\ell} = q_i^\top \left(\sum_j V_{j,\ell} k_j k_j^\top \right) q_i = q_i^\top K^\top D_\ell K q_i.$$

The last expression is the i th diagonal entry of $QK^\top D_\ell KQ^\top$, proving Eq. (4). Multiplication by D_ℓ amounts to scaling key rows and does not require storing a diagonal matrix. For the grouped kernel, expanding $M = LL^\top$ gives

$$\begin{aligned} \kappa_{\text{OAL}}(q, k) &= \epsilon + M_{00} + 2 \sum_g M_{0g} s_g + \sum_{g,h} M_{gh} s_g s_h \\ &= a + \frac{1}{\sqrt{d_h}} \sum_r b_r q_r k_r + \frac{1}{d_h} \sum_{r,t} C_{rt} q_r q_t k_r k_t, \end{aligned} \quad (24)$$

which yields the coefficient mapping in Section 4.3. Substituting Eq. (5) into Eq. (6) and exchanging sums then gives

$$\begin{aligned} \widetilde{u}_i &= \sum_{j \leq i} \left(a + \frac{1}{\sqrt{d_h}} \sum_r b_r q_{i,r} k_{j,r} + \frac{1}{d_h} \sum_{r,t} C_{rt} q_{i,r} q_{i,t} k_{j,r} k_{j,t} \right) u_j \\ &= \sum_{j \leq i} \kappa_{\text{OAL}}(q_i, k_j) u_j. \end{aligned} \quad (25)$$

The same substitution establishes the identity for κ_{dim} with general coefficients. The argument uses separable query and key factors, not Mercer positive semidefiniteness; pointwise positivity of OAL's weights does not imply the latter property.

Exact state and parameter counts The ordered-pair states in Eq. (5) contain $(1 + d_h + d_h^2)(d_v + 1)$ entries. Since $H_{2,i,rt} = H_{2,i,tr}$ and C is symmetric, the packed-symmetric implementation stores only $(1 + d_h + d_h(d_h + 1)/2)(d_v + 1)$ entries. Diagonal pairs have multiplicity one and off-diagonal pairs

multiplicity two during readout, preserving all interactions. These counts describe one logical state, not total workspace or peak memory; GQA can share KV states across query heads. The lower-triangular factor L has $(G + 1)(G + 2)/2$ learned parameters per head.

E. Runtime and Memory Benchmark

Figure 3 evaluates forward-only causal attention on an NVIDIA A100-SXM4-40GB. The batch size is $B = 1$, with $H_q = 14$ query heads, $H_{kv} = 2$ key/value heads, and $d_h = d_v = 64$. Input tensors use BF16, and sequence lengths range from 2^{10} to 2^{15} . These are attention-operator measurements, not end-to-end model latency or per-token autoregressive decoding latency.

OAL ($G = 5$) and Learned Global ($G = 1$) use the same optimized HD block-GEMM backend, with 128-token blocks, BF16 Tensor Core contractions, and FP32 working coefficients. The implementation materializes intermediate tensors and reuses structural preparation across timed calls. OAL uses a fixed five-group configuration, matching the largest group count observed across the heads used in our model-adaptation experiments, to measure operator scaling. This benchmark does not execute the heterogeneous group configurations of all calibrated heads in a converted model.

Performer, SpectraFormer, DiJiang, and TaylorShift use causal adaptations of their official source code. The source repositories and kernel adaptations are described in Appendix F.1. These methods retain their respective causal execution paths, while OAL and Learned Global share the specialized HD backend. The implementations differ in optimization level; in particular, the TaylorShift adaptation receives no additional kernel optimization.

Internal arithmetic also differs across implementations. Performer and SpectraFormer use FP64 feature computation and prefix accumulation. DiJiang uses FP32 feature computation and prefix accumulation with 128-token blocks, explicitly repeating shared keys and values across query heads. Together with the backend differences above, these precision choices affect the measured runtime and memory, so cross-method gaps cannot be attributed solely to the kernel formulations.

Both softmax references use PyTorch scaled dot-product attention (SDPA). The FlashAttention curve selects the FLASH_ATTENTION backend, whereas the PyTorch SDPA curve selects the MATH backend. Both calls enable grouped-query attention and causal masking, with zero attention dropout. Measurements run with gradient recording disabled, after warmup, and runtime is wall-clock latency with CUDA synchronization. Runtime markers report the mean across trials.

GPU memory is the total peak allocated tensor memory reported by `torch.cuda.max_memory_allocated()`, including resident inputs and method state. It is neither incremental memory above the resident allocation nor allocator-reserved memory. This peak includes workspace and therefore differs from the logical recurrent-state size analyzed in Section 4.4. For both SpectraFormer and math-backend SDPA, the $N = 16384$ case is skipped by a pre-run memory guard because the estimated next peak exceeds 80% of the current free GPU memory. Both sweeps therefore end at $N = 8192$; these omissions reflect the benchmark’s memory budget rather than observed out-of-memory failures.

F. Conversion and LoRA Setup

Attention replacement Figure 7 shows sustained large activations in the intermediate layers of both models, with different concentration patterns across hidden dimensions. Because pronounced activation outliers persist through these intermediate layers, we use this region as the primary target for attention replacement and determine the replacement ranges accordingly.

For Qwen2.5-0.5B, we replace zero-based layers 3–20 (inclusive), retaining softmax attention in the

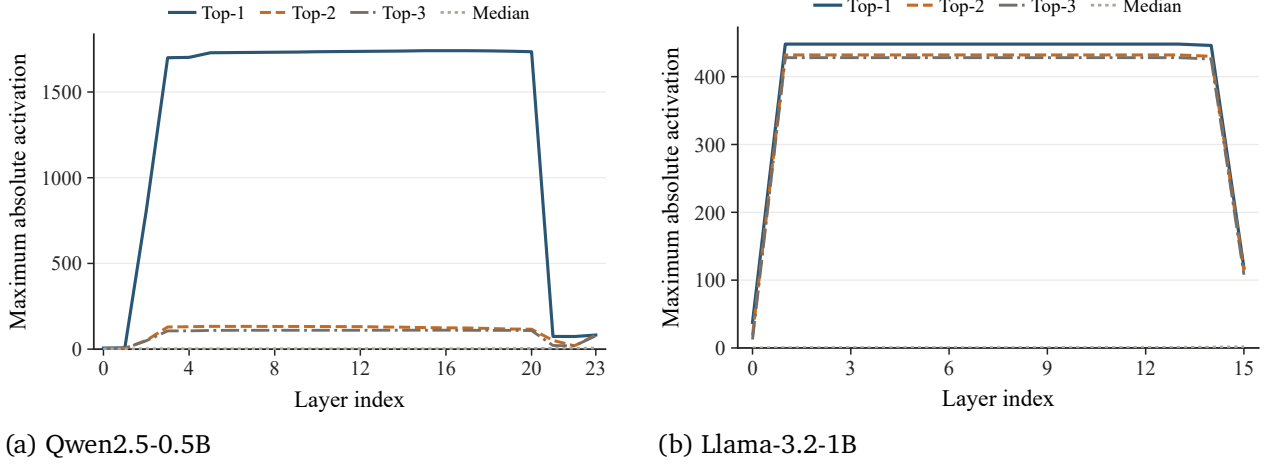


Figure 7 | **Layer-wise activation profiles.** We measure the output residual stream of each decoder block on the same 40 short prompts. For each hidden dimension, we record its maximum absolute activation over all tokens and prompts; Top-1/2/3 denote the three largest dimension-wise maxima, and Median denotes their median across dimensions. Layer indices are zero-based.

first and last three layers. For Llama-3.2-1B, we replace layers 1–14 (inclusive), retaining softmax attention in the first and last layers. The same replacement range is used for all converted methods within each backbone. The softmax reference retains softmax attention throughout. These are hybrid models: the retained softmax layers still contribute quadratic prefill computation, even though the OAL operator is linear in sequence length.

F.1. Baseline Implementations and Adaptations

The baseline implementations in Table 3 are adapted from the official source code of Performer, SpectraFormer, DiJiang, and TaylorShift. We integrate their attention kernels into the same pretrained backbones, retaining the native query, key, value, and output projections, RoPE, and GQA head mapping. Each kernel uses causal prefix aggregation and normalization over the available keys. This setup compares attention approximations under a common conversion protocol rather than reproducing each method’s original full-model architecture and training recipe.

For the frozen-attention evaluation, trainable kernel parameters are first calibrated on a small calibration set while the pretrained Q/K/V states remain fixed. This calibration is separate from the subsequent LoRA runs, where the trainable kernel parameters are optimized jointly with LoRA from the start.

Performer We adapt the official FAVOR+ implementation¹ using positive orthogonal random features with 250 features per query head. Gaussian orthogonal projections are sampled independently for each layer and query head and remain fixed throughout adaptation; only the LoRA parameters are trained. Queries and keys are scaled by $d_h^{-1/4}$ before feature evaluation. To stabilize exponential features, query log-features are shifted by their per-token maximum, while key log-features use a shared maximum over tokens and features within each head. These common factors cancel in normalized attention. Feature computation and causal prefix accumulation use FP64, with the output cast back to the model dtype.

¹<https://github.com/google-research/google-research/tree/master/performer>

SpectraFormer We adapt the OPRF-FastFoodL variant from the official repository², using 65 positive features for $d_h = 64$. SpectraFormer supports multiple combinations of feature functions and projection constructions. We select OPRF-FastFoodL as a lightweight learnable representative: its structured projection requires only $3d_h$ trainable kernel parameters per layer. Each replacement layer has three trainable FastFood vectors $B, G, S \in \mathbb{R}^{64}$ shared across its query heads, with a fixed permutation. This gives 192 kernel parameters per layer, or 3,456 for Qwen2.5-0.5B and 2,688 for Llama-3.2-1B over their respective replacement layers, excluding LoRA. The vectors remain trainable but are initialized from the standard FastFood distributions used in the upstream fixed branch. This avoids the nearly token-independent features observed with the small-normal initialization of the learnable branch. We retain the upstream feature scaling and per-token row-maximum stabilization followed by an additive 10^{-4} feature offset. The attention aggregation is adapted to causal prefixes, using FP64 feature and accumulation arithmetic and FP32 trainable parameters.

DiJiang We adapt the DCT projection and feature-wise softmax construction from the official repository³. The fixed projection combines an orthonormal DCT with diagonal inverse standard-normal samples. After scaling queries and keys by $d_h^{-1/4}$, we apply the projection and a softmax over the 64 feature coordinates. We add positive feature weights $D = \exp(\theta)$ to both query and key features, with a separate 64-dimensional θ for each layer and query head, initialized to zero and trained jointly with LoRA. We normalize the causal weighted-value sum by the corresponding sum of kernel weights. The released model’s positional decay factors, GroupNorm, and output gate are omitted to retain the pretrained attention interface. Features and prefix states use FP32. Thus, the DiJiang row reports this normalized causal adaptation of the released feature map.

TaylorShift We adapt the official implementation⁴ by retaining its query/key ℓ_2 normalization, fixed polynomial $1 + s + \frac{1}{2}s^2$, and learnable per-head temperature. Here, $s = \tau \hat{q}^\top \hat{k}$, where $\hat{q}$ and $\hat{k}$ are unit-normalized vectors. Each layer and query head has its own temperature τ , initialized to one and trained jointly with LoRA. We replace noncausal aggregation with causal prefix evaluation. We disable the optional output-amplitude normalization (output_normalized=False). Its full-sequence factor $\sqrt{N/d_h}$ would make a fixed prefix’s output depend on the total input length, while replacing N with the current prefix length would introduce an additional position-dependent scaling rule. We therefore use no extra output scaling and retain the usual attention numerator/denominator normalization.

Learned global quadratic This controlled reference uses the positive global parameterization in Appendix D. It shares OAL’s output-calibration protocol, and its calibrated kernel parameters are jointly optimized with LoRA. It isolates the effect of global versus contribution-based coefficient sharing within the quadratic family.

Scope of the comparison Our comparison focuses on fixed feature maps and kernels with a small number of trainable parameters, excluding the shared LoRA adapters. Hedgehog instead learns dense MLP feature maps (Zhang et al., 2024). An affine map from d_h input coordinates to m output coordinates contains $d_h m + m$ parameters; for $d_h = m = 64$, this is 4,160 parameters per map, before accounting for replication across heads and layers. This differs from the fixed projections, trainable vectors, and scalar or grouped coefficients used in our evaluated kernels. Its original attention-weight

²<https://github.com/dukenguyenxyz/spectraformer>

³<https://github.com/YuchuanTian/DiJiang>

⁴<https://github.com/tobna/TaylorShift>

distillation objective also compares query–key pairs, with quadratic scaling in sequence length when evaluated over all causal pairs; this training cost does not imply quadratic inference complexity. We do not evaluate Hedgehog under the present protocol, and our results therefore do not establish OAL’s relative performance against MLP-based feature maps. Extending the comparison would require specifying their parameter budget and adaptation procedure, including whether to retain attention-weight distillation or use output matching.

F.2. LoRA Adaptation

For all methods, LoRA is applied to the query, key, value, and output projections in every layer, including layers that retain softmax attention. We use rank 8, scaling parameter $\alpha = 16$, and dropout 0.05, keeping the pretrained backbone weights frozen. Each method is adapted for 512 optimizer updates on WikiText-103 with sequence length 4096, batch size 1, and no gradient accumulation, corresponding to 2,097,152 input tokens. All runs use a single training seed of 42 and the same training-block order within each backbone. OAL starts from training-side calibrated groups and kernel coefficients; its coefficients are optimized jointly with LoRA while the groups remain fixed.

We use AdamW for LoRA and trainable kernel parameters, with learning rate 5×10^{-4} , weight decay 0.01, $(\beta_1, \beta_2) = (0.9, 0.999)$, and optimizer $\epsilon = 10^{-8}$. The learning rate follows a linear warmup over the first 5% of updates, followed by cosine decay. The gradient norm is clipped to 1.0. Every 128 updates, we compute next-token NLL on the same fixed 32 WikiText validation blocks and select the checkpoint with the lowest validation NLL. Final language-modeling and downstream evaluations use this selected checkpoint; test results are not used for checkpoint selection.

F.3. Evaluation Protocol

Within each backbone, all methods use the same evaluation protocol and tokenizer, without chat templates or added BOS/EOS tokens.

Perplexity We evaluate Salesforce/wikitext (wikitext-103-raw-v1, test split). Text entries are joined in source order with two newlines, tokenized, and divided into independent, non-overlapping 4096-token blocks, discarding the incomplete final block. PPL is the exponential of the mean negative log-likelihood over all scored tokens, with 4095 next-token targets per block.

Downstream tasks We report zero-shot accuracy on PIQA validation (lighteval/piqa, 1,838 examples), HellaSwag validation (Rowan/hellaswag, 10,042 examples), and ARC-Easy test (allenai/ai2_arc, 2,376 examples). PIQA and ARC-E use Question: {question}, followed by a newline and Answer:. HellaSwag uses the activity label and context, with the second context segment capitalized, title markers replaced by periods, and bracketed artifacts removed. Candidates are prefixed with a space and ranked by their conditional log-likelihood divided by candidate token count for PIQA/HellaSwag or option-text character count for ARC-E (excluding the added space). Only candidate tokens are scored.