

A. MNIST-like Experiments

The plots presented in this appendix support the conclusions made in the main text and provide an overview of the experiments conducted on MNIST-like datasets.

Each figure shows the compression-accuracy trade-off of a particular method and input features for *SimpleConvModel* and *TwoLayerDenseModel* models for all four of the studied datasets (described in the main text): EMNIST-Letters on the *top-left*, KMNIST – *top-right*, Fashion MNIST – *bottom-left*, and MNIST on the *bottom-right*. Figures 7, 8, 9, and 10 present \mathbb{R} and \mathbb{C} models with *the same intermediate feature sizes*.

We compare \mathbb{R} networks against $\frac{1}{2}\mathbb{C}$ with half the number of parameters for raw input features on figures 13, and 14, and $2\mathbb{R}$ with double the number of parameters against \mathbb{C} for Fourier input features on figures 11 and 12.

B. Complex-valued Local Reparameterization

In this section we show (11).

By $e_i \in \mathbb{R} \hookrightarrow \mathbb{C}$ we denote the i -th unit vector of dimensionality *conforming* to the matrix-vector expression it is used in, $[M]$ denotes *row-major* flattening of a matrix M into a vector, i.e. in lexicographic order of its indices. Furthermore $\text{diag}(\cdot)$ embeds vectors into matrices with zeros everywhere except the diagonal, and \otimes is the Kronecker product, for which we note the following identities $[PQR] = (P \otimes R^\top)[Q]$, $(P \otimes Q)^\top = (P^\top \otimes Q^\top)$, and $(P \otimes R)(C \otimes S) = PQ \otimes RS$ (Petersen and Pedersen, 2012).

If we assume a factorized \mathbb{C} -Gaussian approximation (10) for $W \in \mathbb{C}^{n \times m}$, then $[W]$ is \mathbb{C} -Gaussian vector with

$$[W] \sim \mathcal{CN}_{nm}([\mu], \text{diag}[\Sigma], \text{diag}[C]), \quad (17)$$

where with $C_{ij} = \Sigma_{ij}\xi_{ij}$, $\Sigma_{ij} \geq 0$, and $|C_{ij}|^2 \leq \Sigma_{ij}$. Then for any $x \in \mathbb{C}^m$ and $b \in \mathbb{C}^n$ we have $y = Wx + b = (I_n \otimes x^\top)[W] + b$, whence the covariance and relation matrices of y are

$$\begin{aligned} (I_n \otimes x^\top)\text{diag}[\Sigma](I_n \otimes x^\top)^\top &= \sum_{ij} (I_n \otimes x^\top) \left((e_i \otimes e_j) \Sigma_{ij} (e_i \otimes e_j)^\top \right) (I_n \otimes x^\top)^\top \\ &= \sum_{ij} (e_i \otimes x^\top e_j) \Sigma_{ij} \overline{(e_i \otimes x^\top e_j)^\top} \\ &= \sum_{i=1}^n (e_i e_i^\top) \left\{ \sum_{j=1}^m \Sigma_{ij} |x_j|^2 \right\}, \end{aligned} \quad (18)$$

$$\begin{aligned} (I_n \otimes x^\top)\text{diag}[C](I_n \otimes x^\top)^\top &= \sum_{ij} (I_n \otimes x^\top) \left((e_i \otimes e_j) C_{ij} (e_i \otimes e_j)^\top \right) (I_n \otimes x^\top)^\top \\ &= \sum_{i=1}^n (e_i e_i^\top) \left\{ \sum_{j=1}^m C_{ij} x_j^2 \right\}. \end{aligned} \quad (19)$$

Since (18) and (19) are diagonal, the vector y has independent univariate \mathbb{C} -Gaussian components, whence (11) follows.

C. Backpropagation through \mathbb{C} -networks

Wirtinger (\mathbb{CR}) calculus relies on the natural identification of \mathbb{C} with \mathbb{R}^2 , and regards $f: \mathbb{C} \rightarrow \mathbb{C}$ as an algebraically equivalent function $F: \mathbb{R}^2 \rightarrow \mathbb{C}$ defined $f(z) = f(u + jv) = F(u, v)$. It enables general treatment of functions of vector \mathbb{C} -argument that possess partial derivatives with respect to real and imaginary parts, yet are not required to satisfy Cauchy-Riemann conditions. In \mathbb{CR} calculus the complex argument z and its conjugate \bar{z} act as independent variables and $f(z)$ is treated as $f(z, \bar{z})$ by way of geometric transformations $z = u + jv$ and $\bar{z} = u - jv$.

Wirtinger partial derivative operators are formally defined as $\frac{\partial}{\partial z} = \frac{1}{2} \left(\frac{\partial}{\partial u} - j \frac{\partial}{\partial v} \right)$ and $\frac{\partial}{\partial \bar{z}} = \frac{1}{2} \left(\frac{\partial}{\partial u} + j \frac{\partial}{\partial v} \right)$ and differentials are $dz = du + jdv$ and $d\bar{z} = du - jdv$. In this paradigm The usual rules of calculus, like chain and product rules, follow

directly from the definition of the operators, e.g.

$$\frac{\partial(f \circ g)}{\partial z} = \frac{\partial f(g(z))}{\partial g} \frac{\partial g(z)}{\partial z} + \frac{\partial f(g(z))}{\partial \bar{g}} \frac{\partial \bar{g}(z)}{\partial z}.$$

The total differential of f at $z = u + jv \in \mathbb{C}$ is

$$\begin{aligned} df(z) &= \frac{\partial f}{\partial z} dz + \frac{\partial f}{\partial \bar{z}} d\bar{z} \\ &= \frac{1}{2} \left(\frac{\partial F}{\partial u} du - j \cancel{\frac{\partial F}{\partial v} du} + j \cancel{\frac{\partial F}{\partial u} dv} + \frac{\partial F}{\partial v} dv \right) + \frac{1}{2} \left(\frac{\partial F}{\partial u} du + j \cancel{\frac{\partial F}{\partial v} du} - j \cancel{\frac{\partial F}{\partial u} dv} + \frac{\partial F}{\partial v} dv \right) \\ &= dF(u, v), \end{aligned}$$

At the same time the Cauchy-Riemann conditions $-j \frac{\partial F}{\partial v} = \frac{\partial F}{\partial u}$ can be expressed as $\frac{\partial f}{\partial \bar{z}} = 0$. Thus \mathbb{CR} calculus subsumes the usual \mathbb{C} -calculus of holomorphic functions, since $f(z) = f(z, \bar{z})$ is constant with respect to \bar{z} in the latter.

In optimization-related tasks the objective is $f: \mathbb{C} \rightarrow \mathbb{R}$, meaning that if it were to satisfy the Cauchy-Riemann conditions, then it necessarily should have been constant. Nevertheless, the expression of the \mathbb{CR} gradient is compatible with what is expected, when f is treated like a \mathbb{R}^2 function. For such f we have $\bar{f} = f$, which implies $\frac{\partial f}{\partial \bar{z}} = \frac{\partial \bar{f}}{\partial \bar{z}} = \frac{\partial f}{\partial z}$, whence

$$df = \frac{\partial f}{\partial z} dz + \frac{\partial f}{\partial \bar{z}} d\bar{z} = \frac{\partial f}{\partial z} dz + \overline{\frac{\partial f}{\partial z} dz} = 2\Re\left(\frac{\partial f}{\partial z} dz\right).$$

Therefore the gradient of f at z is given by $\nabla_{\bar{z}} f(z) = \overline{\frac{\partial f}{\partial z}} = \frac{\partial F}{\partial u} + j \frac{\partial F}{\partial v}$. The identification $\mathbb{C} \simeq \mathbb{R}^2$, backed by Wirtinger calculus, and emulation of \mathbb{C} -arithmetic in computational graphs with \mathbb{R} -valued operations makes it possible to reuse \mathbb{R} back-propagation and existing auto-differentiation frameworks.

D. Gradient of the KL-divergence in \mathbb{R} case

In this appendix we study the approximation proposed by Molchanov et al. (2017) for the KL divergence term (4) for \mathbb{R} Sparse Variational Dropout. Following the logic of Lapidoth and Moser (2003) we derive the expression for $\frac{d}{d \log \alpha} K(\alpha)$. Acknowledging that the same result was obtained by Hron et al. (2018, eq. (5)), we provide this appendix for the sake of completeness.

For $(z_i)_{i=1}^m \sim \mathcal{N}(0, 1)$ iid and $(\mu_i)_{i=1}^m \in \mathbb{R}$, the random variable $W = \sum_i (\mu_i + z_i)^2$ has non-central χ^2 distribution with shape m and non-centrality parameter $\lambda = \sum_i \mu_i^2$, i.e. $W \sim \chi_m^2(\lambda)$. Therefore, the divergence (4) has the form

$$K(\alpha) \propto \frac{1}{2} \mathbb{E}_{W \sim \chi_1^2(\frac{1}{\alpha})} \log W. \quad (4')$$

W can alternatively be represented as a Poisson mixture of ordinary χ^2 distributions: if $Z|J \sim \chi_{m+2J}^2$ for $J \sim \text{Pois}(\frac{\lambda}{2})$ then $W \sim Z$. Therefore, expanding the conditional expectation gives

$$\mathbb{E}_{W \sim \chi_m^2(\lambda)} \log W = \mathbb{E}\left(\mathbb{E}(\log W | J)\right) = \mathbb{E}_{J \sim \text{Pois}(\frac{\lambda}{2})}\left(\mathbb{E}_{W \sim \chi_{m+2J}^2} \log W\right). \quad (20)$$

Since χ_ν^2 is Gamma distribution $\Gamma(\frac{\nu}{2}, \frac{1}{2})$, it can be shown that the logarithmic moment $\mathbb{E}_{W \sim \chi_\nu^2} \log W$ is $\psi(\frac{\nu}{2}) - \log \frac{1}{2}$, where ψ is the digamma function ($\psi(x) = \frac{d}{dx} \log \Gamma(x)$). By expanding expectation of a Poisson random variable we get $\mathbb{E}_{W \sim \chi_m^2(\lambda)} \log W = \log 2 + g_m(\frac{\lambda}{2})$, where

$$g_m(x) = e^{-x} \sum_{j \geq 0} \frac{x^j}{j!} \psi\left(\frac{m+2j}{2}\right). \quad (21)$$

Making use of the property $\psi(z+1) = \psi(z) + \frac{1}{z}$ of the digamma function for $z > 0$, we conclude that the power series in (21) converges for any $x \geq 0$. Therefore the derivative of (21) is given by

$$\frac{d}{dx} g_m(x) = -g_m(x) + e^{-x} \sum_{j \geq 0} \frac{x^j}{j!} \left(\psi\left(\frac{m+2j}{2}\right) + \frac{2}{m+2j} \right). \quad (22)$$

By manipulating the partial sums within (22) we get

$$\frac{d}{dx}g_m(x) = e^{-x} \sum_{j \geq 0} \frac{x^j}{j!} \frac{1}{j + \frac{m}{2}} = e^{-x} x^{-\frac{m}{2}} \sum_{j \geq 0} \frac{1}{j!} \int_0^x t^{j+\frac{m}{2}-1} dt. \quad (23)$$

Furthermore, the functions $t \mapsto \sum_{j=0}^J \frac{1}{j!} t^{j+\frac{m}{2}-1}$ are non-decreasing on $(0, x)$ with growing J and converge to $t^{\frac{m}{2}-1} e^t$, which implies by the Monotone Convergence Theorem that

$$\frac{d}{dx}g_m(x) = e^{-x} x^{-\frac{m}{2}} \int_0^x \sum_{j \geq 0} \frac{1}{j!} t^{j+\frac{m}{2}-1} dt = e^{-x} x^{-\frac{m}{2}} \int_0^x t^{\left(\frac{m}{2}-1\right)} e^t dt. \quad (24)$$

Substituting $u^2 = t$ on $[0, \infty]$ with $2udu = dt$ and letting $I_m: x \mapsto e^{-x^2} \int_0^x u^{m-1} e^{u^2} du$ yields

$$\frac{d(20)}{d\lambda} = \frac{1}{2} \frac{d}{dx}g_m(x) \Big|_{x=\frac{\lambda}{2}} = e^{-x} x^{-\frac{m}{2}} \int_0^{\sqrt{x}} u^{m-1} e^{u^2} du \Big|_{x=\frac{\lambda}{2}} = \left(\sqrt{\frac{2}{\lambda}}\right)^m I_m\left(\sqrt{\frac{\lambda}{2}}\right). \quad (25)$$

Since α is non-negative, it is typically parameterized via its logarithm, whence the derivative of (4') with respect to $\log \alpha$ follows from (25) for $m = 1$ and $\lambda = \frac{1}{\alpha}$:

$$\frac{dK(\alpha)}{d\log \alpha} = -\frac{1}{\sqrt{2\alpha}} I_1\left(\frac{1}{\sqrt{2\alpha}}\right). \quad (26)$$

We compute the Monte-Carlo estimate of (4) on a sample of 10^7 draws over an equally spaced grid of $\log \alpha$ in $[-12, +12]$ of size 4096. The approximation proposed by Molchanov et al. (2017) is given in (27), with coefficients $k_1 = 0.63576$, $k_2 = 1.8732$, and $k_3 = 1.48695$. The derivative of the approximation with respect to $\log \alpha$ follows (26) within 4% of relative tolerance, see fig. 15.

$$(4) \approx \frac{1}{2} \log(1 + e^{-\log \alpha}) + k_1 \sigma(-(k_2 + k_3 \log \alpha)), \quad (27)$$

Similarly, the forward difference estimate of the derivative (26) very closely (up to sampling error). For sake of completeness, we compute a similar Monte-Carlo estimate for the KL divergence term in (13') for \mathbb{C} -valued Variational Dropout with $\beta = 2$, fit the best approximation (27), and compare it against the exact $\log \alpha$ derivative $\frac{d(13')}{d\log \alpha} = e^{-\frac{1}{\alpha}} - 1$.

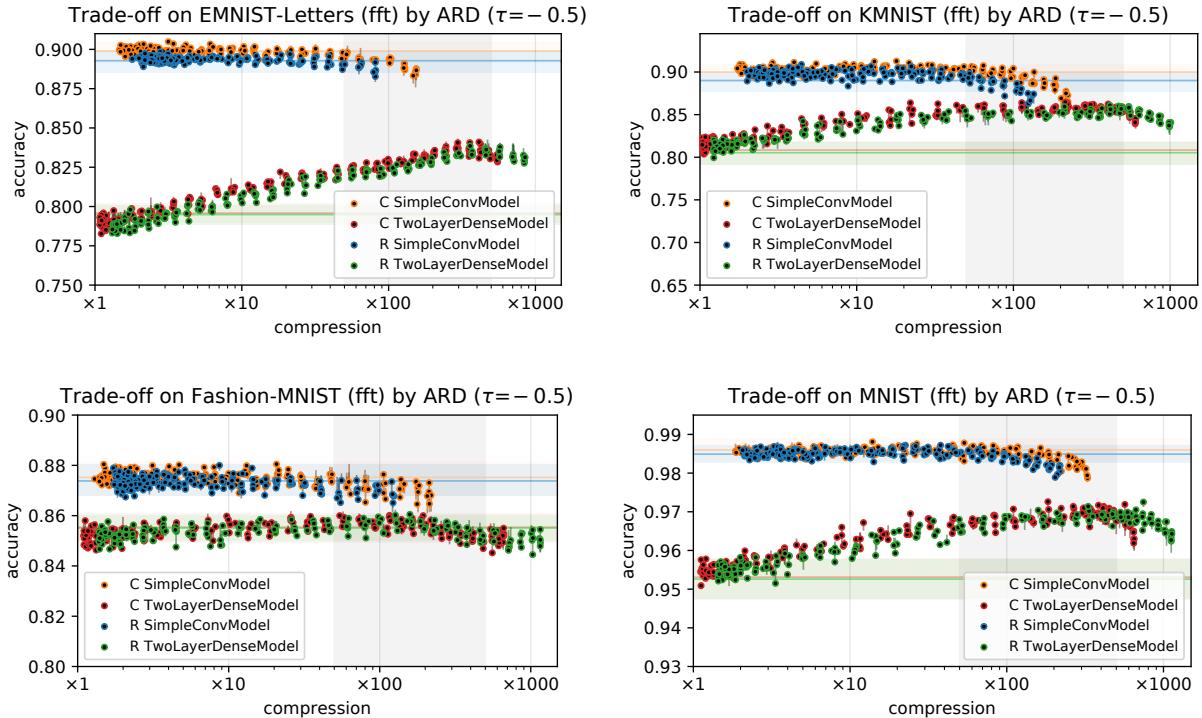


Figure 7. The trade-off of ARD method for \mathbb{R} and \mathbb{C} models using Fourier features.

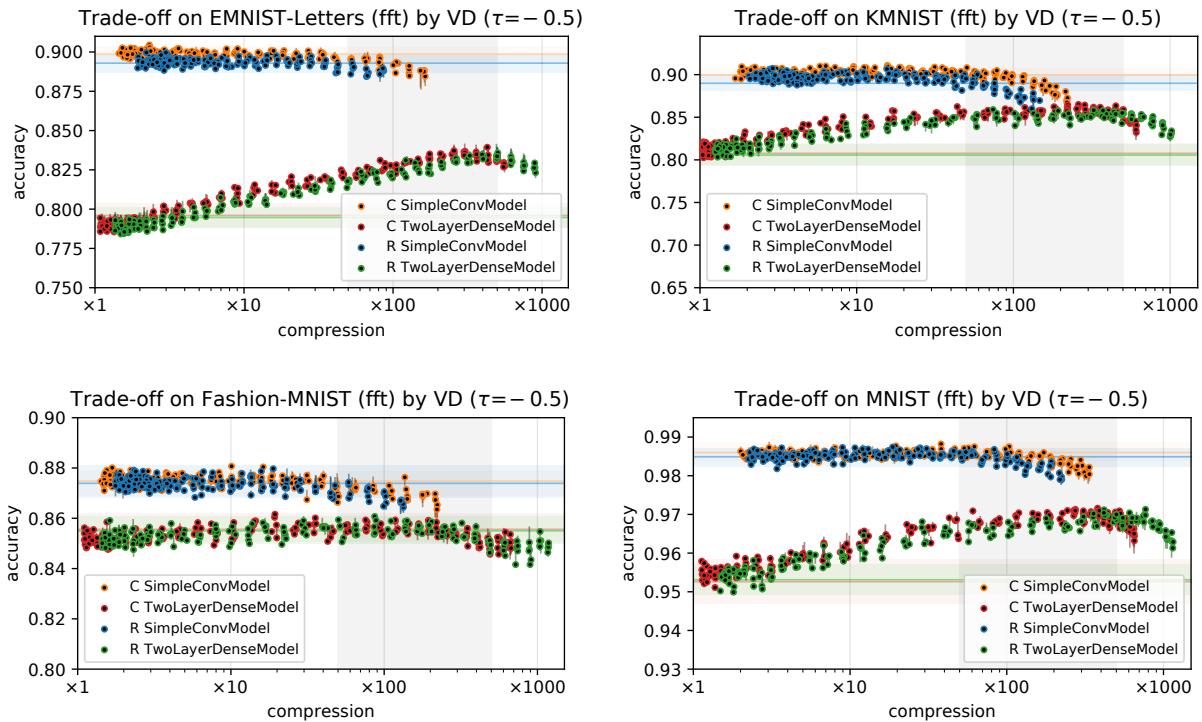


Figure 8. The trade-off of VD method for \mathbb{R} and \mathbb{C} models using Fourier features.

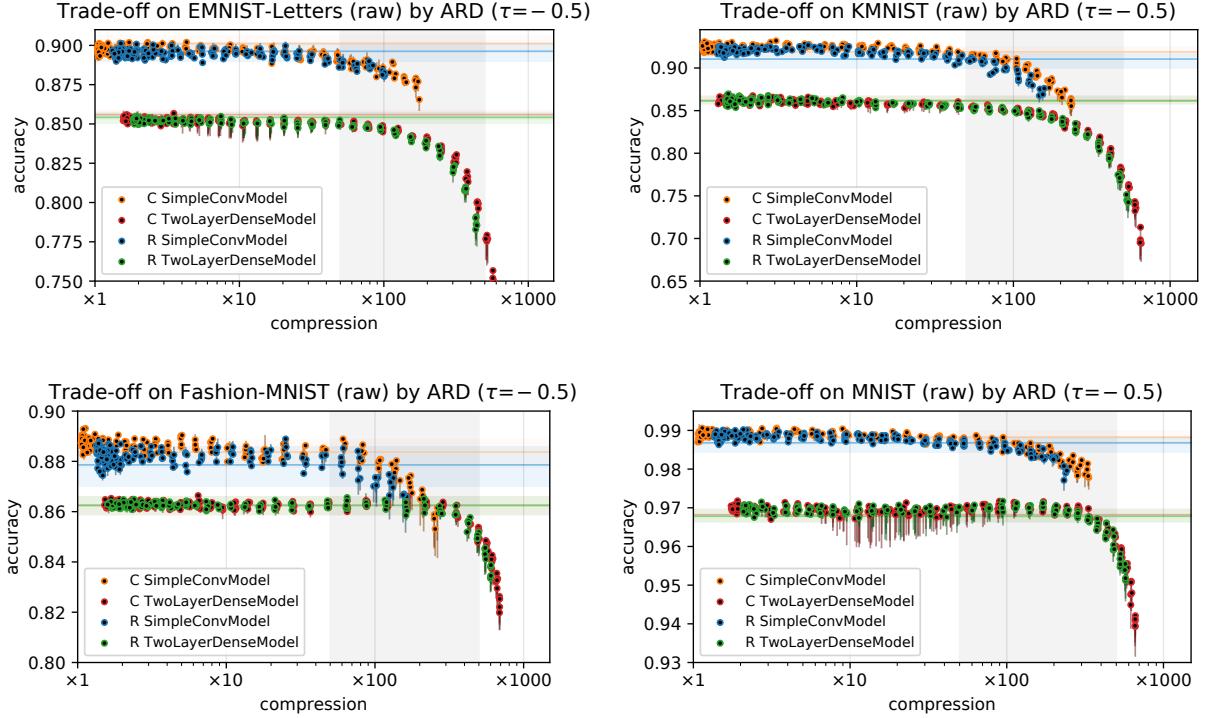


Figure 9. The trade-off of ARD method for \mathbb{R} and \mathbb{C} models using raw features.

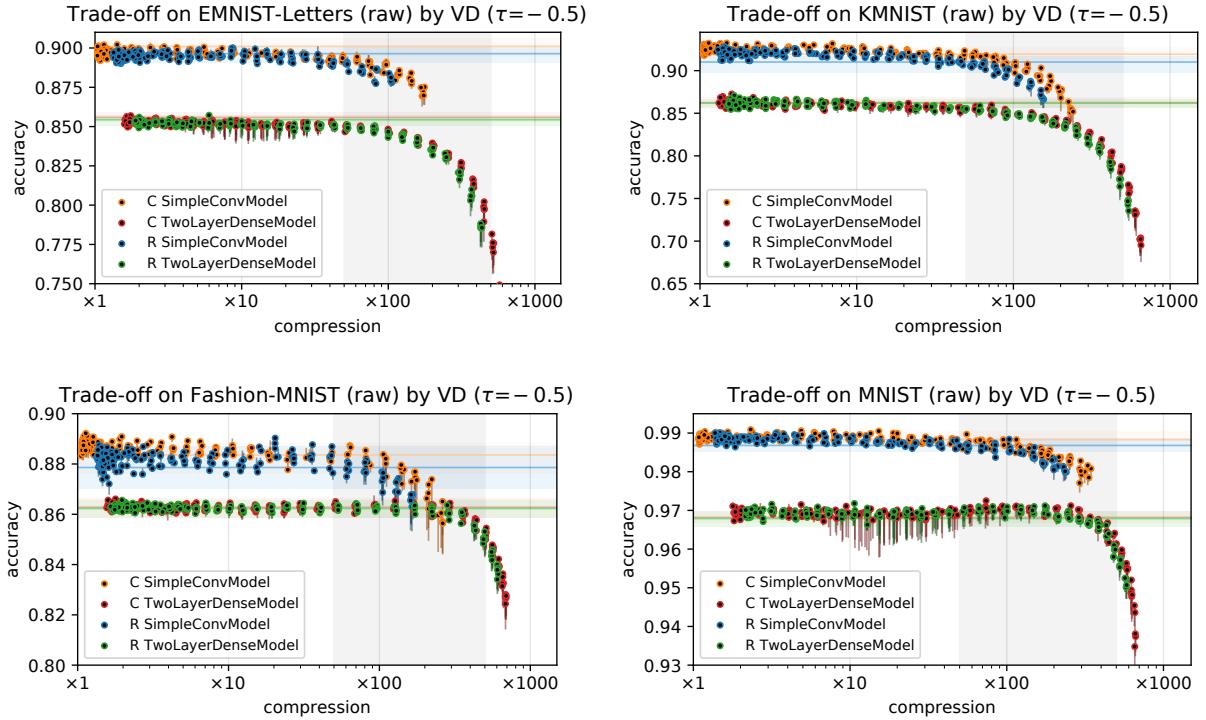


Figure 10. The trade-off of VD method for \mathbb{R} and \mathbb{C} models using raw features.

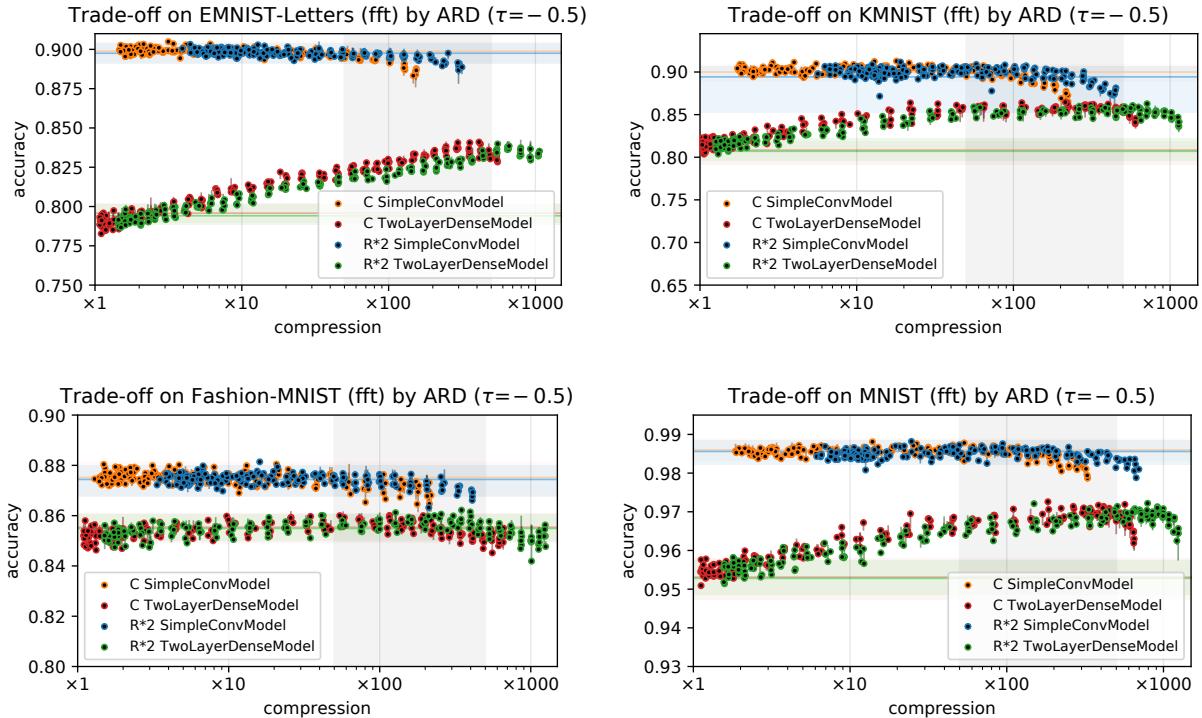


Figure 11. The trade-off of ARD method for $2\mathbb{R}$ and \mathbb{C} models using Fourier features.

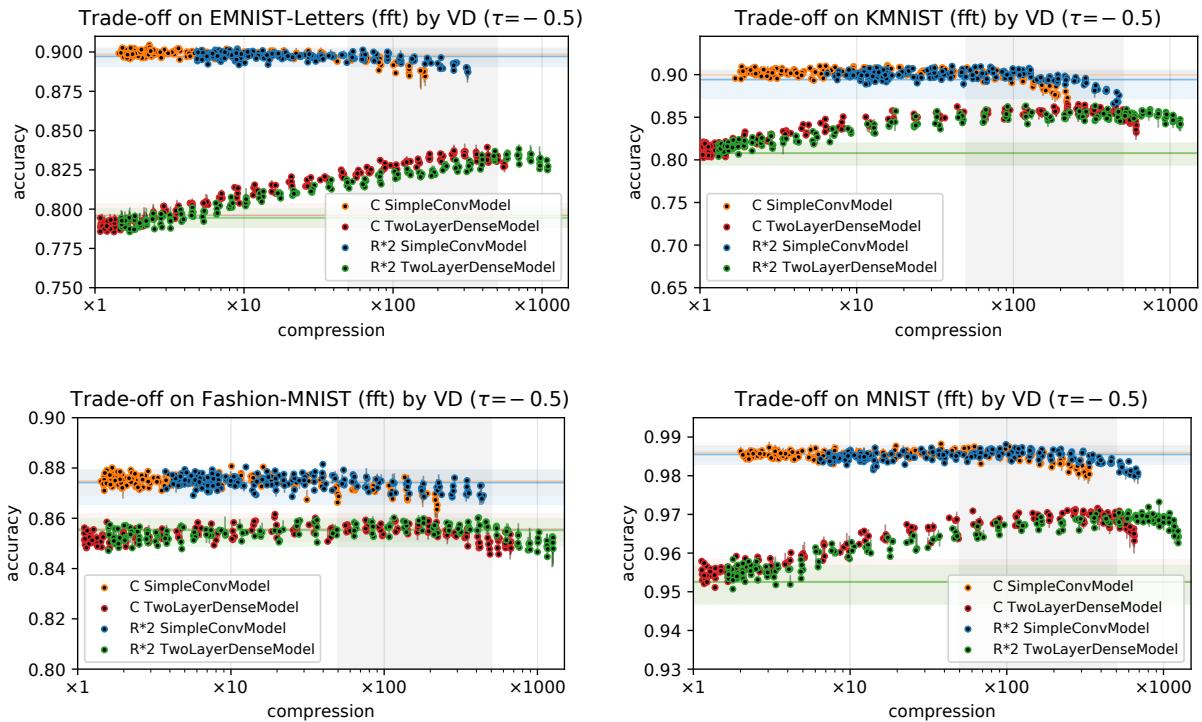


Figure 12. The trade-off of VD method for $2\mathbb{R}$ and \mathbb{C} models using Fourier features.

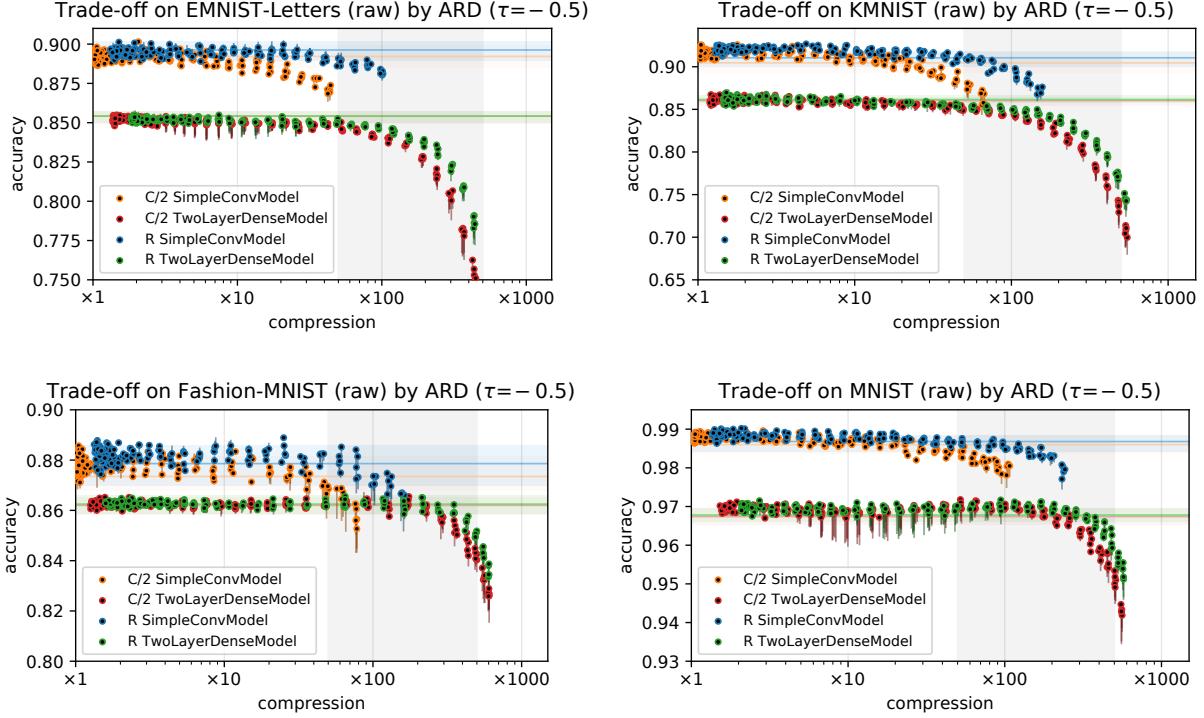


Figure 13. The trade-off of ARD method for \mathbb{R} and $\frac{1}{2}\mathbb{C}$ models using raw features.

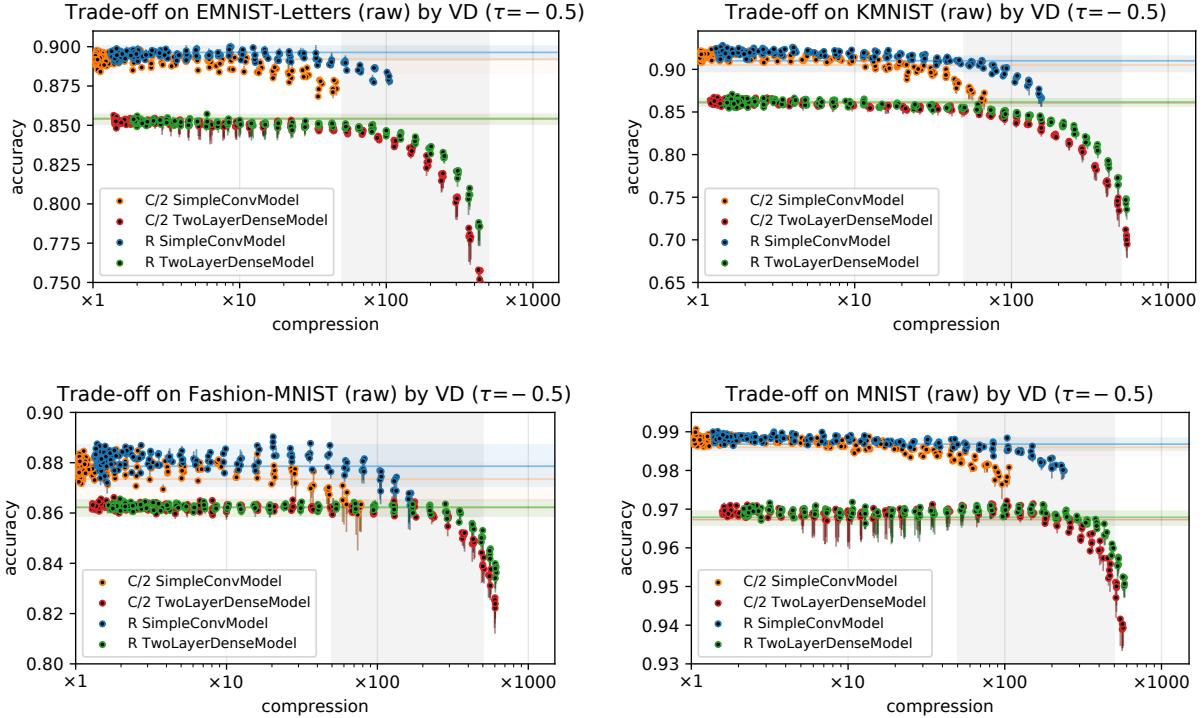


Figure 14. The trade-off of VD method for \mathbb{R} and $\frac{1}{2}\mathbb{C}$ models using raw features.

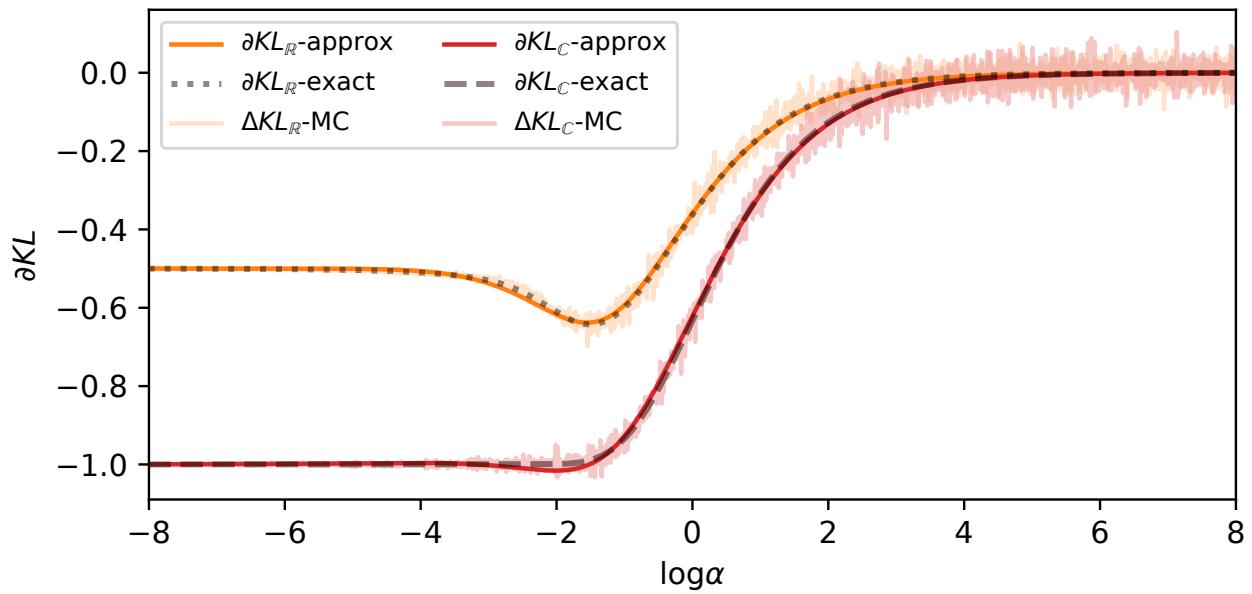


Figure 15. $\frac{dK(\cdot)}{d\log \alpha}$ of the approximation (27), MC estimate of (4), and the exact derivative using (26).