

Supplementary Material to:
 On Computational Power and the Order-Chaos Phase
 Transition in Reservoir Computing

1 Computational Performance for Further Example Tasks

In this section the performance measure p_{exp} defined in eq. (1) in the main text [1] for two further example tasks is shown. In Fig. 1 $p_{\text{exp}}(C, \text{RAND}_5)$ is plotted, averaged over 20 randomly chosen 5-bit functions RAND_5 , circuits C , initial conditions and input streams. A uniform distribution over all 2^{2^5} functions $\text{RAND}_5 : \{-1, +1\}^5 \rightarrow \{-1, 1\}$ was applied. In Fig. 2 the quantity $p_{\text{exp}}(C, \text{AND}_5)$ for the 5-bit AND-function is plotted, i.e. the target output at time t is $\text{AND}_{5,\tau}(u, t) = \max_{i \in \{1, \dots, 5\}} \{u(t - \tau - i)\}$ for delay τ . Shown results are averages over 10 circuits C , initial conditions and input streams. The results shown in Fig. 1 and 2 are qualitatively similar to the results on the PAR-task reported in [1].

2 Definition and Calculation of p_∞

In this section we give a detailed definition of the heuristic performance measure p_∞ and outline an approach to calculate it efficiently.

2.1 Notation

Without loss of generality we may set the readout time t to 0 as the input random process $u(\cdot)$ is assumed to be stationary. Hence the target task of the n last input bits (delay $\tau = 0$) is a function f_T of $u(-1), \dots, u(-n)$. We introduce the following useful notation for the input:

$$\begin{aligned} u^i &:= (u^i(-1), u^i(-2), \dots) \in \{-1, 1\}^{\mathbb{N}} \\ u^{i,t} &:= (u^i(t-1), u^i(t-2), \dots) \in \{-1, 1\}^{\mathbb{N}} \end{aligned}$$

For a vector $x = (x_1, \dots, x_N) \in \mathbb{R}^N$ we use the the p -norm $\|\cdot\|_p$ with $p = 1$ (the Manhattan norm) for measuring distances:

$$\|x\|_1 = \sum_{i=1}^N |x_i|.$$

2.2 Definition of p_∞

In the following we consider two instances N1 and N2 of the same network which solely differ in the input streams $u^1(\cdot)$ and $u^2(\cdot)$ they receive. Let $d(k)$ be the normalized expected distance between the two network states $x^1(0)$ and $x^2(0)$ of N1 and N2 at time 0 for the case where the inputs $u^1(\cdot)$ and $u^2(\cdot)$ only differ at the single time step $t = -k$. Hence for given u^1 , the sequence u^2 is determined by:

$$\begin{aligned} u^2(-i) &= u^1(-i) \quad \forall i \in \mathbb{N} \wedge i \neq k \\ u^2(-k) &= -u^1(-k). \end{aligned}$$

$\forall k \in \mathbb{N}$ the inputs $u^1(-k) \in \{-1, +1\}$ are iid. with probability $p(u^1(-k) = 1) =: 0.5$. $d(k)$ is then formally defined as:

$$d(k) := \frac{1}{N} \langle \langle \|x^1(0) - x^2(0)\|_1 \rangle_W \rangle_u. \quad (1)$$

The average $\langle \cdot \rangle_u$ is taken over all inputs $u^1(\cdot), u^2(\cdot)$ from the ensemble defined above and all initial conditions of the network; $\langle \cdot \rangle_W$ denotes the average over all weight matrices W . If the network is not in the fading memory regime (or equivalently: if it does not have the echo state property, for a formal definition see [2] or [3]) $d(k)$ defined in (1) might well depend on the distribution of initial conditions from which the evolution of x^1 and x^2 starts; we address this subtle point below. We define p_∞ in the following way:

$$p_\infty = \max\left\{ \lim_{k \rightarrow \infty} (d(2) - d(k)), 0 \right\} = \max\{d(2) - d(\infty), 0\}.$$

2.3 The Annealed Approximation

We want to calculate the distance $d(k)$ defined in (1) for large networks $N \rightarrow \infty$ using the annealed approximation (AA) introduced in [4]. The AA consists of approximating the original dynamics of the network by assuming that the weight matrix W is drawn iid. at every time step. First we notice that in the AA and the limit $N \rightarrow \infty$ the following holds:

$$\begin{aligned} d(k) &= \left\langle \left\langle \frac{1}{N} \sum_{b=1}^N |x_b^1(0) - x_b^2(0)| \right\rangle_W \right\rangle_u = \langle \langle |x_a^1(0) - x_a^2(0)| \rangle_W \rangle_u \quad \forall a \in \mathbb{N} \\ d(k) &\xrightarrow{N \rightarrow \infty} \left\langle \left| \sum_{i,j=0}^{2^m-1} p(x_a^1(t) = s_i, x_a^2(t) = s_j | u^{1,t}, u^{2,t}) (s_i - s_j) \right| \right\rangle_u \quad \forall a \in \mathbb{N}. \end{aligned} \quad (2)$$

Here $p(x_a^1(t) = s_i, x_a^2(t) = s_j | u^{1,t}, u^{2,t})$ denotes the joint probability of finding $x_a^1(t)$ in the state s_i and $x_a^2(t)$ in the state s_j given the input $u^{1,t}, u^{2,t}$. Due to the AA this probability is independent of the node index a . Moreover $p(x_a^1(t) = s_i, x_a^2(t) = s_j | u^{1,t}, u^{2,t})$ determines the distribution of the recurrent feedback for the next time step. Hence, in the AA and for $N \rightarrow \infty$ the state of the network is completely described by the joint distribution of a pair of coordinates $x_a^1(t)$ and $x_a^2(t)$ in the state space $\mathcal{S} \times \mathcal{S}$. We therefore define the probability q in the following way:

$$\begin{aligned} q_{ij}(t, u^{1,t}, u^{2,t}) &:= p(x_a^1(t) = s_i, x_a^2(t) = s_j | u^{1,t}, u^{2,t}) \\ q(t, u^{1,t}, u^{2,t}) &:= (q_{ij}(t, u^{1,t}, u^{2,t}))_{i,j \in \{0, \dots, 2^m-1\}}. \end{aligned}$$

Thus $q(t, u^{1,t}, u^{2,t})$ is the $2^m \times 2^m$ matrix whose entry $q_{i,j}$ is the joint probability of finding $x_a^1(t)$ in the state s_i and $x_a^2(t)$ in s_j after applying the input $u^{1,t}$ and $u^{2,t}$ respectively. Using q we can rewrite (2) as:

$$\begin{aligned} d(k) &= \left\langle \left| \sum_{i,j=0}^{2^m-1} q_{ij}(0, u^1, u^2) (s_i - s_j) \right| \right\rangle_u \\ &= \left\langle \left| 2^{1-m} \sum_{i,j=0}^{2^m-1} q_{ij}(0, u^1, u^2) (i - j) \right| \right\rangle_u. \end{aligned}$$

We need to calculate $q_{ij}(t, u^{1,t}, u^{2,t})$ at time $t = 0$ in order to determine $d(k)$. This can be achieved iteratively via the mapping S representing the transition from time step t to $t + 1$ by applying the input pair $u^1(t)$ and $u^2(t)$:

$$q(t+1, u^{1,t+1}, u^{2,t+1}) = S(q(t, u^{1,t}, u^{2,t}), u^1(t), u^2(t)).$$

The k -fold composition of S with the inputs u^1 and u^2 is denoted by $S_{u^1, u^2}^{(k)}$:

$$\begin{aligned} q(t, u^{1,t}, u^{2,t}) &= S_{u^1, u^2}^{(k)}(q(t-k, u^{1,t-k}, u^{2,t-k})) \\ q(t, u^{1,t}, u^{2,t}) &= \lim_{k \rightarrow \infty} \left(S_{u^1, u^2}^{(k)}(q(t-k, u^{1,t-k}, u^{2,t-k})) \right). \end{aligned}$$

For the sake of fast numerical evaluation, we calculate $q(0, u^1, u^2)$ by starting at $t = -n$ with the initial condition q^* :

$$q(0, u^1, u^2) = S_{u^1, u^2}^{(n)}(q^*).$$

The initial condition $q^* := 2^{-m} \text{id}_{2^m \times 2^m}$ (where $\text{id}_{n \times m}$ is the $n \times m$ identity matrix) was chosen because of the following relation:

$$q^* = \langle q(t, u^1, u^1) \rangle_{u^1}.$$

2.4 Separation approximation

Unfortunately the complexity to calculate q scales like $\mathcal{O}(2^{2m})$ (capital \mathcal{O} notation) with the number of bits m . This basically renders the approach described above useless for $m > 5$. However, this can be circumvented by generalizing the AA in the following way. Since the state $x_a^i(t)$ of neuron a is quantized with m bits, we can write it in the following way:

$$\begin{aligned} x_a^i(t) &= \sum_{l=0}^{m-1} 2^{-l} (b_l^i(t) - 1/2), \quad b_l^i(t) \in \{0, 1\} \\ B_l(x_a^i(t)) &:= b_l^i(t). \end{aligned} \quad (3)$$

According to (3) there is a unique binary representation of $x_a^i(t)$ given by $(b_0^i(t), \dots, b_{m-1}^i(t))$; the mapping $B_l(\cdot)$ maps a state to its l^{th} bit. Equation (3) can be interpreted as effectively replacing unit a with m binary units of states $b_l^i(t)$ whose outputs are reduced by $1/2$ and multiplied by 2^{-l} and finally summed up; this is still exact. Now we assume that each of these m units receives input drawn independently from the input distribution and has different weights drawn independently from $N(0, \sigma^2)$ every time step; hence this approach generalizes the AA. Therefore, for given presynaptic input the $b_l^i(t)$ are independent for different i and l under this approximation. Thus:

$$\begin{aligned} q_{ij}(t, u^{1,t}, u^{2,t}) &\approx \prod_{l=0}^{m-1} q_{B_l(s_i), B_l(s_j)}^l(t, u^{1,t}, u^{2,t}) \\ q^l(t, u^{1,t}, u^{2,t}) &= (q_{b_1^1, b_1^2}^l(t, u^{1,t}, u^{2,t}))_{b_1^1, b_1^2 \in \{0, 1\}}. \end{aligned}$$

$q^l(t, u^{1,t}, u^{2,t})$ is the 2×2 matrix, whose entry $q_{b_1^1, b_1^2}^l(t, u^{1,t}, u^{2,t})$ is the joint probability of finding the bit number l of unit $x_a^1(t)$ in state $b_1^1 \in \{0, 1\}$ and of unit $x_a^2(t)$ in state $b_1^2 \in \{0, 1\}$. Under this approximation we only have to calculate $4m$ matrix entries instead of the 2^{2m} entries, which is a considerable reduction of complexity.

We denote the update mapping for q^l by S^l :

$$q^l(t+1, u^{1,t+1}, u^{2,t+1}) = S^l(q^l(t, u^{1,t}, u^{2,t}), u^1(t), u^2(t)).$$

An explicit form for S^l can be derived in the following way. First we condition the probability $q^l(t+1, u^{1,t+1}, u^{2,t+1})$ on the presynaptic input $h^1 = (h_1^1, \dots, h_K^1)$ for network N1 and $h^2 = (h_1^2, \dots, h_K^2)$ for network N2 and on the weight matrix W for this time step (which is the same for N1 and N2). The pairs (h_i^1, h_i^2) , $i \in 1, \dots, K$ are iid. according to $q(t, u^{1,t}, u^{2,t})$. This yields:

$$\begin{aligned} q^l(t+1, u^{1,t+1}, u^{2,t+1}) &= \langle \langle q^l(t+1, u^{1,t+1}, u^{2,t+1} | h^1, h^2, w) \rangle_W \rangle_{h^1, h^2} \\ &= \sum_{h^1, h^2} p(h^1, h^2) \int dW p(W) q^l(t+1, u^{1,t+1}, u^{2,t+1} | h^1, h^2, W). \end{aligned} \quad (4)$$

We used the following abbreviations :

$$\begin{aligned}
\langle X(h^1, h^2) \rangle_{h^1, h^2} &:= \sum_{h^1, h^2} p(h^1, h^2) X(h^1, h^2) \\
&= \sum_{h_1^1, h_1^2} \cdots \sum_{h_K^1, h_K^2} \left(\prod_{i=1}^K q_{h_i^1, h_i^2}(t, u^{1,t}, u^{2,t}) \right) X(h^1, h^2) \\
\langle X(W) \rangle_W &:= \int dW p(W) X(W) = \int_{\mathbb{R}} \cdots \int_{\mathbb{R}} X(W) p(w_1, \dots, w_K) dw_1 \cdots dw_K.
\end{aligned} \tag{5}$$

Here w_1, \dots, w_K denote the K presynaptic weights. Conditioned on the input and the weights, the network realizations N1 and N2 are independent:

$$\begin{aligned}
q_{ij}^l(t+1, u^{1,t+1}, u^{2,t+1} | h^1, h^2, w) &= p(b_i^1(t+1) = i | h^1, w) p(b_i^2(t+1) = j | h^2, w) \\
&= \delta(B_l(f_m(w^T h^1 + u^1(t))), i) \delta(B_l(f_m(w^T h^2 + u^2(t))), j) \\
w^T h^i &:= \sum_{\alpha=1}^K w_{\alpha} h_{\alpha}^i.
\end{aligned}$$

$\delta(\cdot, \cdot)$ denotes the Kronecker-Delta of its two arguments. The K -fold integral over the weights appearing in (4) can be reduced to a single integral:

$$\begin{aligned}
\int dW p(W) q_{ij}^l(t+1, u^{1,t+1}, u^{2,t+1} | h^1, h^2, W) &= F_{ij}^l(h^1, h^2, u^1(t), u^2(t)) \\
F_{ij}^l(h^1, h^2, u^1(t), u^2(t)) &= \frac{1}{(8\pi \det(C) \Sigma_{22})^{1/2}} \sum_{\alpha=0}^{2^l-1} \int_{I_{\alpha, i} - u^1(t)} dv \exp\left(-\frac{v^2}{2} \left(\Sigma_{11} - \frac{\Sigma_{12}^2}{\Sigma_{22}}\right)\right) \times \\
&\quad \sum_{\beta=0}^{2^l-1} \left[\operatorname{erf}\left(\frac{(I_{\beta, j}^+ - u^2(t)) \Sigma_{22} + \Sigma_{21} v}{(2\Sigma_{22})^{1/2}}\right) - \operatorname{erf}\left(\frac{(I_{\beta, j}^- - u^2(t)) \Sigma_{22} + \Sigma_{21} v}{(2\Sigma_{22})^{1/2}}\right) \right]
\end{aligned}$$

Here we use the following definitions:

$$\begin{aligned}
C &= \sigma^2 \begin{pmatrix} (h^1)^T h^1 & (h^1)^T h^2 \\ (h^2)^T h^1 & (h^2)^T h^2 \end{pmatrix} \\
\Sigma_{ij} &= (C^{-1})_{ij} \\
\bigcup_{\alpha=0}^{2^l-1} I_{\alpha, i} &= \operatorname{supp}(\delta(B_l(f_m(\cdot)), i)) \\
I_{\alpha, i} &= [I_{\alpha, i}^-, I_{\alpha, i}^+] \\
I_{\alpha, i}^- &:= \tanh^{-1}(2^{-l+1}\alpha - 1) \\
I_{\alpha, i}^+ &:= \tanh^{-1}(2^{-l+1}(\alpha + 1) - 1).
\end{aligned}$$

The support of the function $\delta(B_l(f_m(\cdot)), i)$ is the union of 2^l disjoint intervals $I_{\alpha, i}$ with lower bound $I_{\alpha, i}^-$ and upper bound $I_{\alpha, i}^+$.

Using the expression F_{ij}^l the update can finally be written as:

$$q_{ij}^l(t+1, u^{1,t+1}, u^{2,t+1}) = \langle F_{ij}^l(h^1, h^2, u^1(t), u^2(t)) \rangle_{h^1, h^2}.$$

For the sake of computational tractability for larger m and K , we do not evaluate the $2K$ sums explicitly involved in the average over the presynaptic input $\langle \cdot \rangle_{h^1, h^2}$. Instead we determine this expectation value by a finite number of samples from the joint distribution $p(h^1, h^2)$; this sampling is easily realizable since $p(h^1, h^2)$ is of the product form given in (5); sample size for all experiments was chosen to be 150.

References

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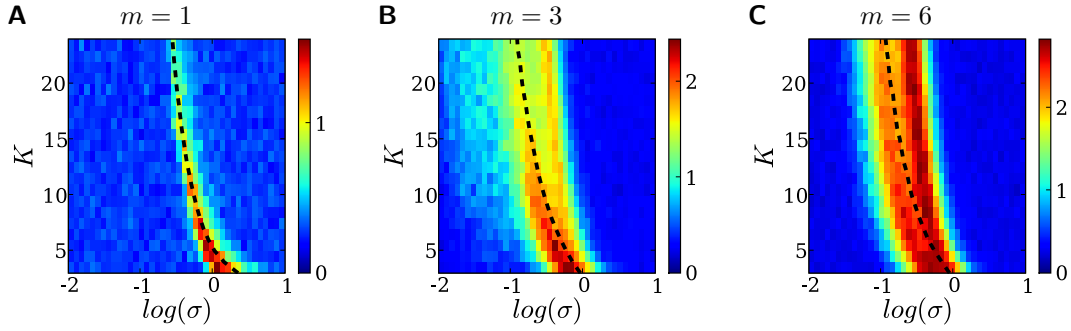


Figure 1: The performance $p_{\text{exp}}(C, \text{RAND}_5)$ for three different quantization levels $m = 1, 3, 6$, averaged over 20 randomly drawn functions of 5 bits, circuits C , initial conditions and input streams. $p_{\text{exp}}(C, \text{RAND}_5)$ is plotted as a function of the network in-degree K and the weight STD σ . The networks size is $N = 150$. The input time series have length 10000. The solid line represents the numerically found critical line.

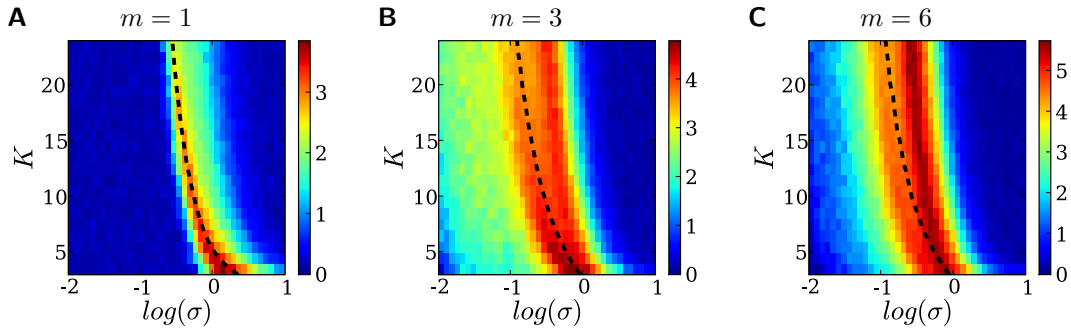


Figure 2: Same figure as Fig. 1 showing the performance $p_{\text{exp}}(C, \text{AND}_5)$ for the 5-bit AND-task averaged over 10 circuits C , initial conditions and input streams.