Machine Learning Algorithms

LMS Analysis

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• Let's consider the steepest descent iteration:

$$\underline{\hat{w}}_{k} = \underline{\hat{w}}_{k-1} + \mu \left(\underline{r}_{\underline{\mathbf{x}}\mathbf{d}}^{*} - R_{\underline{\mathbf{x}}\underline{\mathbf{x}}}^{*} \underline{\hat{w}}_{k-1}\right), k = 1, 2, \dots$$

- Obviously, the knowledge of the cross-correlation vector and the ACF matrix is required.
- If this knowledge is not present, we could operate with estimates. The most simple ones are:

$$\begin{split} \hat{R}_{\underline{\mathbf{x}}\underline{\mathbf{x}}} &= \underline{\mathbf{x}}_k \, \underline{\mathbf{x}}_k^H; \qquad \hat{\underline{r}}_{\underline{\mathbf{x}}\mathbf{d}} = \underline{\mathbf{x}}_k \mathbf{d}_k^* \\ \underline{\mathbf{x}}_k^T &= \left[\mathbf{x}(k), \mathbf{x}(k-1), ..., \mathbf{x}(k-M+1) \right] \\ \text{Univ.-Prof. Dr.-Ing. Markus Rupp} \end{split}$$



- They are called instantaneous estimates.
- Assuming stationary processes, also other estimates are possible, for example

$$\hat{R}_{\underline{\mathbf{x}}\underline{\mathbf{x}}} = \frac{1}{N} \sum_{l=0}^{N-1} \ \underline{\mathbf{x}}_{k-l} \underline{\mathbf{x}}_{k-l}^{H}; \qquad \hat{\underline{r}}_{\underline{\mathbf{x}}\mathbf{d}} = \frac{1}{N} \sum_{l=0}^{N-1} \ \underline{\mathbf{x}}_{k-l} \mathbf{d}_{k-l}^{*}$$

- We will see in the next chapter that this choice leads to a special form, the so-called RLS algorithm.
- Plugging in the steepest descent iterations the instantaneous estimates, we obtain:

$$\underline{\hat{\mathbf{w}}}_{k} = \underline{\hat{\mathbf{w}}}_{k-1} + \mu \underline{\mathbf{x}}_{k}^{*} \left(\mathbf{d}_{k} - \underline{\mathbf{x}}_{k}^{T} \underline{\hat{\mathbf{w}}}_{k-1} \right), k = 1, 2, \dots$$



Utilizing such instantaneous values,

$$\underline{\hat{\mathbf{w}}}_{k} = \underline{\hat{\mathbf{w}}}_{k-1} + \mu \underline{\mathbf{x}}_{k}^{*} \underbrace{\left(\mathbf{d}_{k} - \underline{\mathbf{x}}_{k}^{T} \underline{\hat{\mathbf{w}}}_{k-1}\right)}_{\mathbf{e}_{a,k}}, k = 1,2,...$$

- The update equation of the LMS algorithm is obtained.
- The error signal

$$\widetilde{\mathbf{e}}_{a,k} = \mathbf{d}_k - \underline{\mathbf{x}}_k^T \underline{\hat{\mathbf{w}}}_{k-1}$$

- is called the disturbed/distorted a-priori error signal.
- Due to the reference model, there is also an undisturbed a-priori error signal: $\mathbf{e}_{a,k} = \mathbf{x}_k^T \mathbf{w}_o \mathbf{x}_k^T \hat{\mathbf{w}}_{k-1} = \mathbf{x}_k^T \mathbf{\widetilde{w}}_{k-1}$



• Correspondingly, would we use the estimates at time instant k rather than k-1, we call the error a-posteriori error signals:

$$\mathbf{\tilde{e}}_{p,k} = \mathbf{d}_k - \underline{\mathbf{x}}_k^T \underline{\hat{\mathbf{w}}}_k$$

$$\mathbf{e}_{p,k} = \underline{\mathbf{x}}_k^T \underline{\mathbf{w}}_o - \underline{\mathbf{x}}_k^T \underline{\hat{\mathbf{w}}}_k = \underline{\mathbf{x}}_k^T \underline{\widetilde{\mathbf{w}}}_k$$

• Note that due to the nature of the recursive algorithm, the estimates are random processes:

$$\underline{\hat{\mathbf{w}}}_{k} = \underline{\hat{\mathbf{w}}}_{k-1} + \mu \underline{\mathbf{x}}_{k}^{*} \left(\mathbf{d}_{k} - \underline{\mathbf{x}}_{k}^{T} \underline{\hat{\mathbf{w}}}_{k-1} \right), k = 1, 2, \dots$$



• Many different choices for the step-size are possible:

arbitrary μ_k : stochastic gradient algorithm

$$\mu_k = \frac{\alpha}{\left\|\underline{\mathbf{x}}_k\right\|_2^2}$$
 with $\alpha > 0$, (data-) Normalized LMS algorithm = NLMS

$$\mu_k = \frac{\alpha}{1 + \alpha \|\mathbf{x}_k\|_2^2}$$
 with $\alpha > 0$, a - posteriori form of the LMS algorithm

$$\mu_k = \frac{\alpha}{\varepsilon + \|\mathbf{x}_k\|_2^2}$$
 with $\varepsilon > 0$, the so - called $\varepsilon - \text{NLMS}$ algorithm



Its popularity is also deserved by its mathematical simplicity. Not only

Can be minimized but also

$$\mathrm{E}\left[\left|\tilde{\mathrm{e}}_{a,k}\right|^{2}\right]$$

$$\mathrm{E}\left[\left|f\left[\tilde{\mathrm{e}}_{a,k}\right]\right|^{2}\right]$$

• With arbitrary nonlinear functions f[]. For example $f[x]=x^{K/2}$ leads to the well known Least-Mean-K algorithm (LMK):

$$\underline{\hat{\mathbf{w}}}_{k} = \underline{\hat{\mathbf{w}}}_{k-1} + \mu \underline{\mathbf{x}}_{k}^{*} |\widetilde{\mathbf{e}}_{a,k}|^{K-2} \widetilde{\mathbf{e}}_{a,k}$$



Other variants of the LMS algorithm are:

$$: \underline{\hat{\mathbf{w}}}_{k} = \underline{\hat{\mathbf{w}}}_{k-1} + \mu \underline{\mathbf{x}}_{k}^{*} \left| \widetilde{\mathbf{e}}_{a,k} \right|^{2} \widetilde{\mathbf{e}}_{a,k}$$

Least – Mean Mixed Norm:
$$\underline{\hat{\mathbf{w}}}_{k} = \underline{\hat{\mathbf{w}}}_{k-1} + \mu \underline{\mathbf{x}}_{k}^{*} \left(\beta + (1-\beta) |\widetilde{\mathbf{e}}_{a,k}|^{2} \right) \underline{\widetilde{\mathbf{e}}}_{a,k}$$

$$\underline{\hat{\mathbf{w}}}_{k} = \beta \underline{\hat{\mathbf{w}}}_{k-1} + \mu \underline{\mathbf{x}}_{k}^{*} \widetilde{\mathbf{e}}_{a,k}$$

$$\underline{\hat{\mathbf{w}}}_{k} = \underline{\hat{\mathbf{w}}}_{k-1} + \mu \underline{\mathbf{x}}_{k}^{*} \mathbf{sgn} [\widetilde{\mathbf{e}}_{a,k}]$$



- Since the LMS algorithm uses instantaneous estimates of cross correlation vector and ACF matrix, we cannot expect that the algorithm behaves identical to the steepest descent algorithm.
- We therefore have to analyze the algorithm by computing its behavior in mean and mean square.



Independence assumptions

- The observation of d_k originates from a reference model: $\mathbf{d}_k = \underline{\mathbf{w}}_0^T \underline{\mathbf{x}}_k + \mathbf{v}_k$, \mathbf{v}_k and $\underline{\mathbf{x}}_k$ are of zero mean.
- The regression vectors $\underline{\mathbf{x}}_k$ are statistically independent for different time instants: $f_{\mathbf{x}\mathbf{x}}(\underline{\mathbf{x}}_k,\underline{\mathbf{x}}_l) = f_{\mathbf{x}\mathbf{x}}(\underline{\mathbf{x}}_k) f_{\mathbf{x}\mathbf{x}}(\underline{\mathbf{x}}_l)$ for k and l different.
- The input process \mathbf{x}_k is spherically invariant and Gaussian distributed with zero mean.
- The additive noise \mathbf{v}_k is statistically independent of the driving process \mathbf{x}_k .
- Note that through these conditions, $\underline{\mathbf{w}}_k$ is statistically independent of $\underline{\mathbf{x}}_l$; l>k.



Spherically Invariant Processes

- A complex-valued variable z=x+jy is called Gaussian when x and y are joint Gaussian.

• The second moment of **z** is given by:
$$R_{\mathbf{z}\mathbf{z}} = \mathbf{E} \big[\mathbf{z}\mathbf{z}^* \big] = R_{\mathbf{x}\mathbf{x}} + R_{\mathbf{y}\mathbf{y}} + j \big(R_{\mathbf{y}\mathbf{x}} - R_{\mathbf{x}\mathbf{y}} \big)$$

• Note that the information of R_{zz} is not sufficient to conclude back to R_{xx} , R_{yy} and R_{xy} . For this more information is required, for example:

$$R_{\mathbf{z}\mathbf{z}^*} = \mathbf{E}[\mathbf{z}\mathbf{z}] = R_{\mathbf{x}\mathbf{x}} - R_{\mathbf{y}\mathbf{y}} + j(R_{\mathbf{y}\mathbf{x}} + R_{\mathbf{x}\mathbf{y}})$$

- For spherically invariant processes, we have R_{77*}=0, or equivalently $R_{xx}=R_{yy}$ and $R_{xy}=-R_{yx}$.
- The joint density function of a vector \underline{z} is thus given by:

$$f_{\underline{\mathbf{x}},\underline{\mathbf{y}}}(\underline{\mathbf{x}},\underline{\mathbf{y}}) = f_{\underline{\mathbf{z}}}(\underline{\mathbf{z}}) = \frac{1}{\pi^p} \frac{1}{\det(R_{\mathbf{z}\underline{\mathbf{z}}})} \exp\left(-\underline{\mathbf{z}}^H R_{\underline{\mathbf{z}}\underline{\mathbf{z}}}^{-1}\underline{\mathbf{z}}\right)$$



The parameter error vector in the mean:

$$\underline{\widetilde{\mathbf{w}}}_{k} = \underline{w}_{o} - \underline{\widehat{\mathbf{w}}}_{k}$$

$$\underline{\widetilde{\mathbf{w}}}_{k} = \left(I - \mu \underline{\mathbf{x}}_{k}^{*} \underline{\mathbf{x}}_{k}^{T}\right) \underline{\widetilde{\mathbf{w}}}_{k-1} - \mu \underline{\mathbf{x}}_{k}^{*} \mathbf{v}_{k}$$

$$E[\underline{\widetilde{\mathbf{w}}}_{k}] = E[\left(I - \mu \underline{\mathbf{x}}_{k}^{*} \underline{\mathbf{x}}_{k}^{T}\right) \underline{\widetilde{\mathbf{w}}}_{k-1}] - \mu E[\underline{\mathbf{x}}_{k}^{*} \mathbf{v}_{k}]$$

• With the independence assumptions:

$$E[\widetilde{\underline{\mathbf{w}}}_{k}] = (I - \mu R_{\underline{\mathbf{x}}\underline{\mathbf{x}}}^{*}) E[\widetilde{\underline{\mathbf{w}}}_{k-1}]$$
• Behavior identical to steepest descent!



Behavior identical to steepest descent!

$$E\left[\underline{\widetilde{\mathbf{w}}}_{k}\right] = \left(I - \mu R_{\underline{\mathbf{x}}\underline{\mathbf{x}}}^{*}\right) E\left[\underline{\widetilde{\mathbf{w}}}_{k-1}\right]$$

Thus convergence guaranteed for

$$0 < \mu < \frac{2}{\lambda_{\text{max}}}$$



• The parameter error vector in the mean square sense:

$$P_{k} = \mathbf{E} \left[(\underline{w}_{o} - \underline{\hat{\mathbf{w}}}_{k}) (\underline{w}_{o} - \underline{\hat{\mathbf{w}}}_{k})^{H} \right] = \mathbf{E} \left[\underline{\widetilde{\mathbf{w}}}_{k} \underline{\widetilde{\mathbf{w}}}_{k}^{H} \right]$$

$$= \mathbf{E} \left[(I - \mu \underline{\mathbf{x}}_{k}^{*} \underline{\mathbf{x}}_{k}^{T}) P_{k-1} (I - \mu \underline{\mathbf{x}}_{k}^{*} \underline{\mathbf{x}}_{k}^{T})^{H} \right] + \mu^{2} \mathbf{E} \left[\underline{\mathbf{x}}_{k}^{*} \underline{\mathbf{x}}_{k}^{T} |\mathbf{v}_{k}|^{2} \right]$$

$$= P_{k-1} - \mu \mathbf{E} \left[\underline{\mathbf{x}}_{k}^{*} \underline{\mathbf{x}}_{k}^{T} P_{k-1} \right] - \mu \mathbf{E} \left[P_{k-1} \underline{\mathbf{x}}_{k}^{*} \underline{\mathbf{x}}_{k}^{T} \right] + \mu^{2} \mathbf{E} \left[\underline{\mathbf{x}}_{k}^{*} \underline{\mathbf{x}}_{k}^{T} P_{k-1} \underline{\mathbf{x}}_{k}^{*} \underline{\mathbf{x}}_{k}^{T} \right]$$

$$+ \mu^{2} \mathbf{E} \left[\underline{\mathbf{x}}_{k}^{*} \underline{\mathbf{x}}_{k}^{T} |\mathbf{v}_{k}|^{2} \right]$$



Most problematic is the term

• It can be solved explicitly for (spherically) Gaussian processes. We distinguish real-valued processes:

$$E\left[\underline{\mathbf{x}}_{k}\,\underline{\mathbf{x}}_{k}^{T}P_{k-1}\,\underline{\mathbf{x}}_{k}\,\underline{\mathbf{x}}_{k}^{T}\right] = 2R_{\underline{\mathbf{x}}\underline{\mathbf{x}}}P_{k-1}R_{\underline{\mathbf{x}}\underline{\mathbf{x}}} + R_{\underline{\mathbf{x}}\underline{\mathbf{x}}} \operatorname{trace}\left(P_{k-1}R_{\underline{\mathbf{x}}\underline{\mathbf{x}}}\right)$$

Complex-valued processes:

$$E\left[\underline{\mathbf{x}}_{k}^{*}\underline{\mathbf{x}}_{k}^{T}P_{k-1}\underline{\mathbf{x}}_{k}^{*}\underline{\mathbf{x}}_{k}^{T}\right] = R_{\underline{\mathbf{x}}\underline{\mathbf{x}}}^{*}P_{k-1}R_{\underline{\mathbf{x}}\underline{\mathbf{x}}}^{*} + R_{\underline{\mathbf{x}}\underline{\mathbf{x}}}^{*} \operatorname{trace}\left(P_{k-1}R_{\underline{\mathbf{x}}\underline{\mathbf{x}}}^{*}\right)$$



• The procedure can be extended to general spherically invariant processes. In the general case we obtain

$$E\left[\underline{\mathbf{x}}_{k}^{*}\underline{\mathbf{x}}_{k}^{T}P_{k-1}\underline{\mathbf{x}}_{k}^{*}\underline{\mathbf{x}}_{k}^{T}\right] = \gamma R_{\underline{\mathbf{x}}\underline{\mathbf{x}}}^{*}P_{k-1}R_{\underline{\mathbf{x}}\underline{\mathbf{x}}}^{*} + R_{\underline{\mathbf{x}}\underline{\mathbf{x}}}^{*} \operatorname{trace}\left(P_{k-1}R_{\underline{\mathbf{x}}\underline{\mathbf{x}}}^{*}\right)$$

- with some positive constant $\gamma=1$ for complex and $\gamma=2$ for real valued processes. .
- The update recursion for the parameter error covariance matrix reads now:

$$P_{k} = P_{k-1} - \mu R_{\underline{\mathbf{x}}}^{*} P_{k-1} - \mu P_{k-1} R_{\underline{\mathbf{x}}}^{*} + \mu^{2} \gamma R_{\underline{\mathbf{x}}}^{*} P_{k-1} R_{\underline{\mathbf{x}}}^{*}$$

$$+ \mu^{2} R_{\underline{\mathbf{x}}}^{*} \operatorname{trace} \left(P_{k-1} R_{\underline{\mathbf{x}}}^{*} \right) + \mu^{2} R_{\underline{\mathbf{x}}}^{*} \sigma_{\mathbf{v}}^{2}$$



Orthogonalization for the ACF matrix leads to:

$$Q^H R_{\mathbf{x}\mathbf{x}}^* Q = \Lambda; \qquad Q^H P_k Q = C_k$$

We thus obtain:

$$C_{k} = C_{k-1} - \mu \Lambda C_{k-1} - \mu C_{k-1} \Lambda + \mu^{2} \gamma \Lambda C_{k-1} \Lambda$$
$$+ \mu^{2} \Lambda \operatorname{trace} \left[\Lambda C_{k-1} \right] + \mu^{2} \Lambda \sigma_{\mathbf{v}}^{2}$$

But note that:

$$E\left[\left\|\underline{\widetilde{\mathbf{w}}}_{k}\right\|_{2}^{2}\right] = \operatorname{trace}[P_{k}] = \operatorname{trace}[C_{k}]$$



• Thus, only the diagonal terms of C_k are of importance. Order diagonal terms in vector \underline{c}_k :

$$\underline{c}_{k} = B\underline{c}_{k-1} + \mu^{2}\sigma_{\mathbf{v}}^{2}\underline{\lambda}$$

$$B = I - 2\mu\Lambda + \mu^{2}\gamma\Lambda^{2} + \mu^{2}\underline{\lambda}\underline{\lambda}^{T}$$

$$= \begin{cases} 1 - 2\mu\lambda_{i} + \mu^{2}(1+\gamma)\lambda_{i}^{2} & \mathbf{main diagonal} \\ \mu^{2}\lambda_{i}\lambda_{j} & \mathbf{else} \end{cases}$$



• Theorem 3.1: Under the given conditions, the LMS algorithm is convergent in the mean-square sense, when:

$$0 < \mu < \frac{2}{\gamma \lambda_{\text{max}} + \text{trace}(\Lambda)}$$

• <u>Proof:</u> Convergence is guaranteed if the eigenvalues of matrix B are upper bounded by one. Since B is positive definite and symmetrical, all its eigenvalues are real-valued and positive.



• **Proof:** A sufficient condition is that the largest eigenvalue is smaller than one. This is the l₂ norm of the matrix B. An even loser condition is that the l₁ norm of B is smaller than one:

$$\lambda_{\max}^{(B)} = \|B\|_2 \le \|B\|_1$$

• Thus, for an arbitrary row of B, we have:

$$(1-\mu\lambda_i)^2 + \mu^2\lambda_i((\gamma-1)\lambda_i + \operatorname{trace}(\Lambda)) < 1$$

• From which we can follow for μ :

$$0 < \mu \le \frac{2}{\gamma \lambda_{\max} + \operatorname{trace}(\Lambda)} \le \frac{2}{\gamma \lambda_{i} + \operatorname{trace}(\Lambda)}$$

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(parameter vector) error

system mismatch:

$$E\left[\widetilde{\mathbf{w}}_{k}^{H}\widetilde{\mathbf{w}}_{k}\right] = \operatorname{trace}\left(E\left[\widetilde{\mathbf{w}}_{k}\widetilde{\mathbf{w}}_{k}^{H}\right]\right)$$

$$= \operatorname{trace}(P_{k})$$

$$= \underline{1}^{T} \underline{c}_{k} = \sum_{l=1}^{M} \gamma_{l} \lambda_{B,l}^{k}$$

Steady-state system mismatch:

$$\operatorname{trace}(P_{k}) = \sum_{l=1}^{M} \gamma_{l} \lambda_{B,l}^{k}$$

$$\lim_{k \to \infty} \underline{c}_{k} = \underline{c}_{\infty} = B\underline{c}_{\infty} + \mu^{2} \underline{\lambda} \sigma_{v}^{2}$$

$$= [I - B]^{-1} \mu^{2} \underline{\lambda} \sigma_{v}^{2}$$

$$= [2\Lambda - \mu\Lambda^{2}\gamma - \mu\underline{\lambda}\underline{\lambda}^{T}]^{-1} \mu\underline{\lambda} \sigma_{v}^{2}$$

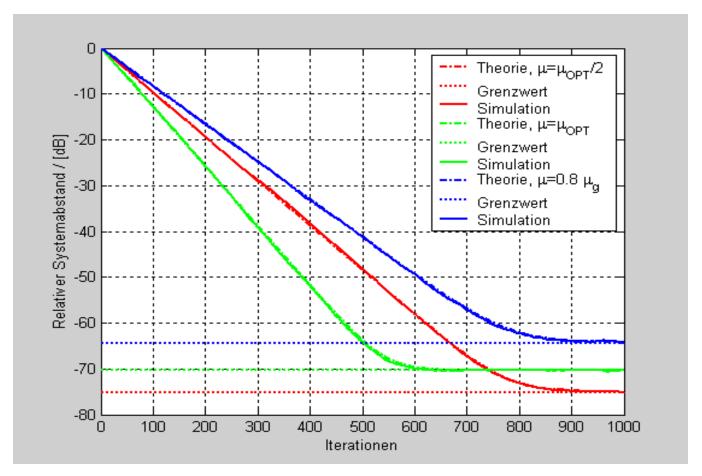
Matrix-Inversion-Lemma

$$\underline{1}^{T} \underline{c}_{\infty} = \mu \sigma_{\mathbf{v}}^{2} \frac{\sum_{l=1}^{M} \frac{1}{2 - \mu \gamma \lambda_{l}}}{1 - \mu \sum_{l=1}^{M} \frac{\lambda_{l}}{2 - \mu \gamma \lambda_{l}}}$$

Small step-sizes μ

$$\underline{1}^T \underline{c}_{\infty} = \frac{\mu \sigma_{\mathbf{v}}^2 M}{2}$$





Relative system mismatch

$$\frac{\underline{1}^T \underline{c}_{\infty}}{\|\underline{w}_o\|}$$

Distorted a-priori error:

$$\widetilde{\mathbf{e}}_{a,k} = \mathbf{d}_k - \underline{\mathbf{x}}_k^T \underline{\hat{\mathbf{w}}}_{k-1}$$

$$E\left[\left|\widetilde{\mathbf{e}}_{a,k}\right|^{2}\right] = E\left[\left|\mathbf{d}_{k} - \underline{\hat{\mathbf{w}}}_{k-1}^{T} \underline{\mathbf{x}}_{k}\right|^{2}\right]$$

$$= E\left[\left|\mathbf{v}_{k} - \underline{\widetilde{\mathbf{w}}}_{k-1}^{T} \underline{\mathbf{x}}_{k}\right|^{2}\right]$$

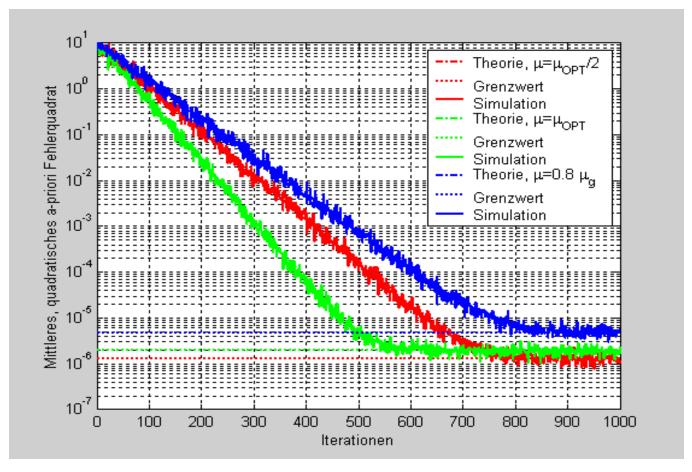
$$= \sigma_{\mathbf{v}}^{2} + E\left[\left|\underline{\widetilde{\mathbf{w}}}_{k-1}^{T} \underline{\mathbf{x}}_{k}\right|^{2}\right]$$

$$= \sigma_{\mathbf{v}}^{2} + E\left[\underline{\widetilde{\mathbf{w}}}_{k-1}^{T} \underline{\mathbf{x}}_{k} \underline{\mathbf{x}}_{k}^{H} \underline{\widetilde{\mathbf{w}}}_{k-1}^{*}\right]$$

$$= \sigma_{\mathbf{v}}^{2} + E\left[\underline{\widetilde{\mathbf{w}}}_{k-1}^{T} R_{\underline{\mathbf{x}}\underline{\mathbf{x}}}^{*} \underline{\widetilde{\mathbf{w}}}_{k-1}^{*}\right]$$

$$= \sigma_{\mathbf{v}}^{2} + \operatorname{trace}\left(E\left[R_{\underline{\mathbf{x}}\underline{\mathbf{x}}}^{*} \underline{\widetilde{\mathbf{w}}}_{k-1}^{*} \underline{\widetilde{\mathbf{w}}}_{k-1}^{T}\right]$$

$$= \sigma_{\mathbf{v}}^{2} + \underline{\lambda}^{T} \underline{c}_{k-1}$$



Excess Mean Square Error

$$g_{\text{ex}} = \underline{\lambda}^{T} \underline{c}_{\infty} = \mu \sigma_{\mathbf{v}}^{2} \frac{\sum_{l=1}^{M} \frac{\lambda_{l}}{2 - \mu \gamma \lambda_{l}}}{1 - \mu \sum_{l=1}^{M} \frac{\lambda_{l}}{2 - \mu \gamma \lambda_{l}}}$$

Misadjustment

$$m_{\mathrm{LMS}} = \frac{g_{\mathrm{ex}}}{g_o} = \mu \frac{\sum_{l=1}^{M} \frac{\lambda_l}{2 - \mu \gamma \lambda_l}}{1 - \mu \sum_{l=1}^{M} \frac{\lambda_l}{2 - \mu \gamma \lambda_l}}$$
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Learning curves

Note that there is some subtleties in averaging learning curves.

• If the outcome of an experiment is $||\mathbf{w}-\mathbf{w}_k||_2^2$ then averaging over N experiments is straightforward:

$$1/N \Sigma ||\underline{\mathbf{w}} - \underline{\mathbf{w}}_{k}||_{2}^{2}$$



Learning curves

- However, how do we average $||\underline{\mathbf{w}}-\underline{\mathbf{w}}_{k}||_{2}^{2}/||\underline{\mathbf{w}}||_{2}^{2}$
- The random variable <u>w</u> now appears in numerator and denominator!
- Jensens inequality: for convex functions:
- (for concave the other way around)

$$f(E[x]) \le E[f(x)]$$



Learning curves

- Therefore, we have to do the following:
- Nu= $1/N \Sigma ||\underline{\mathbf{w}} \underline{\mathbf{w}}_{k}||_{2}^{2}$
- De =1/N $\Sigma ||\underline{\mathbf{w}}||_2^2$
- Sys_{rel}(k)=Nu/De

