

$$\begin{aligned}
 \therefore \Phi_{\text{new}}(x)^T \Phi_{\text{new}}(y) &= \Phi(x)^T \Phi(y) \cdot \exp\left(\frac{-1}{2\sigma^2} x^T x\right) \\
 &\quad \cdot \exp\left(\frac{-1}{2\sigma^2} y^T y\right) \\
 &= \exp\left(\frac{x^T y}{\sigma^2}\right) \cdot \exp\left(\frac{-1}{2\sigma^2} x^T x\right) \\
 &\quad \cdot \exp\left(\frac{-1}{2\sigma^2} y^T y\right) \\
 &= \exp\left(\frac{-1}{2\sigma^2} (\|x - y\|^2)\right) \\
 &= K(x, y)
 \end{aligned}$$

Hence Proved, the Gaussian (rbf) kernel can be expressed in terms of $\Phi^T(x) \cdot \Phi(y)$.
Hence, rbf kernel is a valid kernel.

Q-1.7b). (i). Best $\sigma = 1$ (minimum errors at $\sigma=1$) decrease.

(ii) As we ~~decrease~~ the σ , we basically are shrinking the neighborhood of points for performing our prediction.

Upto $\sigma=1$, the errors decrease as we are shrinking the neighborhood to consider just a small meaningful set of neighbours.

But below $\sigma=1$, our errors increase as ~~we begin~~ our interval shrinks down tremendously and we begin to overfit on our data.

[Plot on the next page].