

## Lagrange interpolation: What nodes?

The setting: Lagrange interpolation in the  $xy$ -plane with nodes  $x_0 < x_1 < \dots < x_n$  and knots  $(x_i, y_i)$ ,  $i = 0:n$ . In many situations the knots are given as input data, and we have no choice. But in some cases we do have such freedom. A common such case is where you have a function subroutine, say `feval.m`, which returns  $y = f(x)$  given an input value  $x$  (in Matlab: `y = feval(x)`). Another common case is an experimental measurement, where the nodes are free to be chosen in the experiment's setup (e.g. "where do I place my sensors?").

Here we think about good nodes and bad nodes. To make things simple we assume  $x_j \in [-1, 1]$ , typically called the *unit interval*. If  $x_0 = -1$  and  $x_n = 1$ , we say the nodes are *closed* in  $[-1, 1]$ . If your nodes, say  $\hat{x}_j$ , are in the interval  $[a, b]$ , then we can linearly map from  $[-1, 1]$  to  $[a, b]$  by  $\hat{x}_j = (a + b + (b - a)x_j)/2$ .

Recall the Lagrange interpolator in Barycentric form

$$P(x) = \left( \sum_{i=0}^n \frac{y_i w_i}{x - x_i} \right) / \left( \sum_{i=0}^n \frac{w_i}{x - x_i} \right), \quad \text{where} \quad 1/w_j = \prod_{i=0, i \neq j}^n (x_j - x_i).$$

It turns out that equidistant nodes (a 'uniform grid') are a very bad choice for interpolation: the relative magnitudes of the  $w_j$  can differ by up to about  $2^n$ . With equidistant nodes, as  $n$  increases small changes in  $y$  (say  $y(1 + \epsilon)$ ) can give large (think  $\epsilon 2^n$ ) changes in  $P$ : *interpolation on a uniform grid is an ill-conditioned problem.*

Rounding errors alone are enough to give big spurious oscillations near the ends of the grid even for  $n < 20$ . We *will* use Lagrange interpolation on uniform grids in this course, but we will keep  $n \lesssim 5$ . For comparison later, here is the  $n=10$  closed uniform grid and  $w$ :

$$x = [-1, - .80, - .60, - .40, - .20, 0, .20, .40, .60, .80, 1]$$
$$w \approx [2.7, -27, 121, -323, 565, -678, 565, -323, 121, -269, 2.7]$$

This isn't true for all node choices for a given interval. There are (infinitely) many sets of nodes on the unit interval for which Lagrange interpolation is well conditioned, even for  $n$  in the thousands. These 'nice' nodes all cluster near the endpoints, and the relative magnitudes between the  $w_j$  are not large. Chebyshev nodes are a common example (we will use a closed set of Chebyshev nodes). These are the real parts of a closed uniform grid on the upper half of the unit circle:  $x_j = -\cos(j\pi/n)$ ,  $j = 0:n$ .

For example, if  $n = 10$  we have the 11 Chebyshev (type 2) nodes and  $w$ :

$$x \approx [-1, - .95, - .81, - .58, - .31, 0, .31, .58, .81, .95, 1]$$
$$w \approx [25.6, -51.2, -51.2, -51.2, -51.2, -51.2, -51.2, -51.2, -51.2, -51.2, 25.6]$$

If we think of Lagrange interpolation input data as an interval and a function  $f$  on that interval, then node selection is an important decision which can strongly effect the quality of the approximation. Choosing 'nice' nodes makes our approximation to  $f$  mostly depend on the smoothness of  $f$ . If  $f$  is smooth, then its frequency content (how 'wiggly' it is) puts a lower bound on  $n$  (see Nyquist frequency, or 'aliasing').