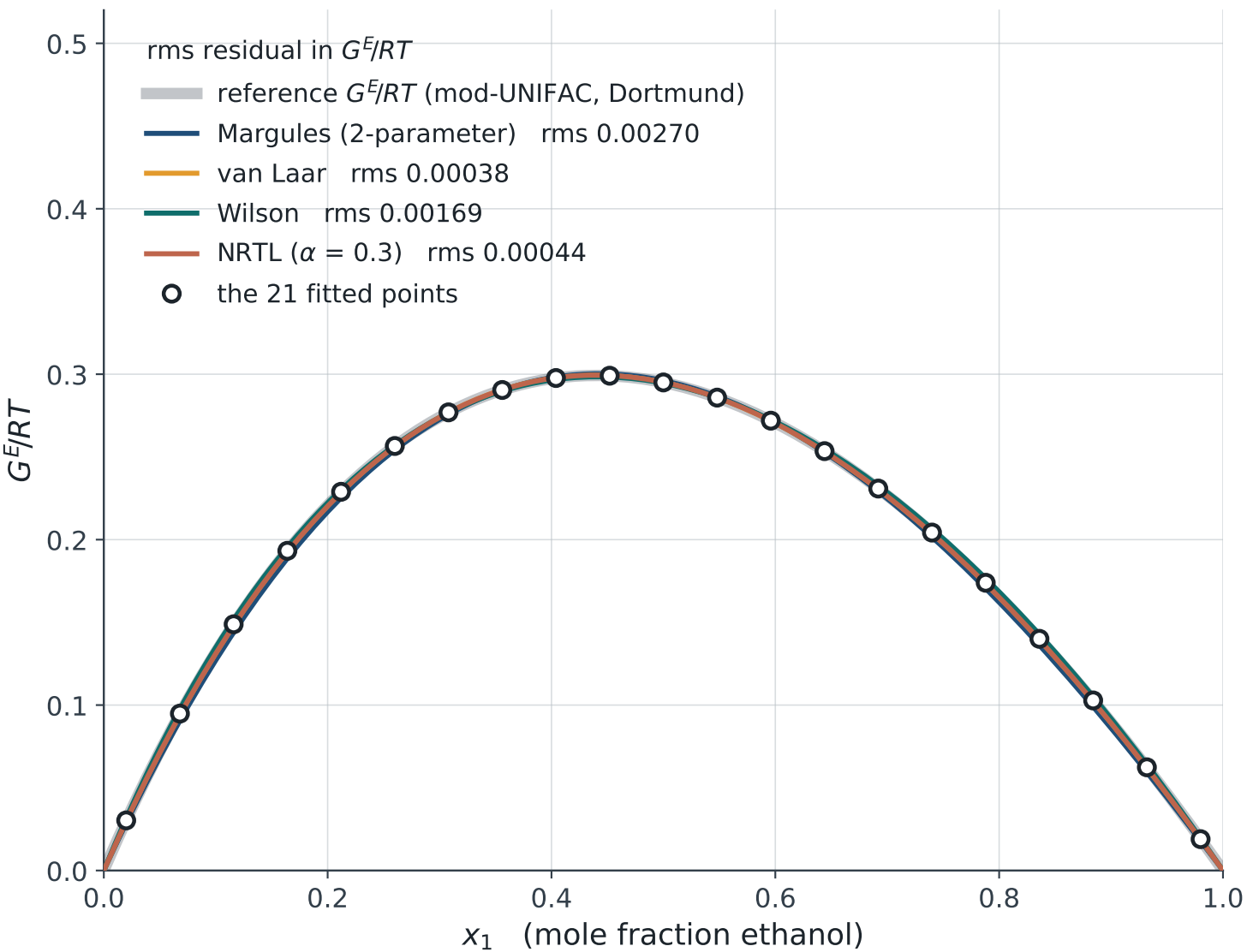


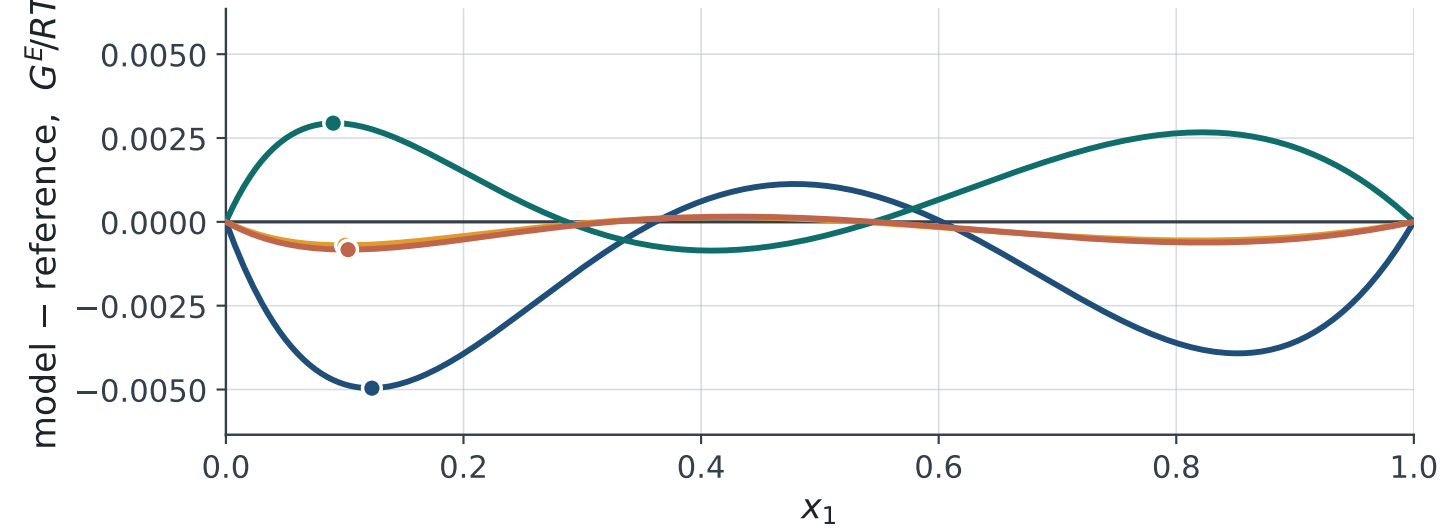
Four functional forms, one dataset: the overlay says they agree, the residual says they do not

ethanol (1) + water (2) at 298.15 K, all four fitted by `vlekit.fit` with Barker's method (objective 'P')

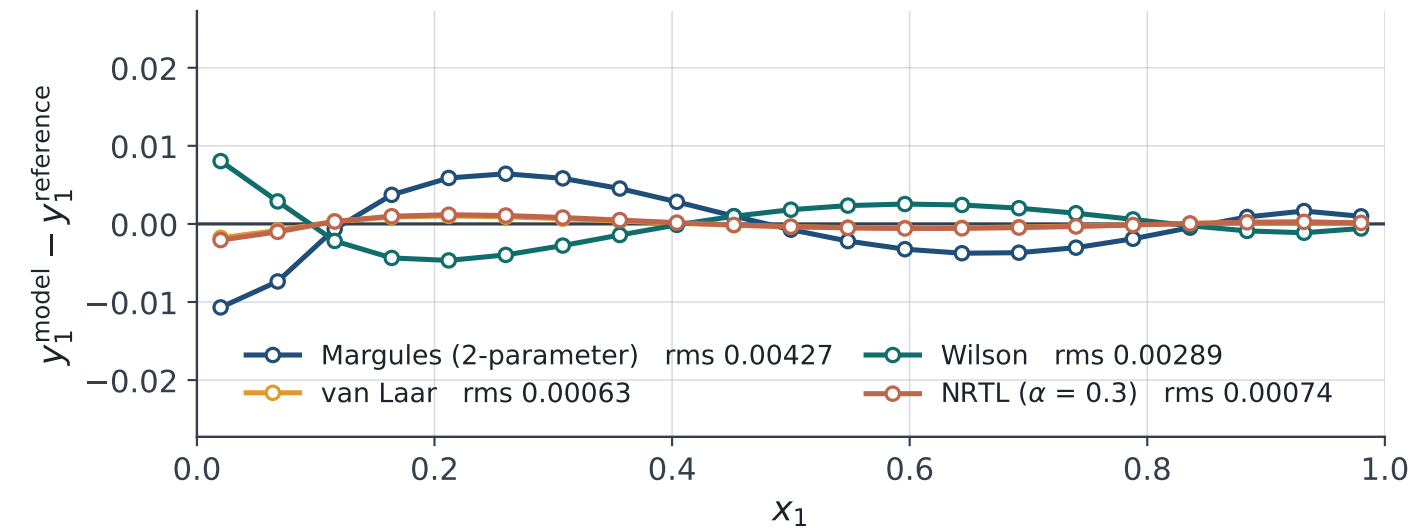
On this axis the four models are the same curve



the same four curves, reference subtracted



and the same four, in the vapour composition



THE DATASET IS NOT MEASURED AND CARRIES NO NOISE. It is generated by `vlekit.generate` from modified UNIFAC (Dortmund) at 298.15 K with an ideal vapour and Antoine constants from `vlekit.data.LIBRARY`, so every residual here is model-form error and nothing else; adding scatter would bury it. The rms values say how well each form can reproduce a realistic G^E , not how well it fits a particular laboratory.

Margules (2-parameter): $A_{12}=+1.4692$, $A_{21}=+0.8996$ | van Laar: $A=+1.5496$, $B=+0.9532$ | Wilson: $L_{12}=+0.2364$, $L_{21}=+0.7949$ | NRTL ($\alpha = 0.3$): $\tau_{12}=-0.0307$, $\tau_{21}=+1.5780$