

Table 3. Binding parameters of GIP₁₋₄₂ at different temperatures

Temperature (K)	K_d (μM)	Stoichiometry n	ΔH (kJ mol^{-1})	$-\text{T}\Delta S$ (kJ mol^{-1})	ΔG (kJ mol^{-1})
283.3	0.3 ± 0.1	1.0 ±	-82.5 ±	-47.3 ±	-35.1 ±
		0.1	0.6	3.6	3.0
288.3	0.6 ± 0.1	1.0 ±	-88.1 ±	-53.8 ±	-34.2 ±
		0.1	0.5	2.3	1.8
293.2	1.1 ± 0.1	1.0 ±	-92.7 ±	-59.2 ±	-33.5 ±
		0.1	0.4	1.5	1.1
298.2	2.2 ± 0.1	1.0 ±	-96.3 ±	-63.9 ±	-32.3 ±
		0.1	0.2	0.7	0.5
303.2	4.1 ± 0.1	1.0 ±	-104.8 ±	-73.5 ±	-31.3 ±
		0.1	0.3	0.7	0.4

Errors correspond to standard deviation of the nonlinear least-squares fit of the data points in the titration curve.