Supporting Information

Ellagitannins with Glucopyranose Cores Have Higher Affinities to Proteins than Acyclic Ellagitannins by Isothermal Titration Calorimetry

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- Figure S2. Examples of thermograms for the interactions of cyclic dimeric and trimeric ellagitannins with BSA (raw heat data, no control experiments subtracted): A) agrimoniin into 30 μ M BSA, B) gemin A into 30 μ M BSA, C) sanguiin H-6 into 30 μ M BSA and D) lambertianin C into 30 μ M BSA.
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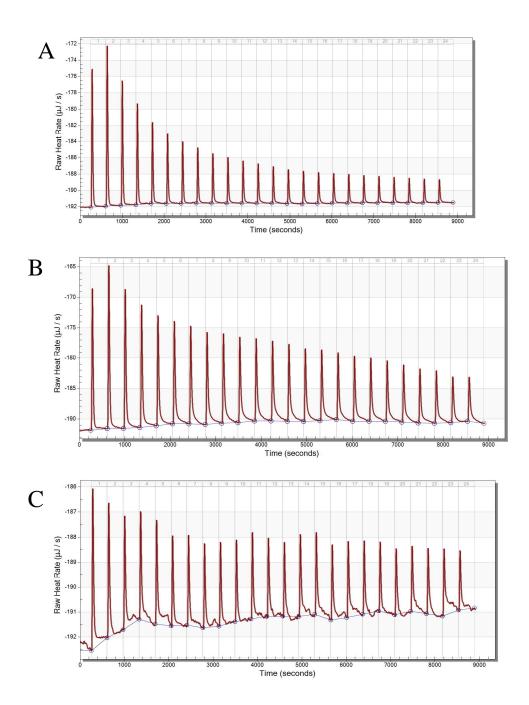


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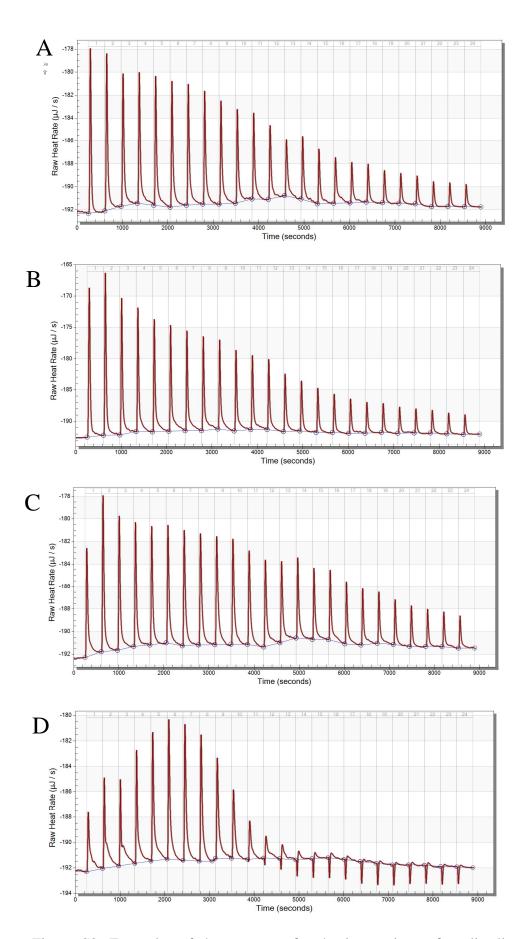


Figure S2. Examples of thermograms for the interactions of cyclic dimeric and trimeric ellagitannins with BSA (raw heat data, no control experiments subtracted): A) agrimoniin into 30 μM BSA, B) gemin A into 30 μM BSA, C) sanguiin H-6 into 30 μM BSA and D) lambertianin C into 30 μM BSA.

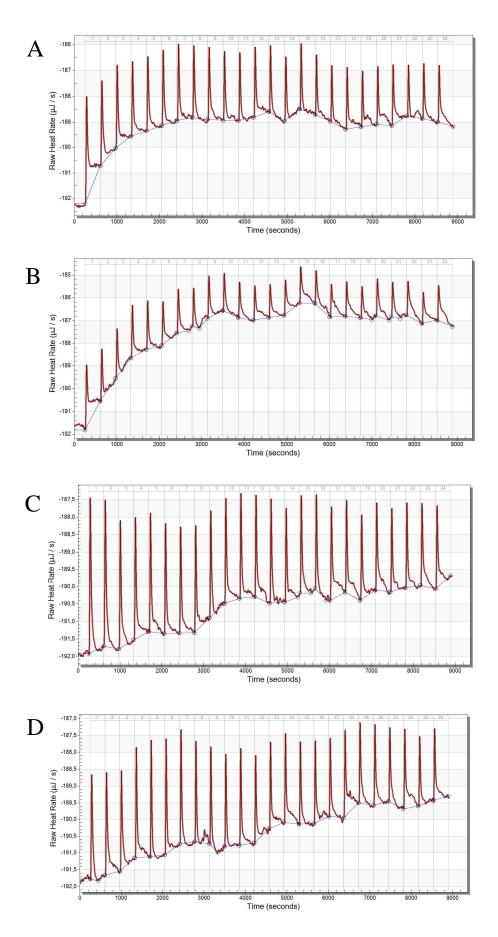


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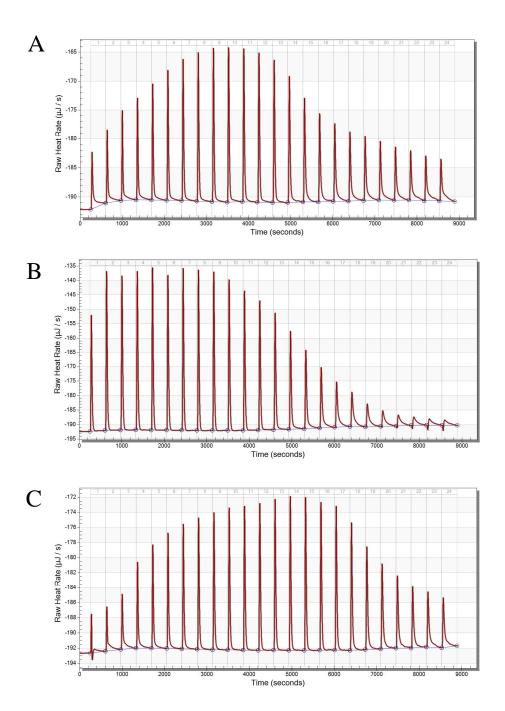


Figure S4. Examples of thermograms for the interactions of cyclic monomeric ellagitannins with gelatin (raw heat data, no control experiments subtracted): A) tellimagrandin I into 15 μM gelatin, B) tellimagrandin II into 20 μM gelatin and C) geraniin into 20 μM gelatin.

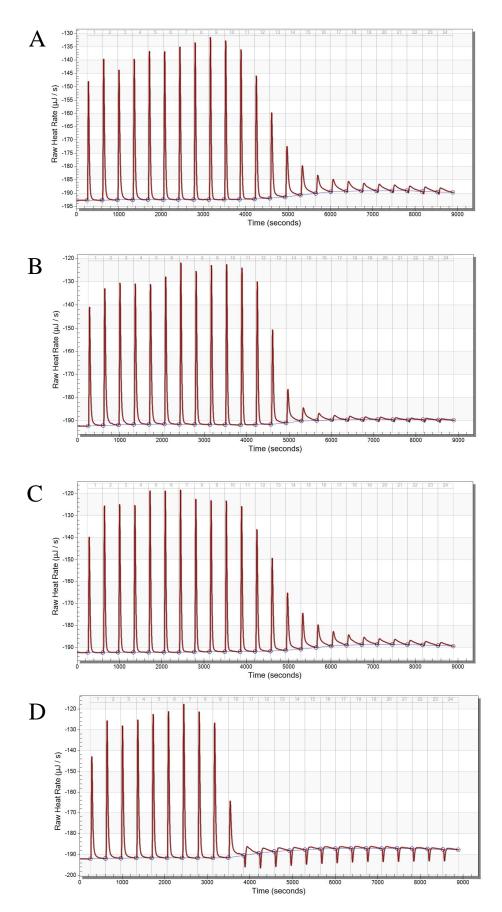


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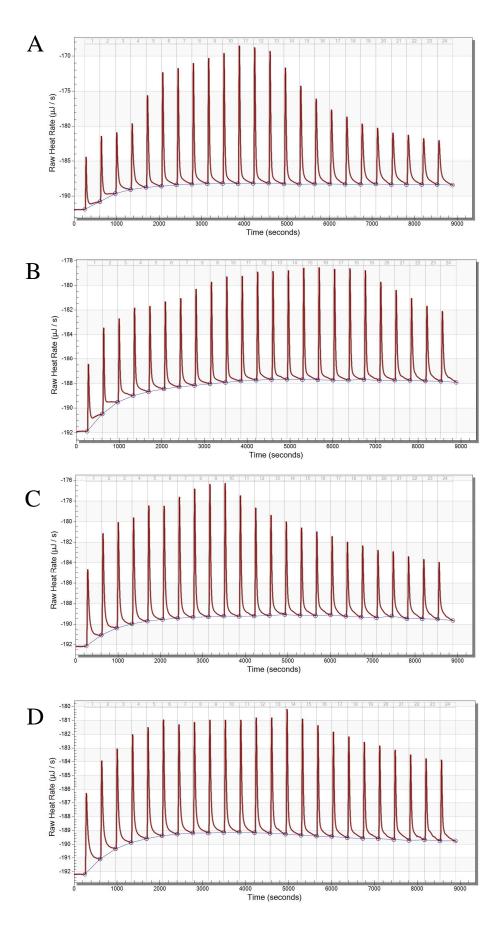


Figure S6. Examples of thermograms for the interactions of acyclic monomeric ellagitannins with gelatin (raw heat data, no control experiments subtracted): A) castalagin into 20 μ M gelatin, B) vescalagin into 20 μ M gelatin, C) castavaloninic acid into 20 μ M gelatin and D) vescavaloninic acid into 20 μ M gelatin.

Table S1. Estimated Entropies for the Interactions of Ellagitannins with BSA and Gelatin Fitted by Two-Site and One-Site Binding Models

	Tellimagrandin I	Tellimagrandin II	Agrimoniin	Gemin A	Sanguiin H-6	Roshenin C	Lambertianin C
BSA	1400		1000				
Two-Site							
ΔS ₁ (J mol ⁻¹ K ⁻¹)	15 ± 14	-38 ± 22	26 ± 17	-45 ± 2	28 ± 4		16 ± 7
ΔS ₂ (J mol ⁻¹ K ⁻¹)	18 ± 22	-64 ± 23	28 ± 4	29 ± 7	38 ± 7		46 ± 13
One-Site							
ΔS ₁ (J mol ⁻¹ K ⁻¹)	2 ± 24	-52 ± 25	1 ± 18	-73 ± 14	-8 ± 10		4 ± 7
Gelatin							
Two-Site							
ΔS ₁ (J mol ⁻¹ K ⁻¹)	53 ± 10	-93 ± 1	-117 ± 17	-99 ± 8	-108 ± 14	-2 ± 13	-185 ± 22
ΔS ₂ (J mol ⁻¹ K ⁻¹)	55 ± 8	64 ± 2	44 ± 7	32 ± 4	49 ± 2	31 ± 4	-28 ± 34
One-Site							
ΔS ₁ (J mol ⁻¹ K ⁻¹)	-57 ± 30	-103 ± 7	-163 ± 17	-165 ± 5	-170 ± 9	-112 ± 27	-194 ± 18