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Supplementary materials to: Effect of ion and binding site on the conformation of chosen glycosaminoglycans at the albumin surface

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Table S1. Binding energies in kcal/mol for 10 realizations of HSA-CS6 and HSA-HA complexes after docking.

n.o.	CS6	HA
1	4.356	5.416
2	4.354	4.984
3	4.348	4.945
4	4.347	4.715
5	4.339	4.69
6	4.322	4.688
7	4.314	4.687
8	4.308	4.684
9	4.274	4.635
10	4.249	4.62

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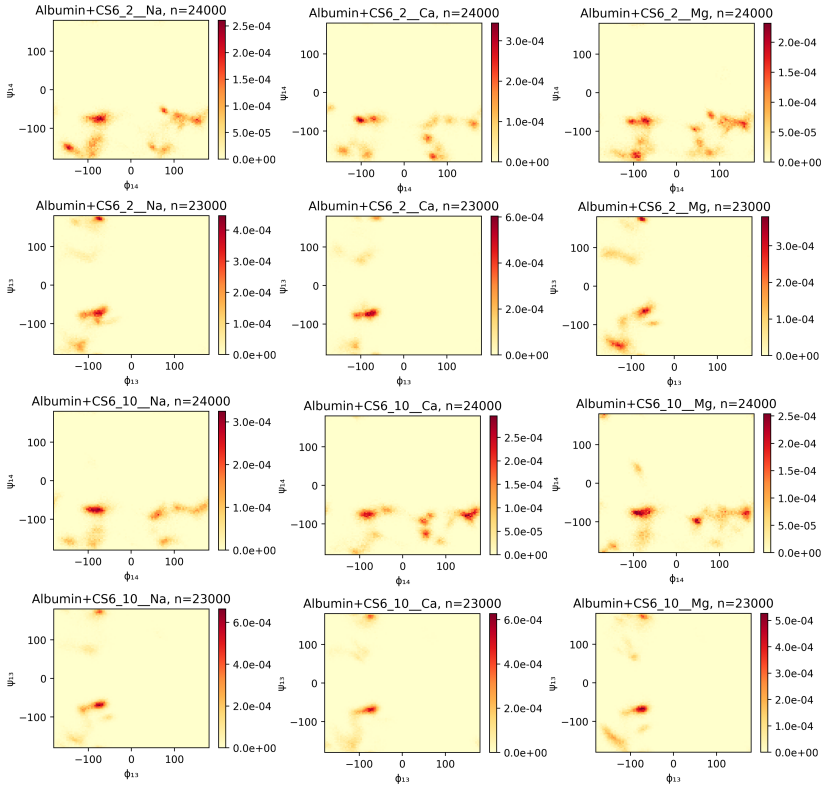

Figure S1. Normalized histograms for different realizations for HSA-CS6 complexes.

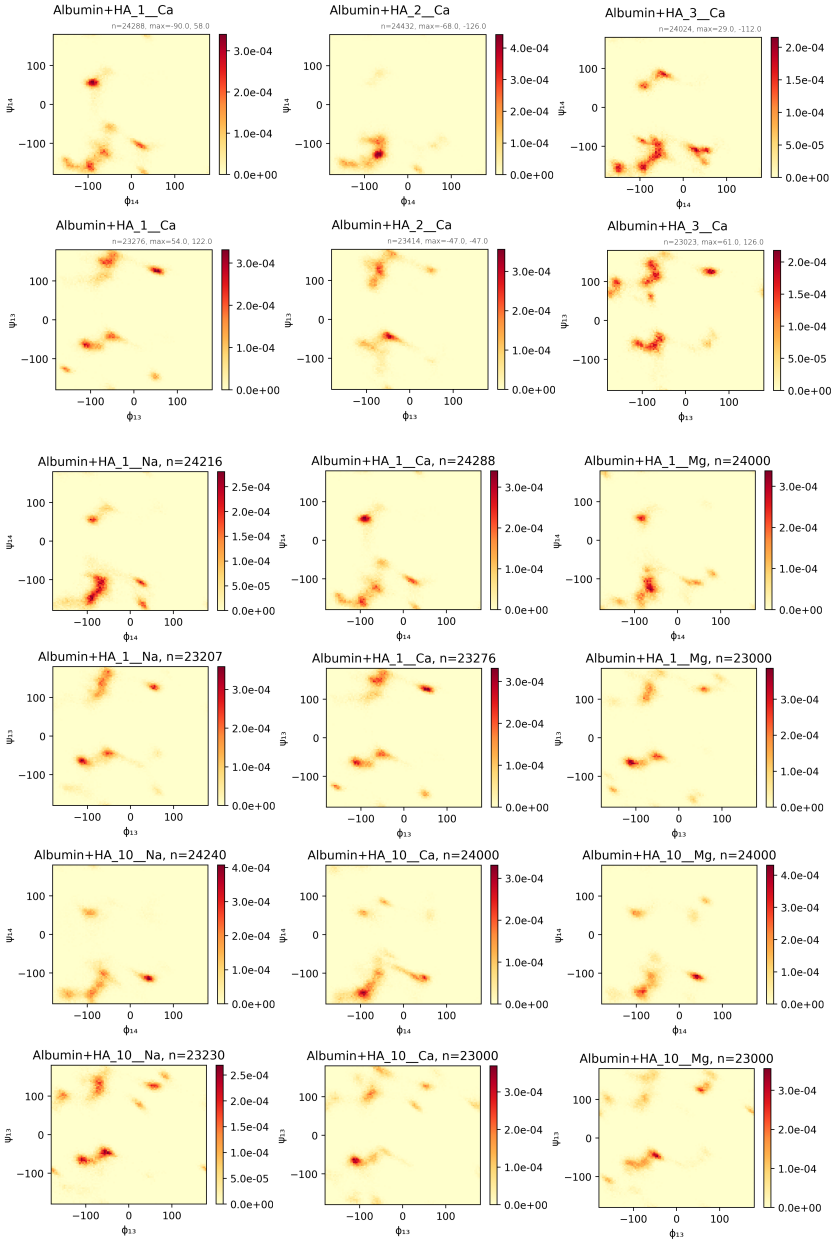

Figure S2. Normalized histograms for different realizations for HSA-HA complexes.