

Supplementary Materials

Topological Water Network Analysis Around Amino Acids

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Table S1. The number of TWNs observed around various atoms of all amino acids in the MD simulations. TWN analysis was carried out on the water molecules which were extracted every 10 ps for each simulated system. Amino acids are ordered from the most hydrophobic one (Ile, on the left hand side) to the most hydrophilic one (Arg, on the right hand side), according to the Kyte-Doolittle scale.

(A) 4-ring TWNs (Backbone + Side chain)

	Ile	Val	Leu	Phe	Cys	Met	Ala	Gly	Thr	Trp	Ser	Tyr	Pro	Hsd	Hse	Gln	Asp	Asn	Glu	Lys	Arg
O	61	52	68	70	69	73	75	66	101	66	112	73	55	67	56	78	105	105	71	79	
N	25	12	20	13	23	26	15	50	19	31	12	20	22	49	55	46	10	51	27	37	43
C	157	132	139	143	80	114	132	118	100	122	95	124	143	120	119	79	74	78	55	120	100
S					47	28															
Total	243	196	227	226	219	241	222	234	220	219	219	217	220	236	230	203	189	234	187	228	222
O,N/Total	0.35	0.33	0.39	0.37	0.42	0.41	0.41	0.50	0.55	0.44	0.57	0.43	0.35	0.49	0.48	0.61	0.61	0.67	0.71	0.47	0.55
C/Total	0.65	0.67	0.61	0.63	0.37	0.47	0.59	0.50	0.45	0.56	0.43	0.57	0.65	0.51	0.52	0.39	0.39	0.33	0.29	0.53	0.45
S/Total					0.21	0.12															

(B) 4-ring TWNs (Backbone)

	Ile	Val	Leu	Phe	Cys	Met	Ala	Gly	Thr	Trp	Ser	Tyr	Pro	Hsd	Hse	Gln	Asp	Asn	Glu	Lys	Arg
O	61	52	68	70	69	73	75	103	67	66	69	67	55	67	56	65	48	63	60	71	79
N	25	12	20	13	23	26	15	65	19	14	12	20	22	21	27	18	10	16	27	26	15
C	73	59	68	67	63	81	78	168	59	69	77	74	68	73	79	71	63	64	48	72	66
Total	159	123	156	150	155	180	168	336	145	149	158	161	145	161	162	154	121	143	135	169	160
O,N/Total	0.54	0.52	0.56	0.55	0.59	0.55	0.54	0.50	0.59	0.54	0.51	0.54	0.53	0.55	0.51	0.54	0.48	0.55	0.64	0.57	0.59
C/Total	0.46	0.48	0.44	0.45	0.41	0.45	0.46	0.50	0.41	0.46	0.49	0.46	0.47	0.45	0.49	0.46	0.52	0.45	0.36	0.43	0.41

(C) 4-ring TWNs (Side chain)

	Ile	Val	Leu	Phe	Cys	Met	Ala	Gly	Thr	Trp	Ser	Tyr	Pro	Hsd	Hse	Gln	Asp	Asn	Glu	Lys	Arg			
O											34		43	6			13	57	42	45				
N											17				28	28	28		35		11	28		
C	84	73	71	76	17	33	54			41	53	18	50	75	47	40	8	11	14	7	48	34		
S								47	28															
Total	84	73	71	76	64	61	54			75	70	61	56	75	75	68	49	68	91	52	59	62		
O,N/Total											0.45	0.24	0.70	0.11			0.37	0.41	0.84	0.84	0.85	0.87	0.19	0.45
C/Total	1.00	1.00	1.00	1.00	0.27	0.54	1.00			0.55	0.76	0.30	0.89	1.00	0.63	0.59	0.16	0.16	0.15	0.13	0.81	0.55		

(G) 6-ring TWNs (Backbone + Side chain)

(H) 6-ring TWNs (Backbone)

	Ile	Val	Leu	Phe	Cys	Met	Ala	Gly	Thr	Trp	Ser	Tyr	Pro	Hsd	Hse	Gln	Asp	Asn	Glu	Lys	Arg
O	5	5	9	7	9	15	10	2	7	3	9	9	4	5	5	10	2	4	3	10	7
N	0	2	3	2	3	2	4	7	4	2	0	1	1	1	2	1	0	5	4	1	2
C	11	9	5	2	5	8	6	5	7	5	4	4	3	3	8	4	5	2	7	3	5
Total	16	16	17	11	17	25	20	14	18	10	13	14	8	9	15	15	7	11	14	14	14
O/N/Total	0.31	0.44	0.71	0.82	0.71	0.68	0.70	0.64	0.61	0.50	0.69	0.71	0.63	0.67	0.47	0.73	0.29	0.82	0.50	0.79	0.64
C/Total	0.69	0.56	0.29	0.18	0.29	0.32	0.30	0.36	0.39	0.50	0.31	0.29	0.38	0.33	0.53	0.27	0.71	0.18	0.50	0.21	0.36

(I) 6-ring TWNs (Side chain)

Table S2. The number of TWNs observed around various atoms of all amino acids in the MD simulations. TWN analysis was carried out on the water molecules which were extracted every 5 ps for each simulated system. Amino acids are ordered from the most hydrophobic one (Ile, on the left hand side) to the most hydrophilic one (Arg, on the right hand side), according to the Kyte-Doolittle scale.

(A) 3-ring TWNs (Backbone + Side chain)

(B) 3-ring TWNs (Backbone)

	Ile	Val	Leu	Phe	Cys	Met	Ala	Gly	Thr	Trp	Ser	Tyr	Pro	Hsd	Hse	Gln	Asp	Asn	Glu	Lys	Arg
O	173	190	174	175	205	194	262	227	198	198	206	213	164	182	204	185	189	198	177	201	202
N	47	59	66	37	38	56	61	128	35	35	43	41	72	53	49	53	45	35	54	54	52
C	223	195	218	214	225	195	275	332	213	227	241	205	182	202	194	197	179	193	202	217	228
Total	443	444	458	426	468	445	598	687	446	460	490	459	418	437	447	435	413	426	433	472	482
O,N/Total	0.50	0.56	0.52	0.50	0.52	0.56	0.54	0.52	0.52	0.51	0.51	0.55	0.56	0.54	0.57	0.55	0.57	0.55	0.53	0.54	0.53
C/Total	0.50	0.44	0.48	0.50	0.48	0.44	0.46	0.48	0.48	0.49	0.49	0.45	0.44	0.46	0.43	0.45	0.43	0.45	0.47	0.46	0.47

(C) 3-ring TWNs (Side chain)

(D) 4-ring TWNs (Backbone + Side chain)

(E) 4-ring TWNs (Backbone)

	Ile	Val	Leu	Phe	Cys	Met	Ala	Gly	Thr	Trp	Ser	Tyr	Pro	Hsd	Hse	Gln	Asp	Asn	Glu	Lys	Arg	
O	117	106	138	131	132	146	158	148	124	141	140	137	110	133	124	114	107	118	139	143	143	
N	44	25	36	41	44	45	41	109	34	32	31	42	49	39	44	44	25	35	42	42	39	
C	135	126	138	136	134	145	156	220	104	140	140	134	129	133	134	138	122	134	105	138	132	
Total	296	257	312	308	310	336	355	477	262	313	311	313	288	305	302	296	254	287	286	323	314	
O,N/Total	0.54	0.51	0.56	0.56	0.57	0.57	0.56	0.54	0.60	0.55	0.55	0.57	0.55	0.56	0.56	0.53	0.52	0.53	0.63	0.57	0.58	
C/Total	0.46	0.49	0.44	0.44	0.43	0.43	0.44	0.46	0.40	0.45	0.45	0.43	0.45	0.44	0.44	0.47	0.48	0.47	0.47	0.37	0.43	0.42

(F) 4-ring TWNs (Side chain)

(G) 5-ring TWNs (Backbone + Side chain)

(H) 5-ring TWNs (Backbone)

	Ile	Val	Leu	Phe	Cys	Met	Ala	Gly	Thr	Trp	Ser	Tyr	Pro	Hsd	Hse	Gln	Asp	Asn	Glu	Lys	Arg
O	52	56	40	49	53	61	49	58	45	54	56	43	45	53	62	56	39	50	32	52	59
N	13	9	22	16	18	24	11	35	18	17	10	15	9	14	17	9	9	14	10	18	19
C	46	46	53	40	55	44	60	73	37	34	30	37	43	38	44	39	26	38	45	45	45
Total	111	111	115	105	126	129	120	166	100	105	96	95	97	105	123	104	74	102	87	115	123
O/N/Total	0.59	0.59	0.54	0.62	0.56	0.66	0.50	0.56	0.63	0.68	0.69	0.61	0.56	0.64	0.64	0.63	0.65	0.63	0.48	0.61	0.63
C/Total	0.41	0.41	0.46	0.38	0.44	0.34	0.50	0.44	0.37	0.32	0.31	0.39	0.44	0.36	0.36	0.38	0.35	0.37	0.52	0.39	0.37

(I) 5-ring TWNs (Side chain)

(J) 6-ring TWNs (Backbone + Side chain)

(K) 6-ring TWNs (Backbone)

	Ile	Val	Leu	Phe	Cys	Met	Ala	Gly	Thr	Trp	Ser	Tyr	Pro	Hsd	Hse	Gln	Asp	Asn	Glu	Lys	Arg
O	11	11	19	12	23	23	19	8	9	11	19	16	8	11	10	17	6	8	13	14	11
N	3	3	4	2	5	5	4	13	6	7	3	4	3	1	5	3	3	8	6	4	5
C	13	12	11	5	8	14	10	16	11	8	10	8	8	5	9	11	11	6	11	9	11
Total	27	26	34	19	36	42	33	37	26	26	32	28	19	17	24	31	20	22	30	27	27
O,N/Total	0.52	0.54	0.68	0.74	0.78	0.67	0.70	0.57	0.58	0.69	0.69	0.71	0.58	0.71	0.63	0.65	0.45	0.73	0.63	0.67	0.59
C/Total	0.48	0.46	0.32	0.26	0.22	0.33	0.30	0.43	0.42	0.31	0.31	0.29	0.42	0.29	0.38	0.35	0.55	0.27	0.37	0.33	0.41

(L) 6-ring TWNs (Side chain)

Table S3. The number of TWNs observed around various atoms of all amino acids in the PDBs. Amino acids are ordered from the most hydrophobic one (Ile, on the left hand side) to the most hydrophilic one (Arg, on the right hand side), according to the Kyte-Doolittle scale.

(A) 4-ring TWNs (Backbone + Side chain)

(B) 4-ring TWNs (Backbone)

(C) 4-ring TWNs (Side chain)

	Ile	Val	Leu	Phe	Cys	Met	Ala	Gly	Thr	Trp	Ser	Tyr	Pro	His	Gln	Asp	Asn	Glu	Lys	Arg	
O									1965		1994	1402			1429	5916	1393	5480			
N										78				755	670		669		1583	1791	
C	114	151	199	98	2	81	125		156	45	44	82	243	163	80	52	41	113	251	167	
S						45	30														
Total	114	151	199	98	47	111	125		2121	123	2038	1484	243	918	2179	5968	2103	5593	1834	1958	
O,N/									0.93	0.63	0.98	0.94			0.82	0.96	0.99	0.98	0.98	0.86	0.91
Total																					
C/Total	1.00	1.00	1.00	1.00	0.04	0.73	1.00		0.07	0.37	0.02	0.06	1.00	0.18	0.04	0.01	0.02	0.02	0.14	0.09	

(E) 5-ring TWNs (Backbone)																				
	Ile	Val	Leu	Phe	Cys	Met	Ala	Gly	Thr	Trp	Ser	Tyr	Pro	His	Gln	Asp	Asn	Glu	Lys	Arg
O	307	447	491	227	53	123	743	669	442	118	401	234	564	169	248	471	319	454	478	397
N	31	34	79	23	2	10	89	81	36	3	47	15	5	23	33	46	66	51	41	32
C	1	0	1	2	0	3	11	22	3	2	8	0	7	6	2	3	3	3	1	2
Total	339	481	571	252	55	136	843	772	481	123	456	249	576	198	283	520	388	508	520	431
O,N/ Total	1.00	1.00	1.00	0.99	1.00	0.98	0.99	0.97	0.99	0.98	0.98	1.00	0.99	0.97	0.99	0.99	0.99	0.99	1.00	1.00
C/Total	0.00	0.00	0.00	0.01	0.00	0.02	0.01	0.03	0.01	0.02	0.02	0.00	0.01	0.03	0.01	0.01	0.01	0.01	0.00	0.00

(G) 6-ring TWNs (Backbone + Side chain)

(H) 6-ring TWNs (Backbone)

	Ile	Val	Leu	Phe	Cys	Met	Ala	Gly	Thr	Trp	Ser	Tyr	Pro	His	Gln	Asp	Asn	Glu	Lys	Arg
O	105	176	173	157	67	36	324	312	220	47	148	92	148	48	96	146	139	140	115	127
N	9	17	17	2	0	3	44	43	12	17	19	11	4	11	17	14	107	12	7	7
C	0	0	0	0	0	2	1	10	1	1	1	0	1	0	1	0	2	1	0	0
Total	114	193	190	159	67	41	369	365	233	65	168	103	153	59	114	160	248	153	122	134
O,N/ Total	1.00	1.00	1.00	1.00	1.00	0.95	1.00	0.97	1.00	0.98	0.99	1.00	0.99	1.00	0.99	1.00	0.99	1.00	1.00	1.00
C/Total	0.00	0.00	0.00	0.00	0.00	0.05	0.00	0.03	0.00	0.02	0.01	0.00	0.01	0.00	0.01	0.00	0.01	0.01	0.00	0.00

(I) 6-ring TWNs (Side chain)

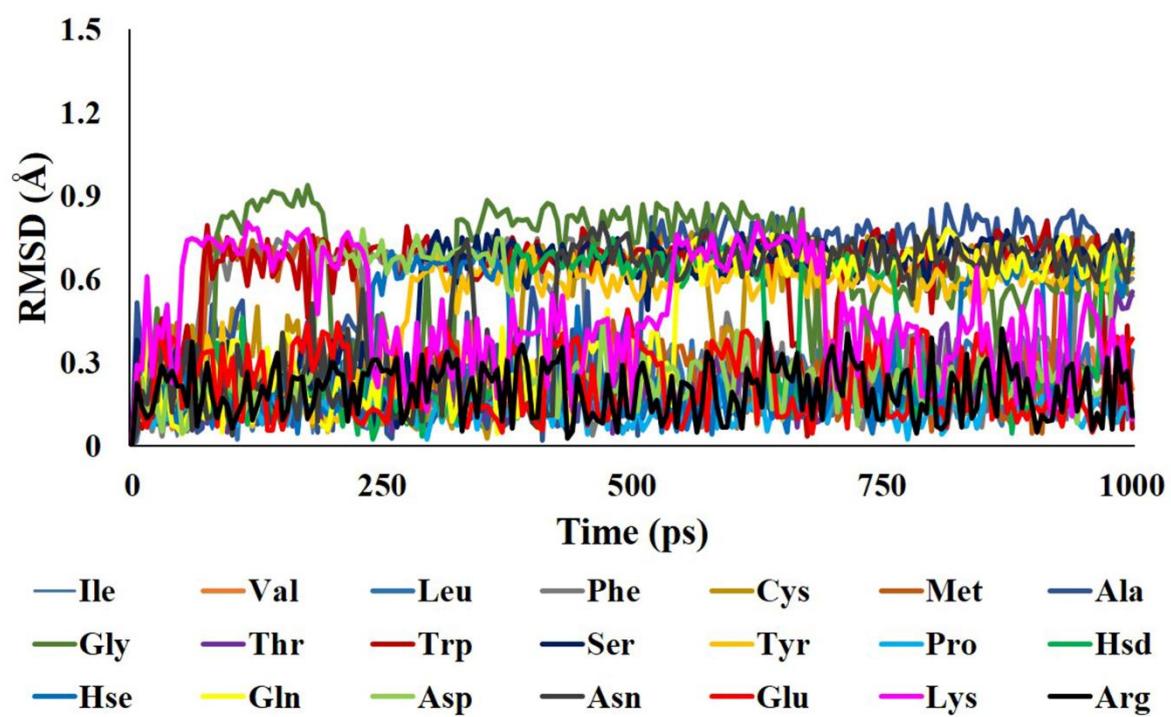


Figure S1. RMSD plot for the backbone atoms of amino acids from the initial structures throughout the 1 ns MD simulation as a function of time.

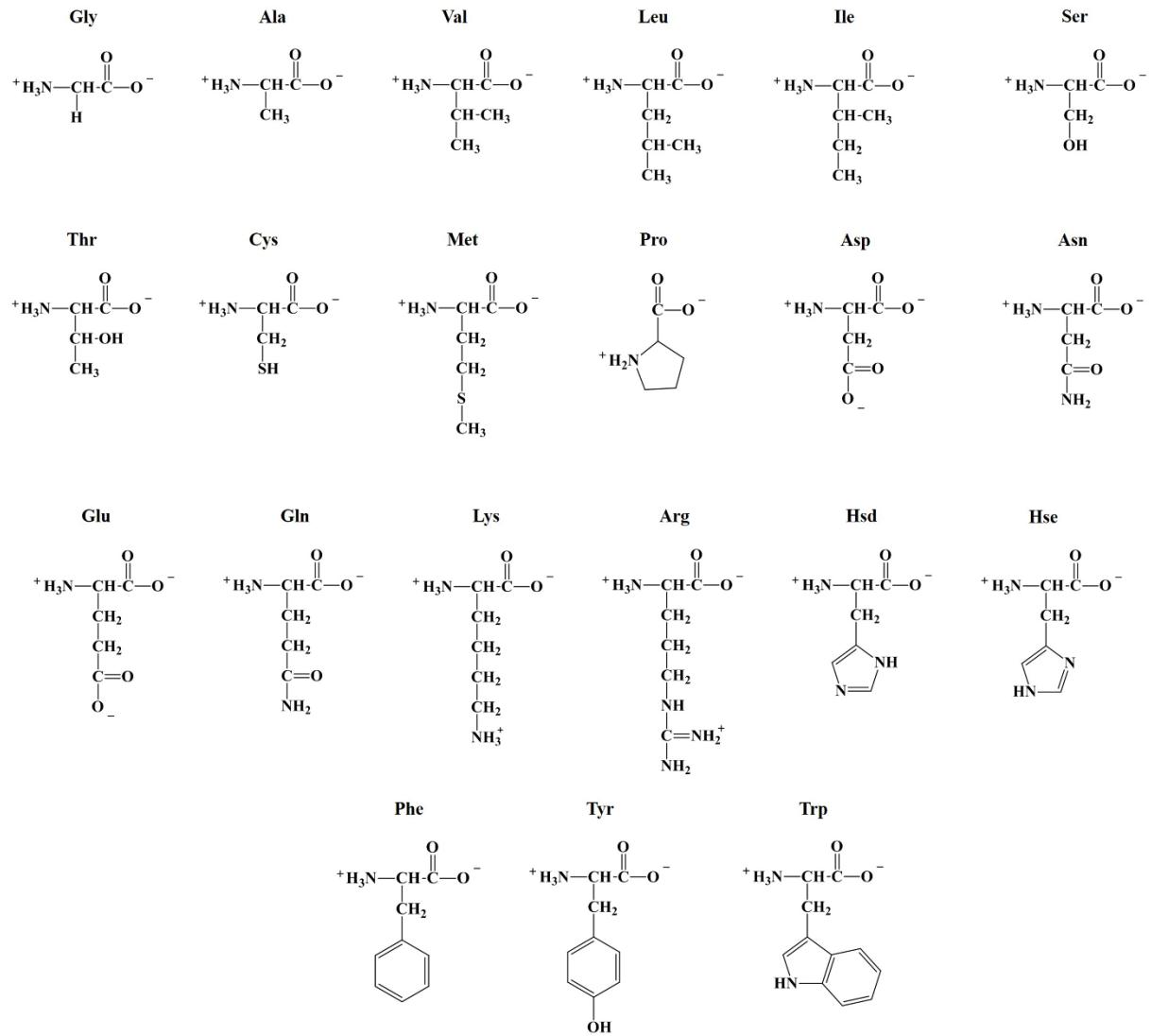


Figure S2. Structures of the amino acids studied in the present work.