



wwPDB EM Validation Summary Report ⓘ

Mar 25, 2026 – 06:40 AM UTC

PDB ID : 8J9H / pdb_00008j9h
EMDB ID : EMD-36107
Title : Cryo-EM structure of Euglena gracilis respiratory complex I, deactive state
Authors : Wu, M.C.; He, Z.X.; Tian, H.T.; Hu, Y.Q.; Han, F.Z.; Zhou, L.
Deposited on : 2023-05-03
Resolution : 3.11 Å(reported)

This is a wwPDB EM Validation Summary Report for a publicly released PDB entry.

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A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

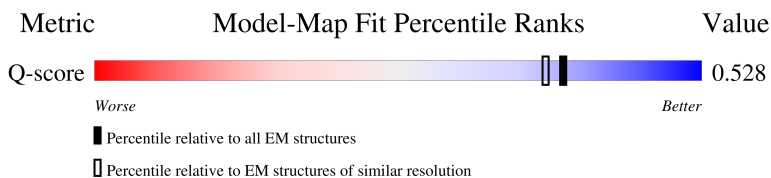
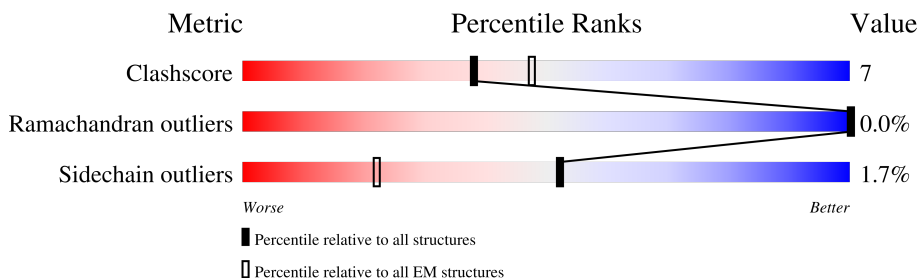
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 3.11 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14465 (2.61 - 3.61)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	1A	385	
2	1B	527	
3	2B	142	
4	4L	171	

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Mol	Chain	Length	Quality of chain
5	A1	141	
6	A2	193	
7	A3	125	
8	A5	184	
9	A6	437	
10	A7	136	
11	A8	223	
12	A9	489	
13	AB	134	
14	AL	281	
15	AM	198	
16	AN	287	
17	B2	145	
18	B3	62	
19	B4	171	
20	B5	140	
21	B6	91	
22	B7	97	
23	B8	176	
24	B9	158	
25	BL	144	
26	BM	112	
27	C4	185	
28	E1	483	
29	E2	467	

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Mol	Chain	Length	Quality of chain
30	E3	434	
31	E4	368	
32	E5	290	
33	E6	371	
34	E8	205	
35	EA	126	
36	EB	101	
37	EC	101	
38	ED	151	
39	FX	325	
40	G1	436	
41	G2	267	
42	G3	261	
43	N1	670	
44	N2	300	
45	N3	293	
45	N6	293	
46	N4	478	
47	N5	584	
48	S2	395	
49	S3	277	
50	S4	208	
51	S5	122	
52	S6	147	
53	S7	207	

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Mol	Chain	Length	Quality of chain
54	S8	212	
55	U1	12	
55	U2	12	
56	V1	526	
57	V2	225	
58	E7	246	
59	AC	134	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
61	SF4	S8	297	-	-	X	-
61	SF4	S8	298	-	-	X	-

2 Entry composition

There are 70 unique types of molecules in this entry. The entry contains 226135 atoms, of which 112544 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called NDUS1A.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
1	1A	352	5501	1753	2700	488	537	23	0	0

- Molecule 2 is a protein called NDUS1B.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
2	1B	525	8357	2679	4159	743	765	11	1	0

- Molecule 3 is a protein called NADH dehydrogenase subunit 5.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
3	2B	140	2059	712	989	172	183	3	0	0

- Molecule 4 is a protein called ND4L.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
4	4L	108	1768	606	878	133	145	6	0	0

- Molecule 5 is a protein called NDUFA1.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
5	A1	137	2097	684	1026	192	192	3	0	0

- Molecule 6 is a protein called NDUFA2.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	S		
6	A2	192	2967	942	1474	267	280	4	0	0

- Molecule 7 is a protein called NDUFA3.

Mol	Chain	Residues	Atoms						AltConf	Trace
7	A3	124	Total	C	H	N	O	S	0	0
			2089	678	1039	191	175	6		

- Molecule 8 is a protein called NDUFA5.

Mol	Chain	Residues	Atoms						AltConf	Trace
8	A5	154	Total	C	H	N	O	S	0	0
			2509	794	1248	221	244	2		

- Molecule 9 is a protein called NDUFA6.

Mol	Chain	Residues	Atoms						AltConf	Trace
9	A6	423	Total	C	H	N	O	S	0	0
			6608	2091	3280	601	632	4		

- Molecule 10 is a protein called NDUFA7.

Mol	Chain	Residues	Atoms						AltConf	Trace
10	A7	136	Total	C	H	N	O	S	0	0
			2272	735	1118	219	194	6		

- Molecule 11 is a protein called NDUFA8.

Mol	Chain	Residues	Atoms						AltConf	Trace
11	A8	223	Total	C	H	N	O	S	0	0
			3548	1160	1726	315	334	13		

- Molecule 12 is a protein called NDUFA9.

Mol	Chain	Residues	Atoms						AltConf	Trace
12	A9	484	Total	C	H	N	O	S	0	0
			7679	2449	3850	662	700	18		

- Molecule 13 is a protein called NDUFAB1-alpha.

Mol	Chain	Residues	Atoms						AltConf	Trace
13	AB	88	Total	C	H	N	O	S	0	0
			1367	437	673	114	139	4		

- Molecule 14 is a protein called NDUFA12.

Mol	Chain	Residues	Atoms						AltConf	Trace
14	AL	265	Total	C	H	N	O	S	0	0
			4409	1439	2172	414	379	5		

- Molecule 15 is a protein called NDUFA13.

Mol	Chain	Residues	Atoms						AltConf	Trace
15	AM	184	Total	C	H	N	O	S	0	0
			2935	953	1448	264	263	7		

- Molecule 16 is a protein called NDUFA11.

Mol	Chain	Residues	Atoms						AltConf	Trace
16	AN	287	Total	C	H	N	O	S	0	0
			4573	1501	2267	396	399	10		

- Molecule 17 is a protein called NDUFB2.

Mol	Chain	Residues	Atoms						AltConf	Trace
17	B2	105	Total	C	H	N	O	S	0	0
			1770	604	857	142	166	1		

- Molecule 18 is a protein called NDUFB3.

Mol	Chain	Residues	Atoms						AltConf	Trace
18	B3	61	Total	C	H	N	O	S	0	0
			758	292	309	88	68	1		

- Molecule 19 is a protein called NDUFB4.

Mol	Chain	Residues	Atoms						AltConf	Trace
19	B4	171	Total	C	H	N	O	S	0	0
			2735	885	1358	250	236	6		

- Molecule 20 is a protein called NDUFB5.

Mol	Chain	Residues	Atoms						AltConf	Trace
20	B5	140	Total	C	H	N	O	S	0	0
			2181	708	1069	207	195	2		

- Molecule 21 is a protein called NDUFB6.

Mol	Chain	Residues	Atoms						AltConf	Trace
21	B6	91	Total	C	H	N	O	S	0	0
			1520	509	747	132	128	4		

- Molecule 22 is a protein called NDUFB7.

Mol	Chain	Residues	Atoms						AltConf	Trace
22	B7	97	Total	C	H	N	O	S	0	0
			1692	536	835	165	149	7		

- Molecule 23 is a protein called NDUFB8.

Mol	Chain	Residues	Atoms						AltConf	Trace
23	B8	147	Total	C	H	N	O	S	0	0
			2351	804	1127	199	213	8		

- Molecule 24 is a protein called NDUFB9.

Mol	Chain	Residues	Atoms						AltConf	Trace
24	B9	151	Total	C	H	N	O	S	0	0
			2443	795	1207	216	222	3		

- Molecule 25 is a protein called NDUFB10.

Mol	Chain	Residues	Atoms						AltConf	Trace
25	BL	144	Total	C	H	N	O	S	0	0
			2406	786	1179	215	216	10		

- Molecule 26 is a protein called NDUFB11.

Mol	Chain	Residues	Atoms						AltConf	Trace
26	BM	112	Total	C	H	N	O	S	0	0
			1737	577	827	164	167	2		

- Molecule 27 is a protein called NDUF2.

Mol	Chain	Residues	Atoms						AltConf	Trace
27	C4	183	Total	C	H	N	O	S	0	0
			3062	1000	1517	268	271	6		

- Molecule 28 is a protein called NDUEG1.

Mol	Chain	Residues	Atoms						AltConf	Trace
28	E1	450	Total	C	H	N	O	S	0	0
			7008	2244	3496	601	654	13		

- Molecule 29 is a protein called NDUEG2.

Mol	Chain	Residues	Atoms						AltConf	Trace
29	E2	466	Total	C	H	N	O	S	0	0
			7103	2286	3540	618	655	4		

- Molecule 30 is a protein called NDUEG3.

Mol	Chain	Residues	Atoms						AltConf	Trace
30	E3	432	Total	C	H	N	O	S	0	0
			6518	2071	3263	565	612	7		

- Molecule 31 is a protein called NDUEG4.

Mol	Chain	Residues	Atoms						AltConf	Trace
31	E4	351	Total	C	H	N	O	S	0	0
			5502	1774	2732	477	504	15		

- Molecule 32 is a protein called NDUEG5.

Mol	Chain	Residues	Atoms						AltConf	Trace
32	E5	276	Total	C	H	N	O	S	0	0
			4046	1265	2069	341	369	2		

- Molecule 33 is a protein called NDUEG6.

Mol	Chain	Residues	Atoms						AltConf	Trace
33	E6	318	Total	C	H	N	O	S	0	0
			5228	1703	2554	477	482	12		

- Molecule 34 is a protein called NDUEG8.

Mol	Chain	Residues	Atoms						AltConf	Trace
34	E8	205	Total	C	H	N	O	S	0	0
			3354	1100	1663	288	292	11		

- Molecule 35 is a protein called NDUEG10.

Mol	Chain	Residues	Atoms						AltConf	Trace
35	EA	124	Total	C	H	N	O	S	0	0
			1793	630	832	172	156	3		

- Molecule 36 is a protein called NDUEG11.

Mol	Chain	Residues	Atoms						AltConf	Trace
36	EB	101	Total	C	H	N	O	S	0	0
			1405	473	631	150	144	7		

- Molecule 37 is a protein called NDUEG12.

Mol	Chain	Residues	Atoms						AltConf	Trace
37	EC	85	Total	C	H	N	O	S	0	0
			1323	424	663	116	118	2		

- Molecule 38 is a protein called NDUEG13.

Mol	Chain	Residues	Atoms						AltConf	Trace
38	ED	138	Total	C	H	N	O	S	0	0
			2273	736	1131	205	196	5		

- Molecule 39 is a protein called NDUFX.

Mol	Chain	Residues	Atoms						AltConf	Trace
39	FX	237	Total	C	H	N	O	S	0	0
			3816	1263	1849	338	359	7		

- Molecule 40 is a protein called NDUCA1.

Mol	Chain	Residues	Atoms						AltConf	Trace
40	G1	403	Total	C	H	N	O	S	0	0
			6146	1979	2999	558	594	16		

- Molecule 41 is a protein called NDUCA2.

Mol	Chain	Residues	Atoms						AltConf	Trace
41	G2	236	Total	C	H	N	O	S	0	0
			3650	1138	1846	323	338	5		

- Molecule 42 is a protein called NDUCA3.

Mol	Chain	Residues	Atoms						AltConf	Trace
42	G3	261	Total	C	H	N	O	S	0	0
			3905	1226	1944	356	373	6		

- Molecule 43 is a protein called ND1.

Mol	Chain	Residues	Atoms						AltConf	Trace
43	N1	310	Total	C	H	N	O	S	0	0
			5331	1783	2726	380	435	7		

- Molecule 44 is a protein called ND2A.

Mol	Chain	Residues	Atoms						AltConf	Trace
44	N2	296	Total	C	H	N	O	S	0	0
			5101	1725	2589	362	418	7		

- Molecule 45 is a protein called ND3.

Mol	Chain	Residues	Atoms						AltConf	Trace
45	N3	121	Total	C	H	N	O	S	0	0
			2094	720	1057	143	172	2		
45	N6	154	Total	C	H	N	O	S	0	0
			2642	857	1385	187	210	3		

- Molecule 46 is a protein called NADH-ubiquinone oxidoreductase chain 4.

Mol	Chain	Residues	Atoms						AltConf	Trace
46	N4	478	Total	C	H	N	O	S	0	0
			8215	2743	4214	582	663	13		

- Molecule 47 is a protein called ND5.

Mol	Chain	Residues	Atoms						AltConf	Trace
47	N5	584	Total	C	H	N	O	S	0	0
			9869	3293	5032	711	808	25		

- Molecule 48 is a protein called NDUFS2.

Mol	Chain	Residues	Atoms						AltConf	Trace
48	S2	394	Total	C	H	N	O	S	0	0
			6274	2041	3101	541	569	22		

- Molecule 49 is a protein called NDUFS3.

Mol	Chain	Residues	Atoms						AltConf	Trace
49	S3	248	Total	C	H	N	O	S	0	0
			3978	1307	1928	346	384	13		

- Molecule 50 is a protein called NDUF54.

Mol	Chain	Residues	Atoms						AltConf	Trace
50	S4	190	Total	C	H	N	O	S	0	0
			3038	956	1502	300	273	7		

- Molecule 51 is a protein called NDUF55.

Mol	Chain	Residues	Atoms						AltConf	Trace
51	S5	122	Total	C	H	N	O	S	0	0
			1886	625	895	173	188	5		

- Molecule 52 is a protein called NDUF56.

Mol	Chain	Residues	Atoms						AltConf	Trace
52	S6	147	Total	C	H	N	O	S	0	0
			2392	759	1192	225	208	8		

- Molecule 53 is a protein called NDUF57.

Mol	Chain	Residues	Atoms						AltConf	Trace
53	S7	201	Total	C	H	N	O	S	0	0
			3045	975	1500	272	284	14		

- Molecule 54 is a protein called NDUF58.

Mol	Chain	Residues	Atoms						AltConf	Trace
54	S8	182	Total	C	H	N	O	S	0	0
			2843	915	1392	245	275	16		

- Molecule 55 is a protein called UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK.

Mol	Chain	Residues	Atoms					AltConf	Trace
55	U1	12	Total	C	H	N	O	0	0
			76	36	16	12	12		
55	U2	12	Total	C	H	N	O	0	0
			76	36	16	12	12		

- Molecule 56 is a protein called NDUFV1.

Mol	Chain	Residues	Atoms						AltConf	Trace
56	V1	504	Total	C	H	N	O	S	0	0
			7724	2463	3827	680	727	27		

- Molecule 57 is a protein called NDUFV2.

Mol	Chain	Residues	Atoms						AltConf	Trace
57	V2	225	Total	C	H	N	O	S	0	0
			3460	1124	1701	299	319	17		

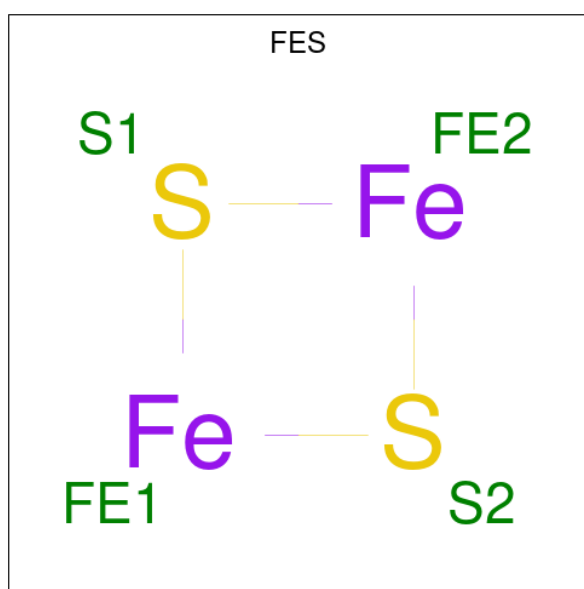
- Molecule 58 is a protein called NDUEG7.

Mol	Chain	Residues	Atoms						AltConf	Trace
58	E7	246	Total	C	H	N	O	S	0	0
			3780	1205	1892	332	344	7		

- Molecule 59 is a protein called NDUFAB1-beta.

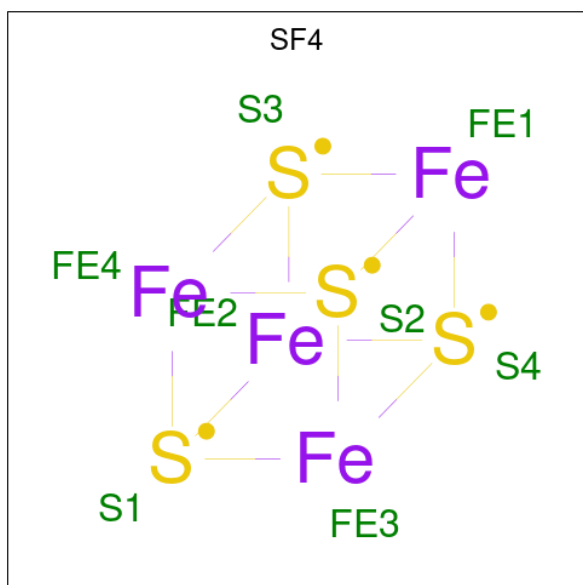
Mol	Chain	Residues	Atoms						AltConf	Trace
59	AC	92	Total	C	H	N	O	S	0	0
			1418	461	697	116	140	4		

- Molecule 60 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe_2S_2).



Mol	Chain	Residues	Atoms			AltConf
60	1A	1	Total	Fe	S	0
			4	2	2	
60	V2	1	Total	Fe	S	0
			4	2	2	

- Molecule 61 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4) (labeled as "Ligand of Interest" by depositor).

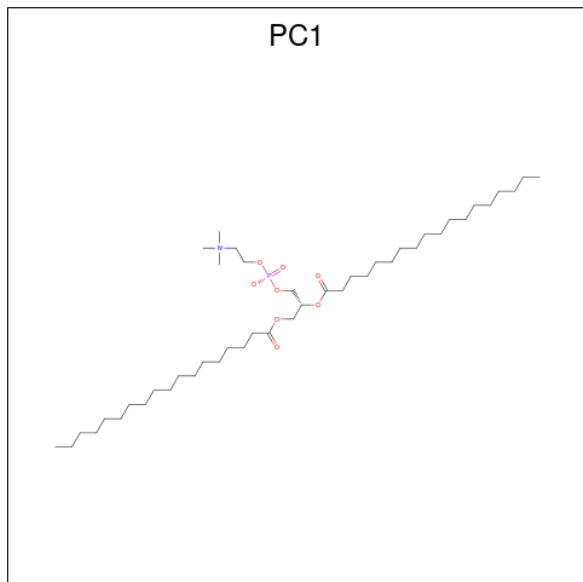


Mol	Chain	Residues	Atoms			AltConf
61	1A	1	Total	Fe	S	0
			8	4	4	
61	1A	1	Total	Fe	S	0
			8	4	4	
61	S7	1	Total	Fe	S	0
			8	4	4	
61	S8	1	Total	Fe	S	0
			8	4	4	
61	S8	1	Total	Fe	S	0
			8	4	4	
61	V1	1	Total	Fe	S	0
			8	4	4	

- Molecule 62 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		AltConf
62	1A	1	Total	K	0
			1	1	

- Molecule 63 is 1,2-DIACYL-SN-GLYCERO-3-PHOSPHOCHOLINE (CCD ID: PC1) (formula: $C_{44}H_{88}NO_8P$).



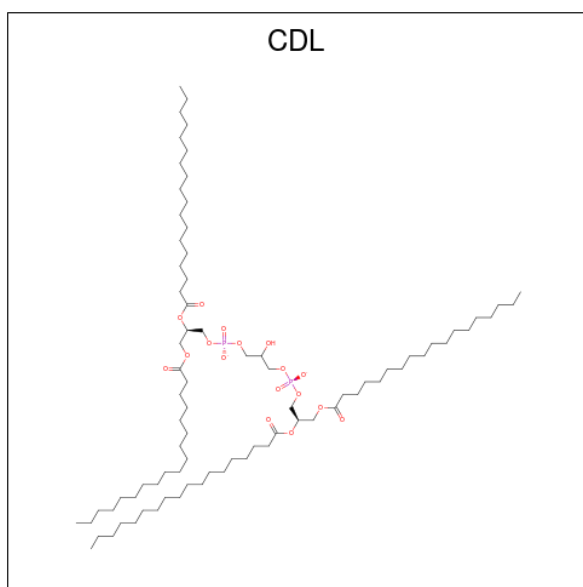
Mol	Chain	Residues	Atoms						AltConf
63	A1	1	Total	C	H	N	O	P	0
			124	39	75	1	8	1	
63	A1	1	Total	C	H	N	O	P	0
			67	21	36	1	8	1	
63	A9	1	Total	C	H	N	O	P	0
			73	23	40	1	8	1	
63	A9	1	Total	C	H	N	O	P	0
			73	23	40	1	8	1	
63	AL	1	Total	C	H	N	O	P	0
			127	40	77	1	8	1	
63	AM	1	Total	C	H	N	O	P	0
			124	39	75	1	8	1	
63	AM	1	Total	C	H	N	O	P	0
			121	38	73	1	8	1	
63	AN	1	Total	C	H	N	O	P	0
			121	38	73	1	8	1	
63	B5	1	Total	C	H	N	O	P	0
			142	44	88	1	8	1	
63	B5	1	Total	C	H	N	O	P	0
			142	44	88	1	8	1	
63	C4	1	Total	C	H	N	O	P	0
			88	28	50	1	8	1	
63	E4	1	Total	C	H	N	O	P	0
			130	41	79	1	8	1	

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Mol	Chain	Residues	Atoms						AltConf
63	E8	1	Total	C	H	N	O	P	0
			142	44	88	1	8	1	
63	E8	1	Total	C	H	N	O	P	0
			142	44	88	1	8	1	
63	E8	1	Total	C	H	N	O	P	0
			73	23	40	1	8	1	
63	E8	1	Total	C	H	N	O	P	0
			64	20	34	1	8	1	
63	ED	1	Total	C	H	N	O	P	0
			142	44	88	1	8	1	
63	N1	1	Total	C	H	N	O	P	0
			124	39	75	1	8	1	
63	N1	1	Total	C	H	N	O	P	0
			94	30	54	1	8	1	
63	N2	1	Total	C	H	N	O	P	0
			85	27	48	1	8	1	
63	N3	1	Total	C	H	N	O	P	0
			103	32	61	1	8	1	
63	N4	1	Total	C	H	N	O	P	0
			91	29	52	1	8	1	
63	N4	1	Total	C	H	N	O	P	0
			73	23	40	1	8	1	
63	N5	1	Total	C	H	N	O	P	0
			142	44	88	1	8	1	
63	N5	1	Total	C	H	N	O	P	0
			97	31	56	1	8	1	
63	N5	1	Total	C	H	N	O	P	0
			82	26	46	1	8	1	

- Molecule 64 is CARDIOLIPIN (CCD ID: CDL) (formula: $C_{81}H_{156}O_{17}P_2$).



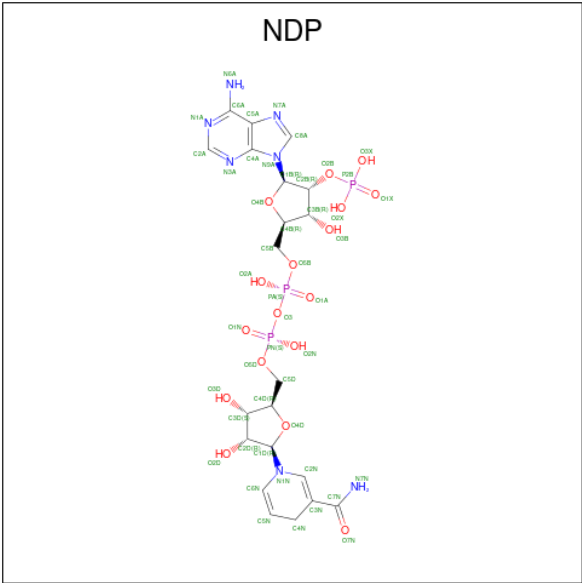
Mol	Chain	Residues	Atoms					AltConf
64	A3	1	Total	C	H	O	P	0
			118	39	60	17	2	
64	AL	1	Total	C	H	O	P	0
			148	49	80	17	2	
64	AL	1	Total	C	H	O	P	0
			136	45	72	17	2	
64	AL	1	Total	C	H	O	P	0
			154	51	84	17	2	
64	AM	1	Total	C	H	O	P	0
			163	53	91	17	2	
64	AM	1	Total	C	H	O	P	0
			163	53	91	17	2	
64	AM	1	Total	C	H	O	P	0
			163	53	91	17	2	
64	B3	1	Total	C	H	O	P	0
			139	46	74	17	2	
64	B5	1	Total	C	H	O	P	0
			118	39	60	17	2	
64	C4	1	Total	C	H	O	P	0
			235	75	141	17	2	
64	C4	1	Total	C	H	O	P	0
			151	50	82	17	2	
64	E6	1	Total	C	H	O	P	0
			136	45	72	17	2	
64	EA	1	Total	C	H	O	P	0
			121	40	62	17	2	
64	EA	1	Total	C	H	O	P	0
			109	36	54	17	2	

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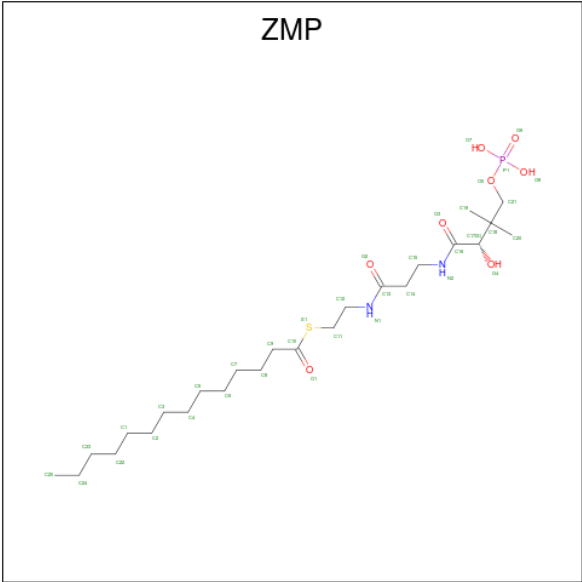
Mol	Chain	Residues	Atoms					AltConf
64	N4	1	Total	C	H	O	P	0
			247	79	149	17	2	
64	N5	1	Total	C	H	O	P	0
			157	51	87	17	2	
64	N5	1	Total	C	H	O	P	0
			229	74	136	17	2	
64	E7	1	Total	C	H	O	P	0
			148	49	80	17	2	

- Molecule 65 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NDP) (formula: C₂₁H₃₀N₇O₁₇P₃).



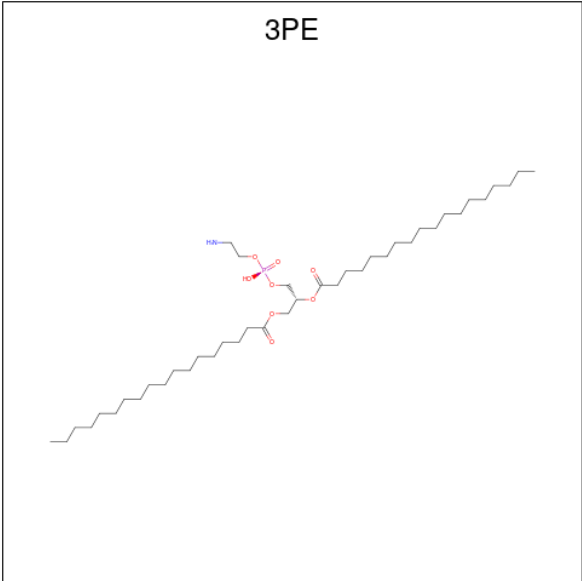
Mol	Chain	Residues	Atoms						AltConf
65	A9	1	Total	C	H	N	O	P	0
			74	21	26	7	17	3	

- Molecule 66 is S-[2-({N-[(2S)-2-hydroxy-3,3-dimethyl-4-(phosphonooxy)butanoyl]-beta-alanyl}amino)ethyl] tetradecanethioate (CCD ID: ZMP) (formula: C₂₅H₄₉N₂O₈PS) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						AltConf
66	AB	1	Total	C	N	O	P	S	0
			36	25	2	7	1	1	
66	AC	1	Total	C	N	O	P	S	0
			36	25	2	7	1	1	

- Molecule 67 is 1,2-Distearoyl-sn-glycerophosphoethanolamine (CCD ID: 3PE) (formula: C₄₁H₈₂NO₈P).



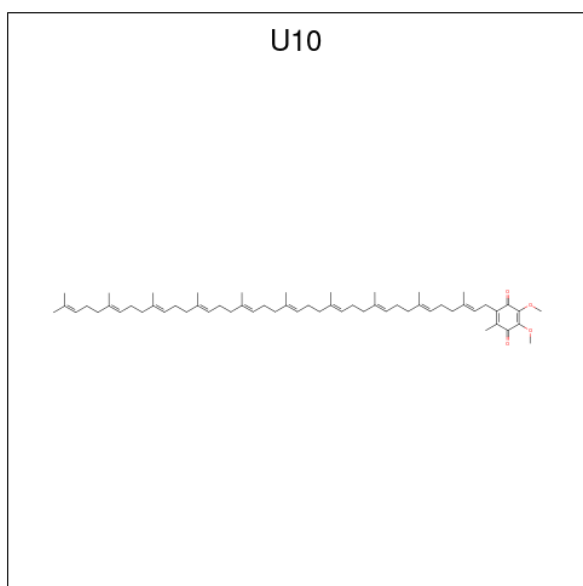
Mol	Chain	Residues	Atoms						AltConf
67	AN	1	Total	C	H	N	O	P	0
			132	41	81	1	8	1	

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Mol	Chain	Residues	Atoms					AltConf
67	G1	1	Total	C	H	N	O	P
			96	30	56	1	8	1
67	N4	1	Total	C	H	N	O	P
			96	31	55	1	8	1
67	N5	1	Total	C	H	N	O	P
			132	41	81	1	8	1

- Molecule 68 is UBIQUINONE-10 (CCD ID: U10) (formula: $C_{59}H_{90}O_4$) (labeled as "Ligand of Interest" by depositor).

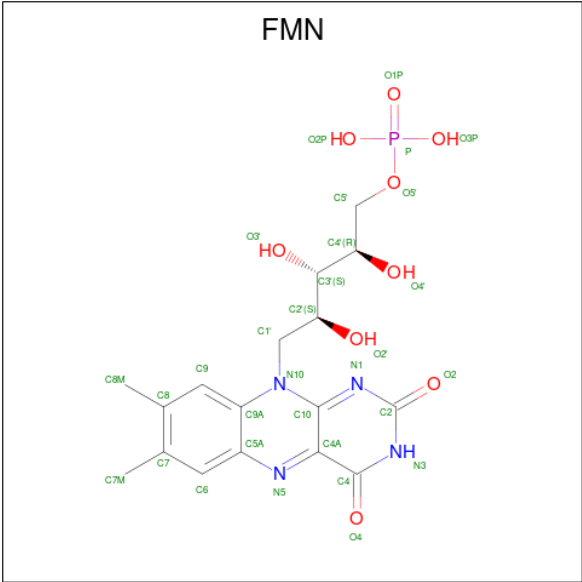


Mol	Chain	Residues	Atoms				AltConf
68	N4	1	Total	C	H	O	
			98	39	55	4	0

- Molecule 69 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
69	S6	1	Total	Zn	
			1	1	0
69	E7	1	Total	Zn	
			1	1	0

- Molecule 70 is FLAVIN MONONUCLEOTIDE (CCD ID: FMN) (formula: $C_{17}H_{21}N_4O_9P$).

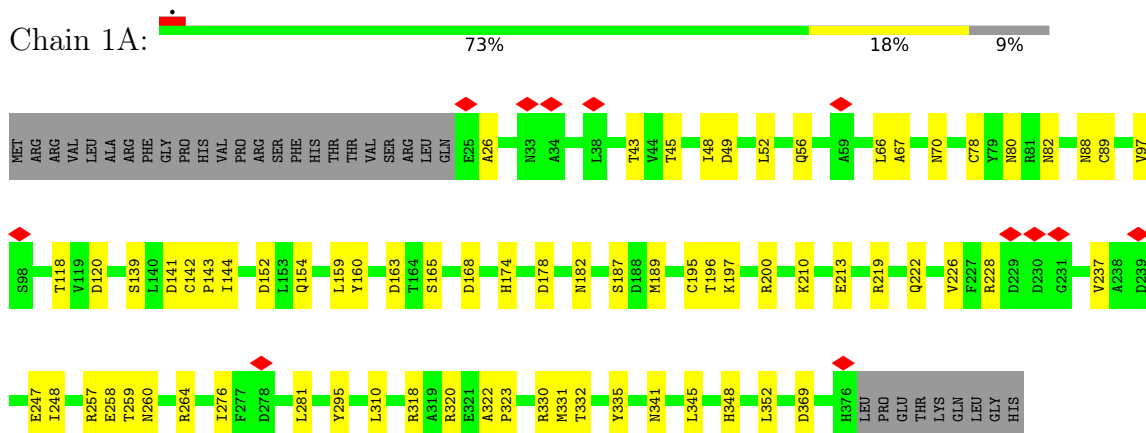


Mol	Chain	Residues	Atoms						AltConf
70	V1	1	Total	C	H	N	O	P	0
			50	17	19	4	9	1	

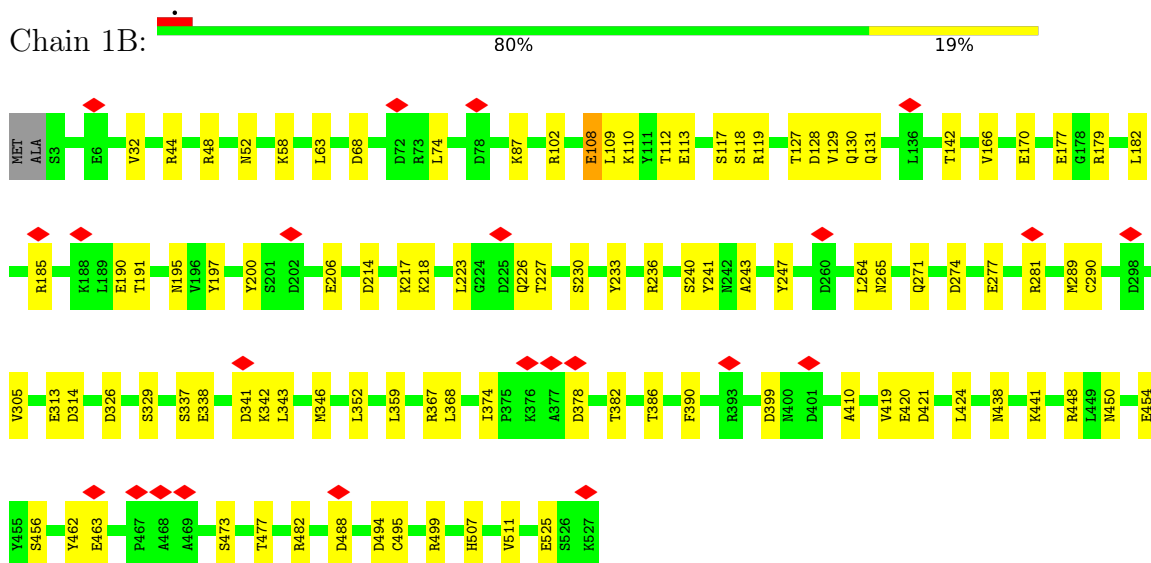
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

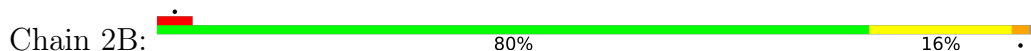
• Molecule 1: NDUS1A

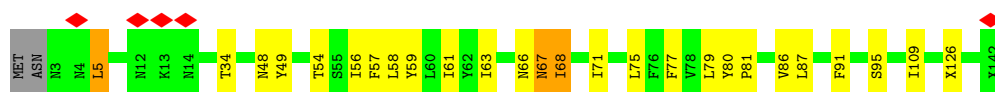


• Molecule 2: NDUS1B

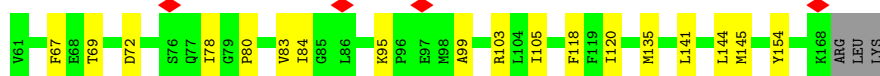
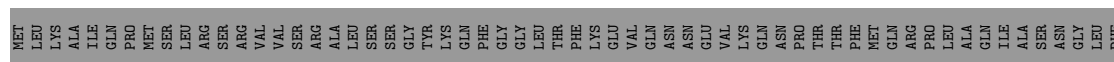


• Molecule 3: NADH dehydrogenase subunit 5

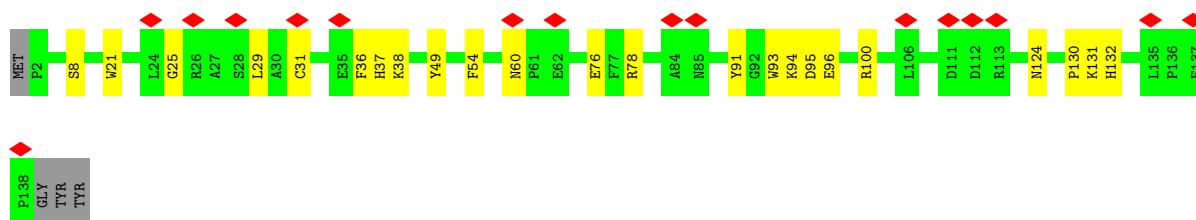
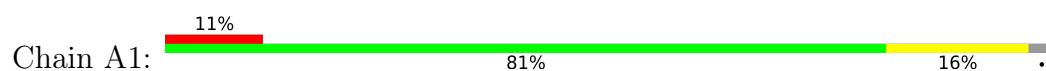




• Molecule 4: ND4L



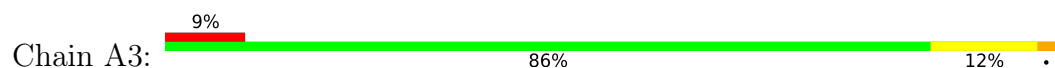
• Molecule 5: NDUFA1



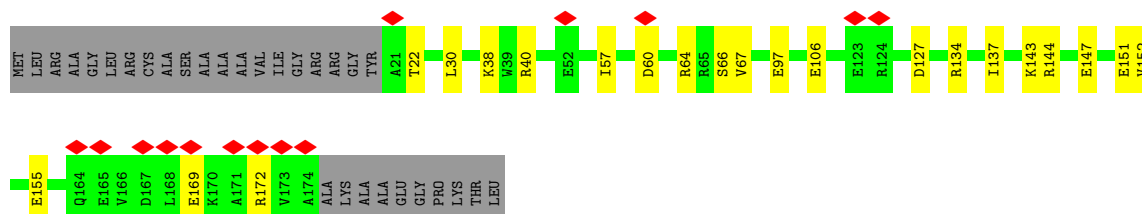
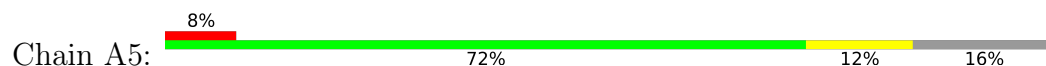
• Molecule 6: NDUFA2



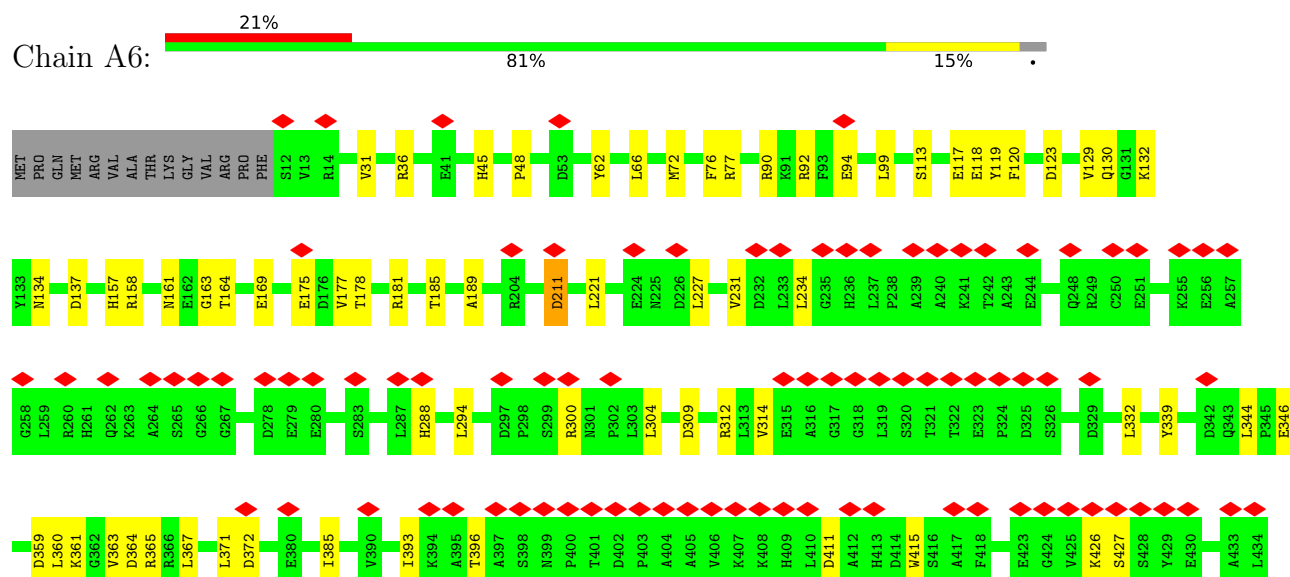
• Molecule 7: NDUFA3



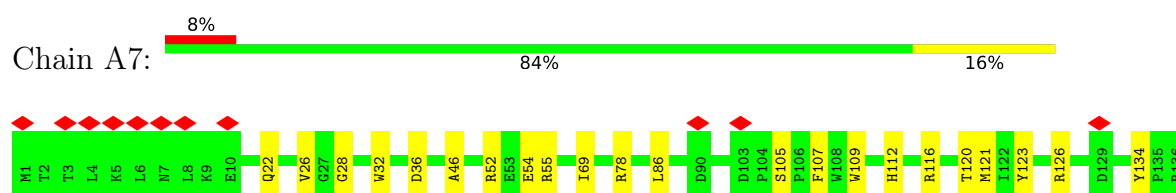
• Molecule 8: NDUFA5



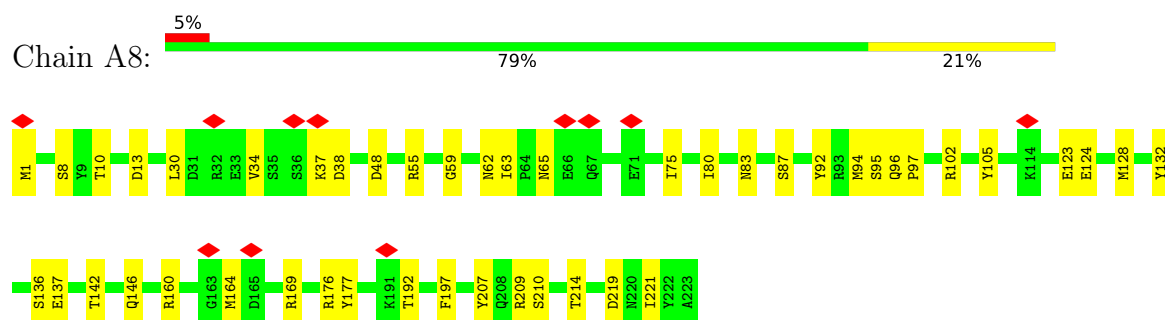
- Molecule 9: NDUFA6



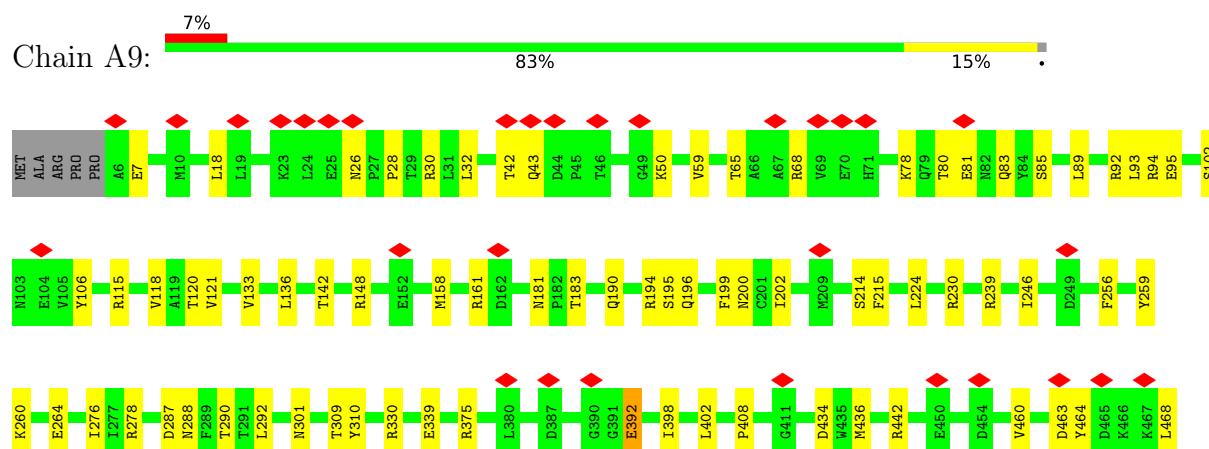
- Molecule 10: NDUFA7

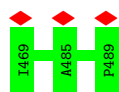


- Molecule 11: NDUFA8

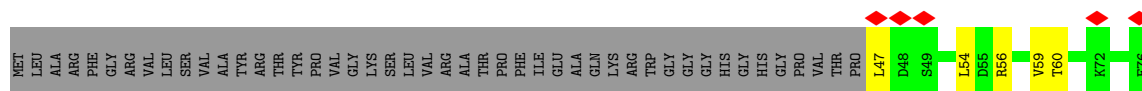


- Molecule 12: NDUFA9

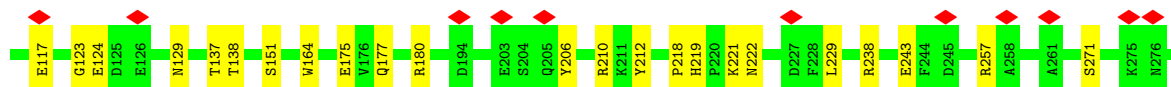
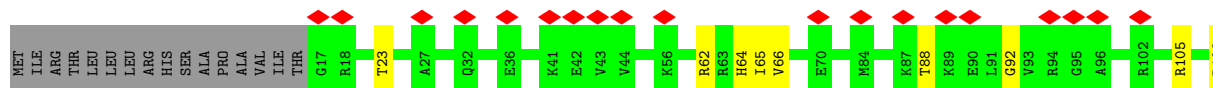
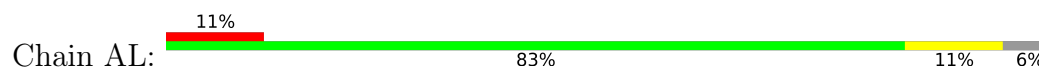




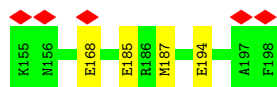
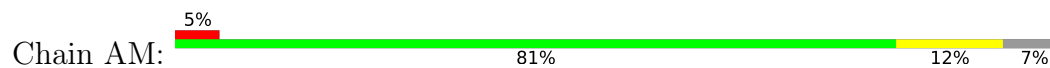
• Molecule 13: NDUFAB1-alpha



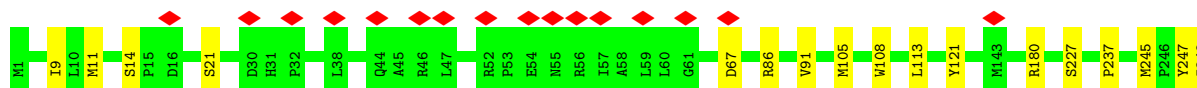
• Molecule 14: NDUFA12

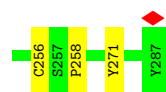


• Molecule 15: NDUFA13

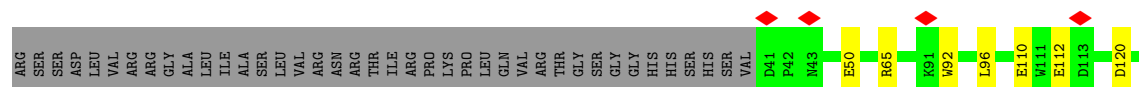


• Molecule 16: NDUFA11

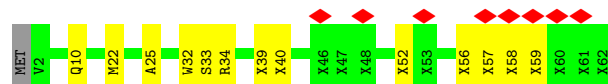
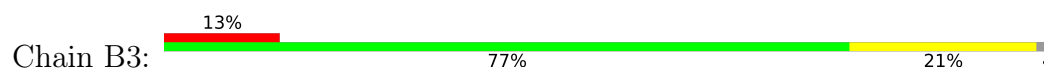




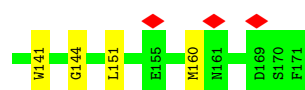
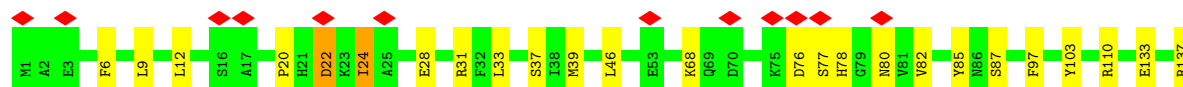
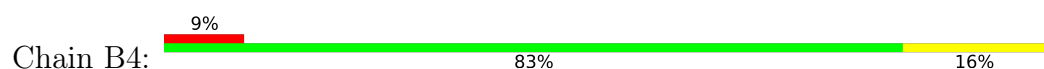
• Molecule 17: NDUFB2



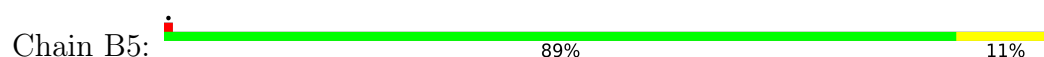
• Molecule 18: NDUFB3



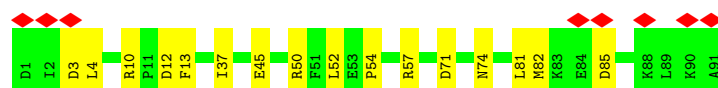
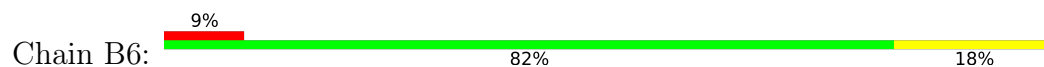
• Molecule 19: NDUFB4



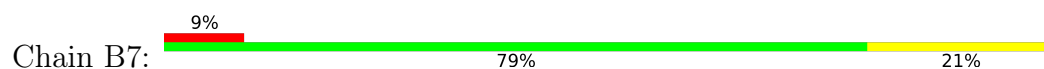
• Molecule 20: NDUFB5

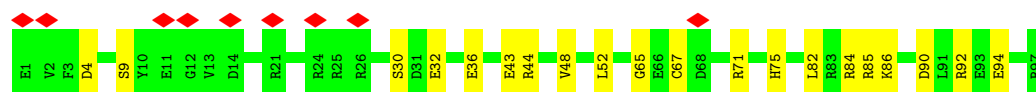


• Molecule 21: NDUFB6

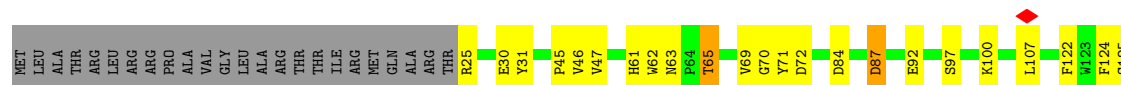


• Molecule 22: NDUFB7

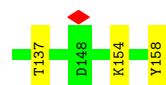
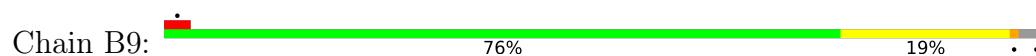




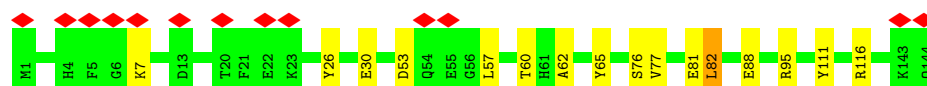
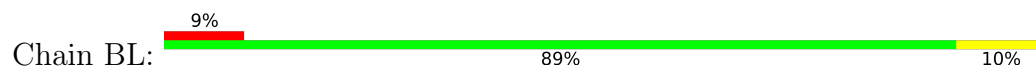
• Molecule 23: NDUFB8



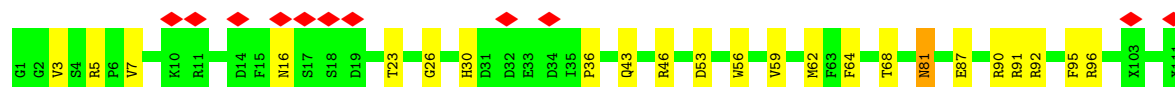
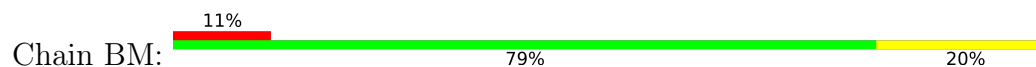
• Molecule 24: NDUFB9



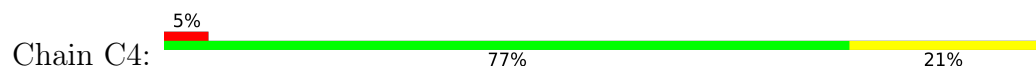
• Molecule 25: NDUFB10

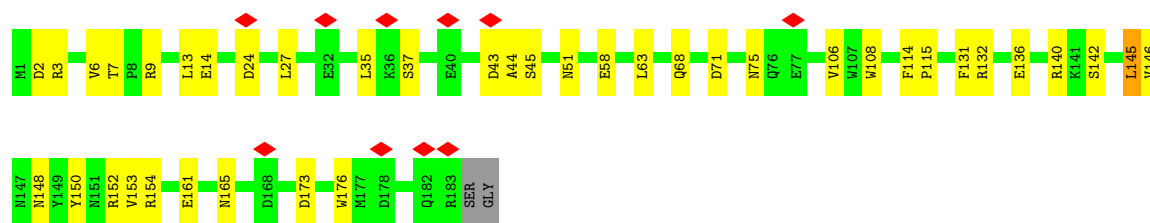


• Molecule 26: NDUFB11

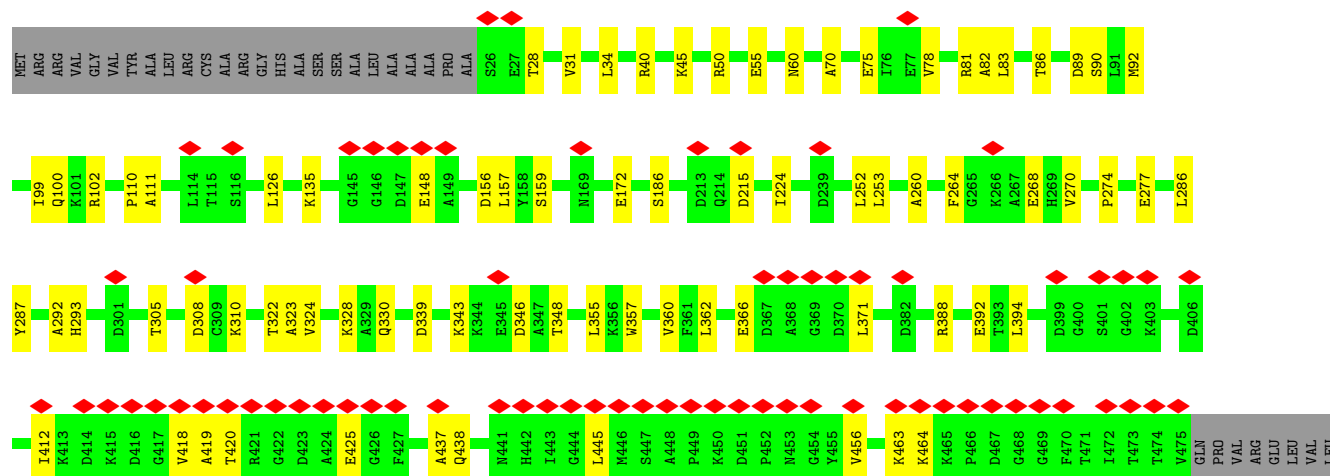
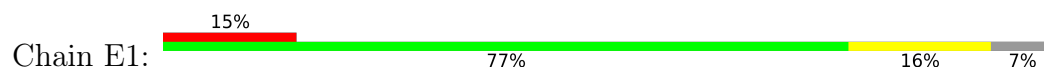


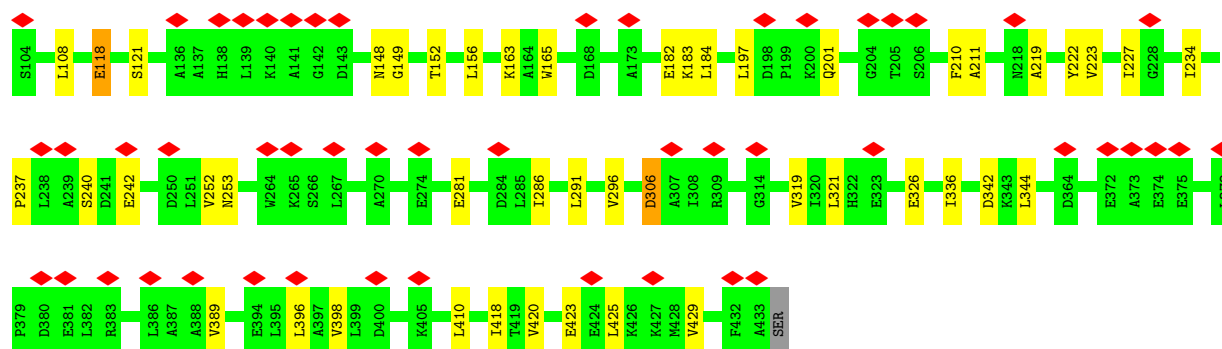
• Molecule 27: NDUFC2



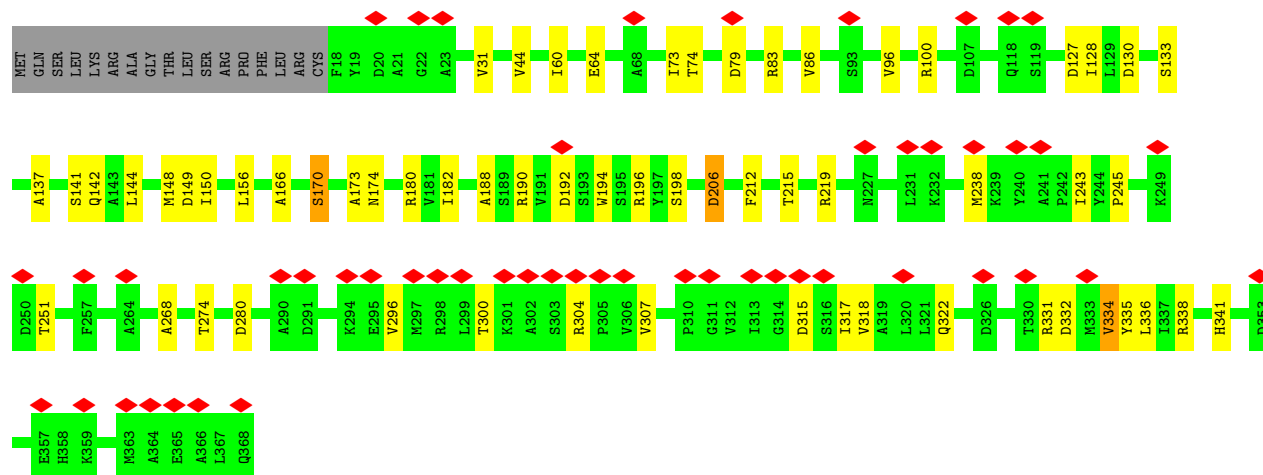
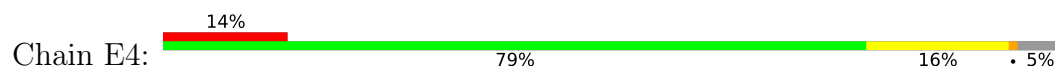


• Molecule 28: NDUEG1

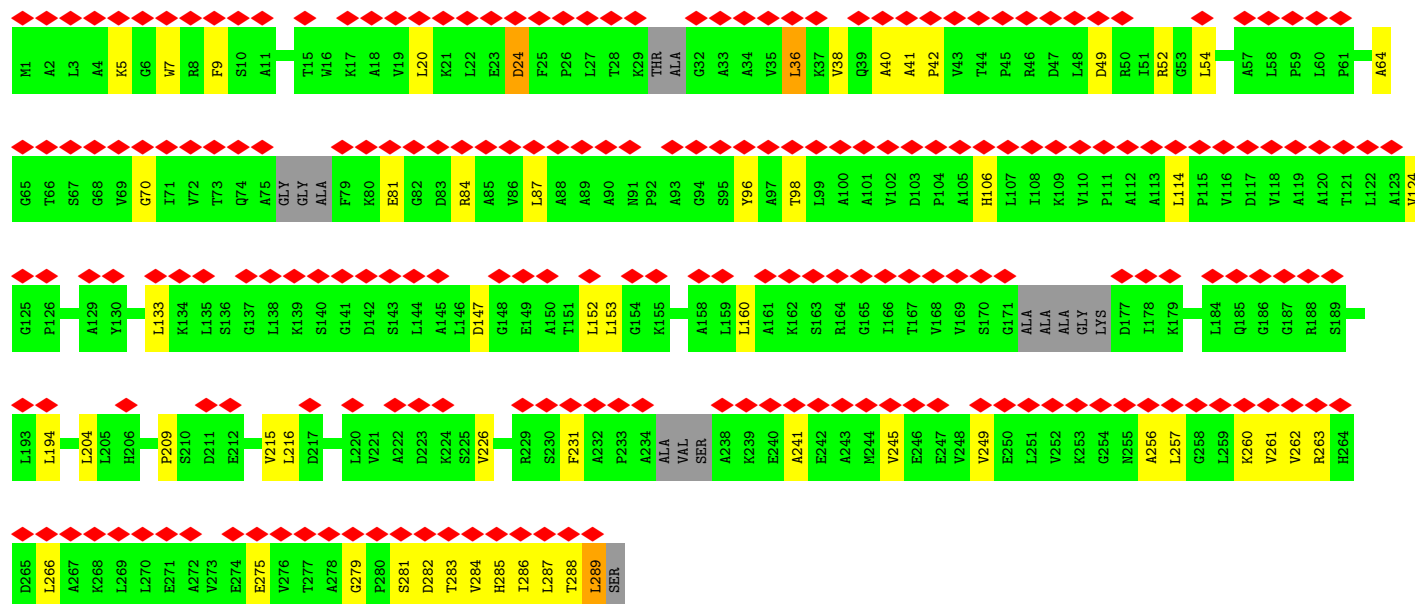
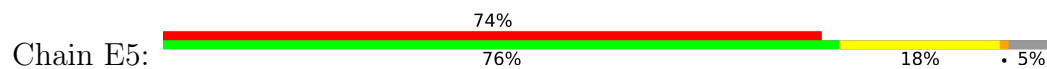




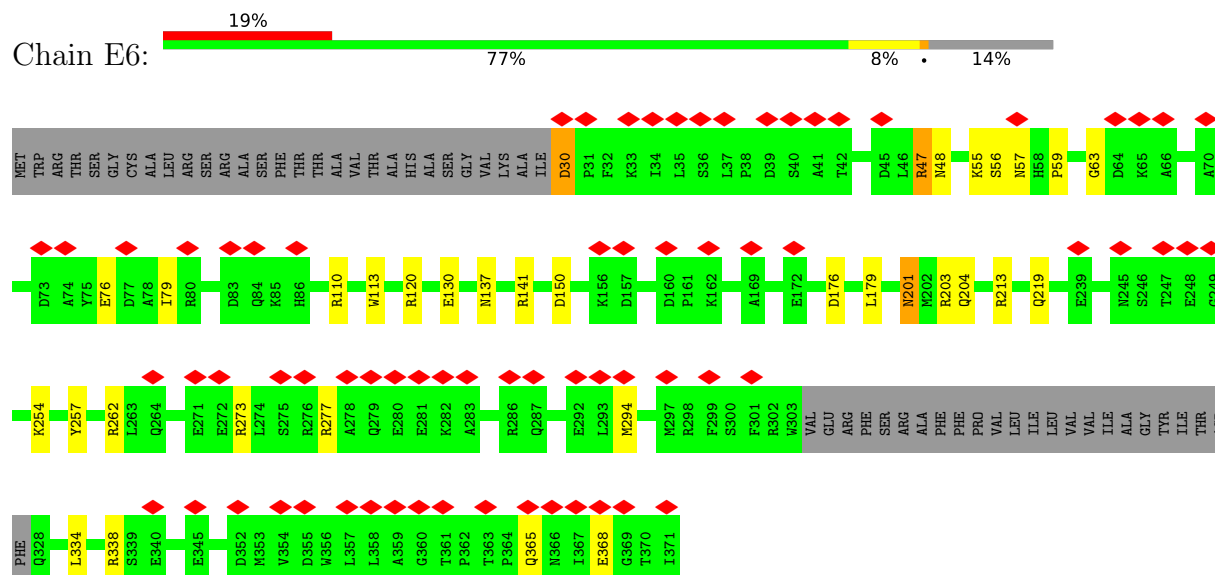
• Molecule 31: NDUEG4



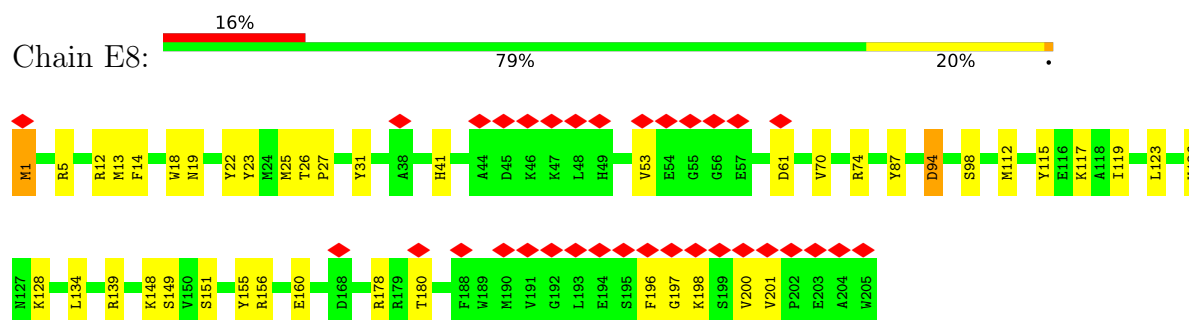
• Molecule 32: NDUEG5



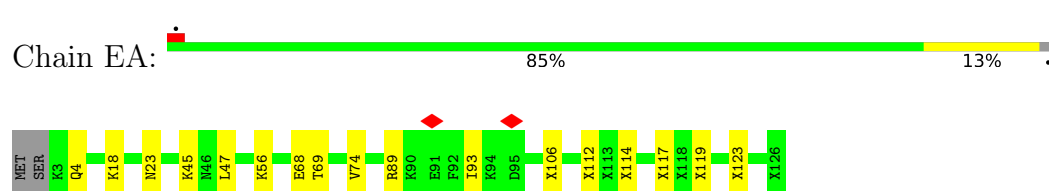
- Molecule 33: NDUEG6



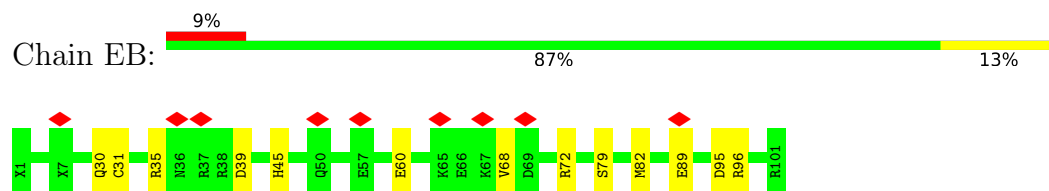
- Molecule 34: NDUEG8



- Molecule 35: NDUEG10

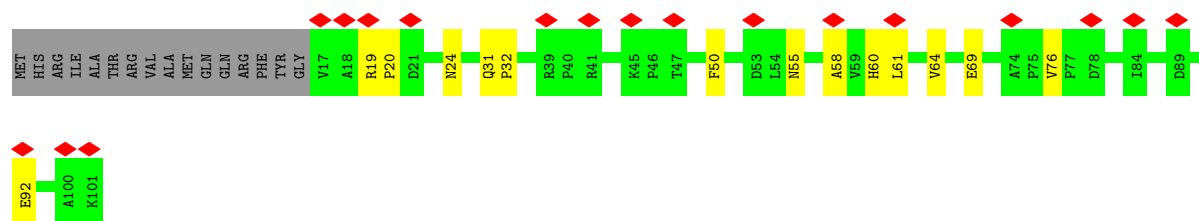


- Molecule 36: NDUEG11

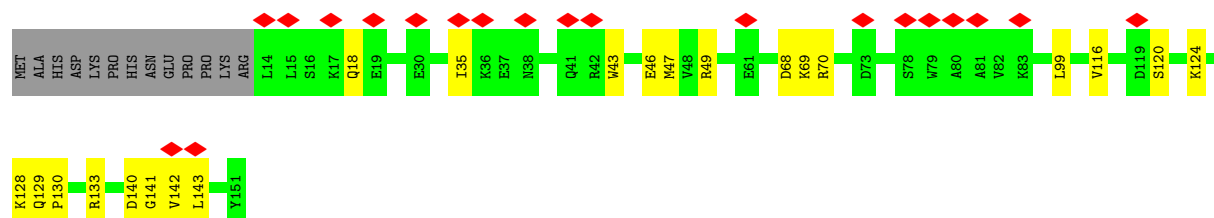


- Molecule 37: NDUEG12

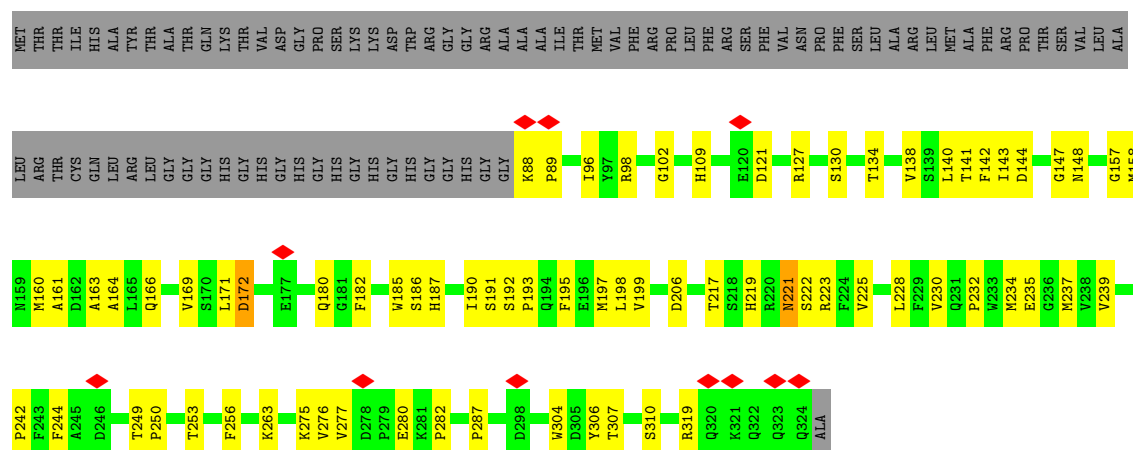




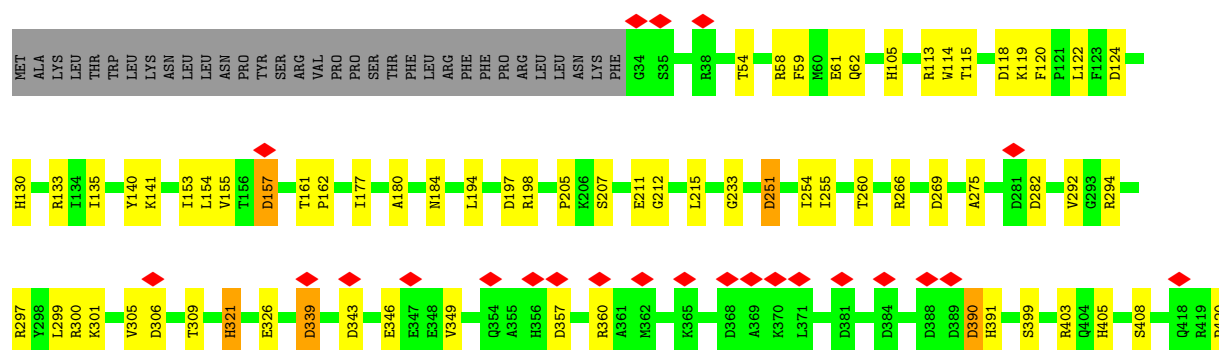
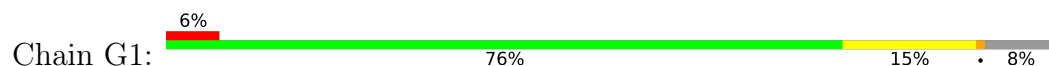
• Molecule 38: NDUEG13



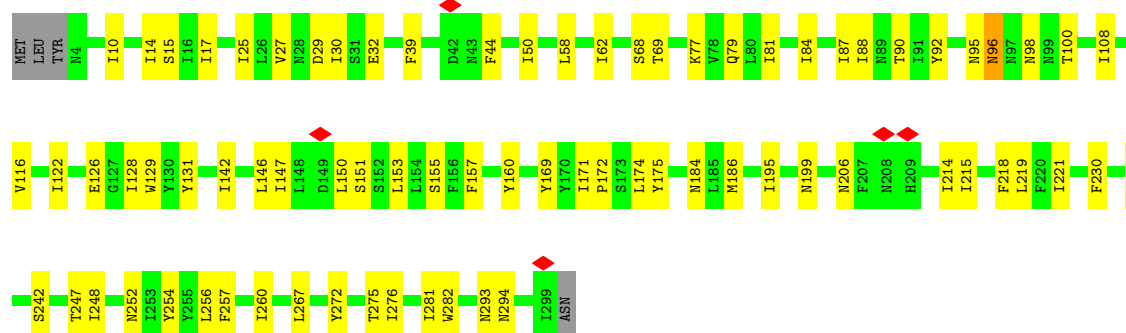
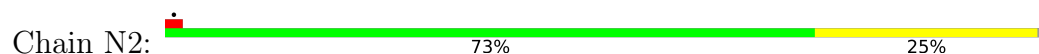
• Molecule 39: NDUFX



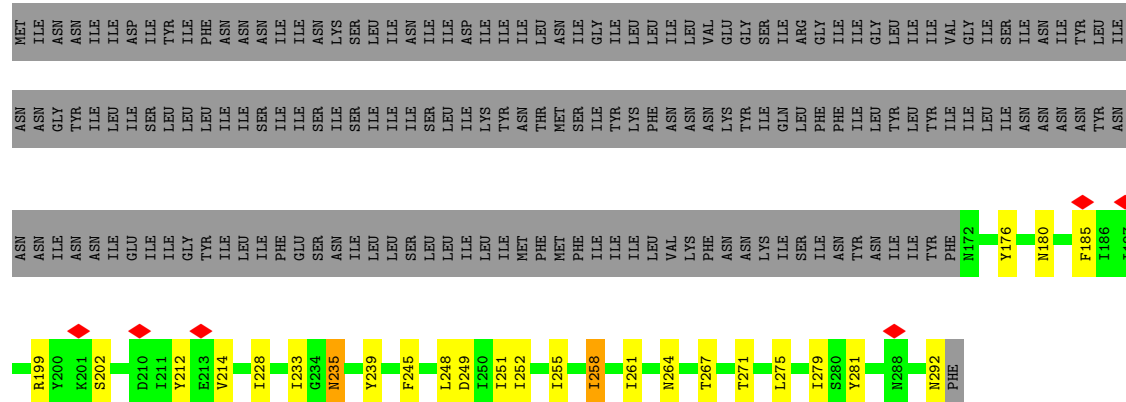
• Molecule 40: NDUCA1



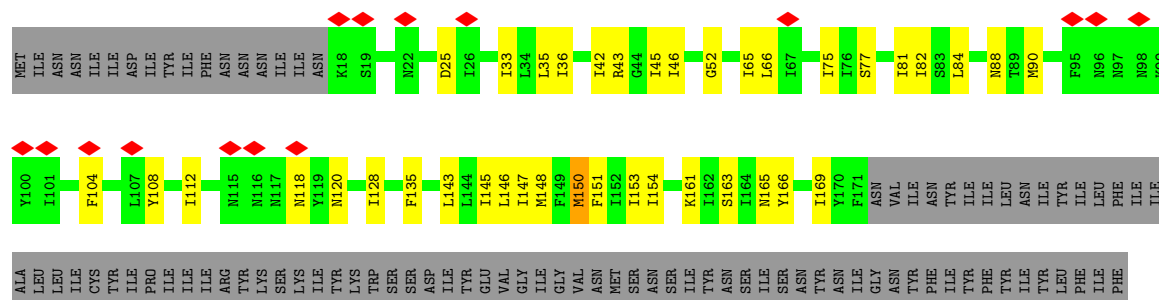
- Molecule 44: ND2A



- Molecule 45: ND3



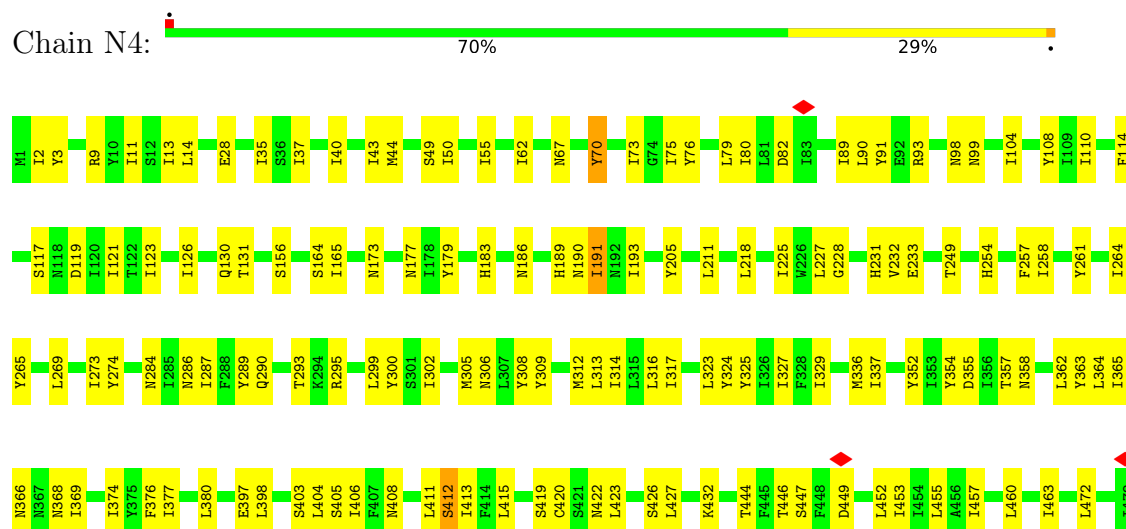
- Molecule 45: ND3



LEU ASP
ILE ILE
ILE ASN
ILE ILE
ILE ILE
ILE LEU
TYR PHE
PHE ILE
ILE ASN
ILE LEU
PHE PHE
THR THR
TYR PHE
PHE THR
LEU LEU
PHE PHE
ILE ILE
ALA ILE
ILE SER
TYR SER
LEU LEU
PHE PHE
SER ASN
ASN PHE

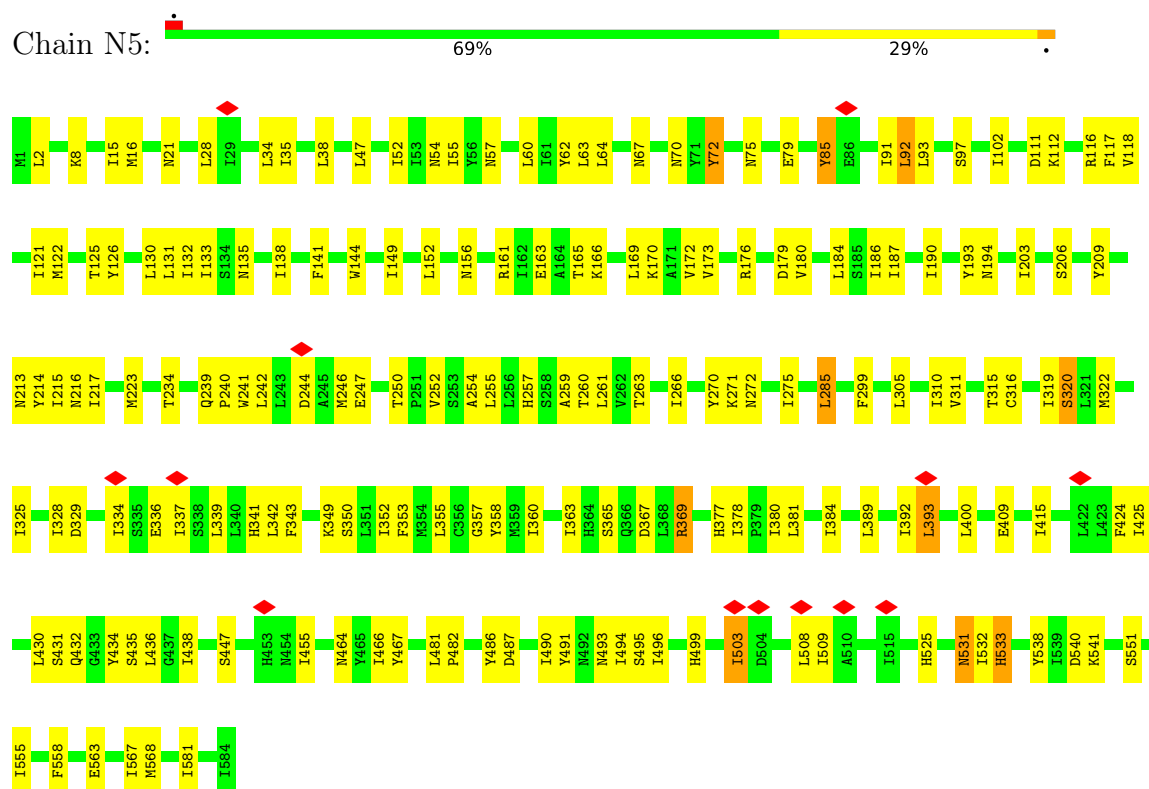
• Molecule 46: NADH-ubiquinone oxidoreductase chain 4

Chain N4:



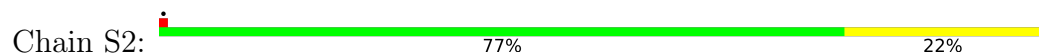
• Molecule 47: ND5

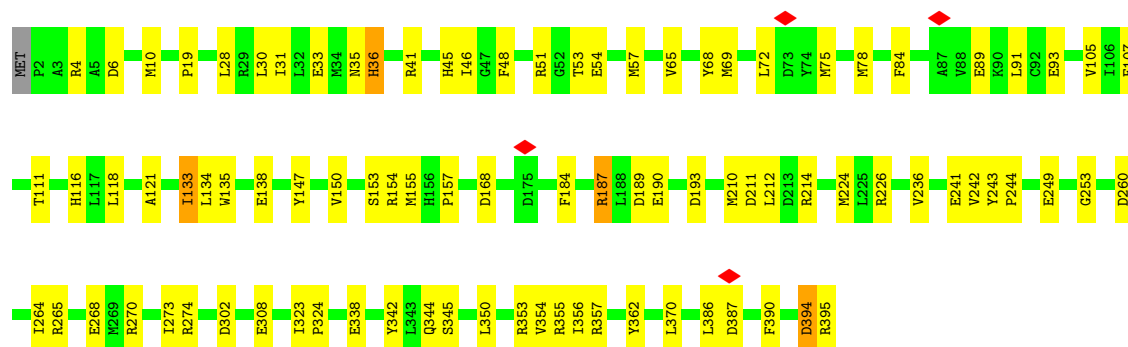
Chain N5:



• Molecule 48: NDUFS2

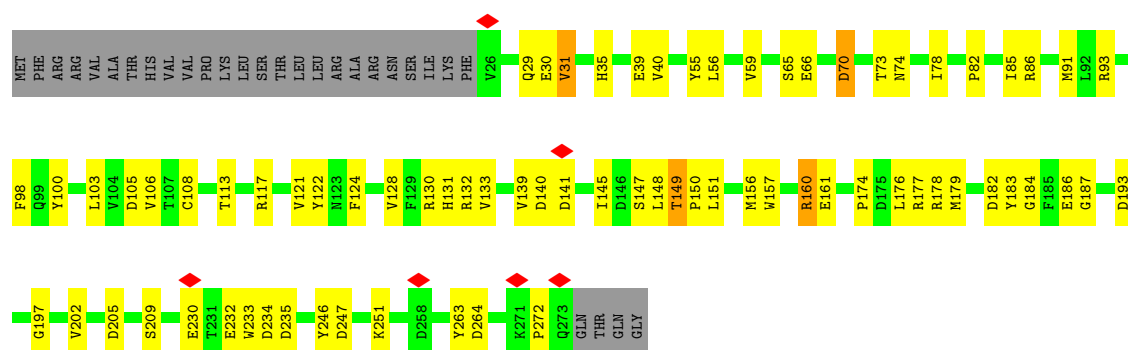
Chain S2:





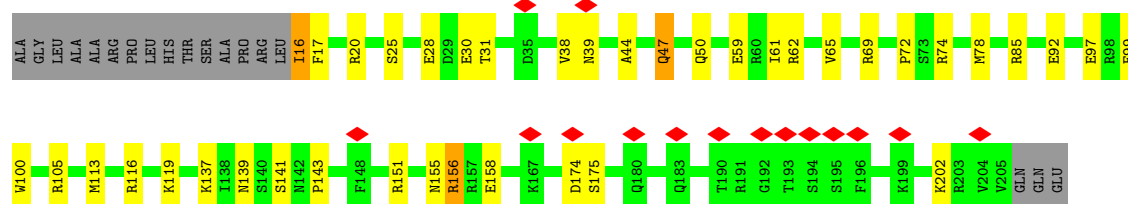
• Molecule 49: NDUFS3

Chain S3: 62% 26% 10%



• Molecule 50: NDUFS4

Chain S4: 72% 18% 9%



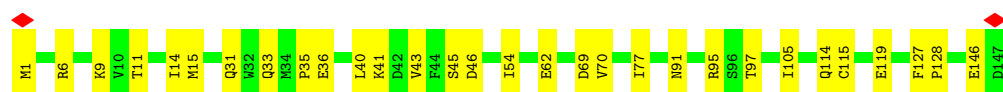
• Molecule 51: NDUFS5

Chain S5: 93% 6% 7%

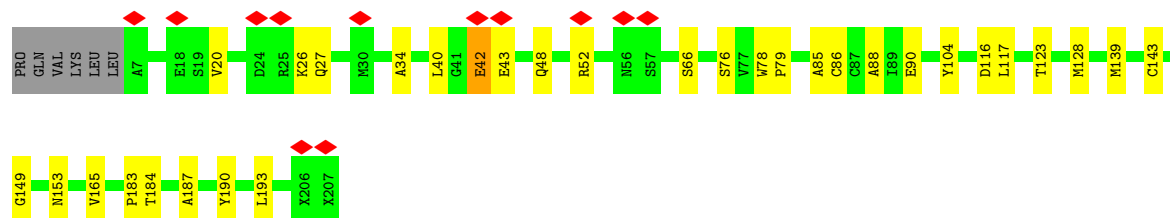
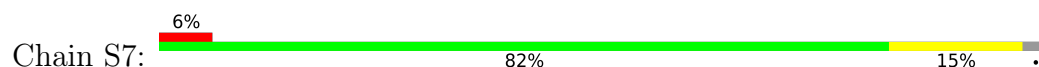


• Molecule 52: NDUFS6

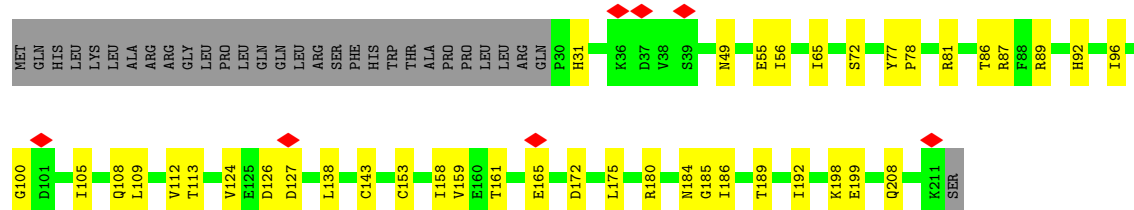
Chain S6: 80% 20%



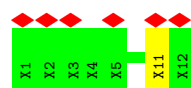
• Molecule 53: NDUF57



• Molecule 54: NDUF58



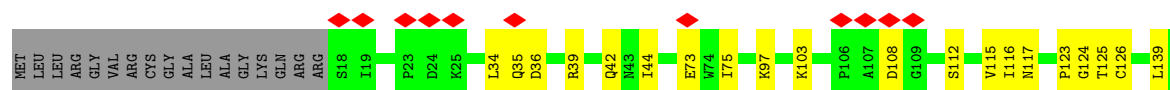
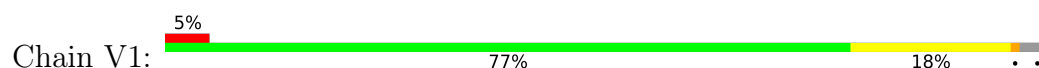
• Molecule 55: UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK

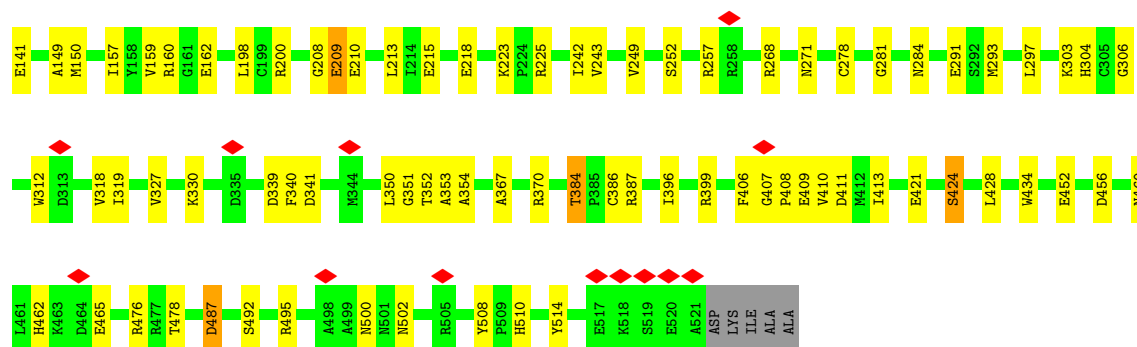


• Molecule 55: UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK

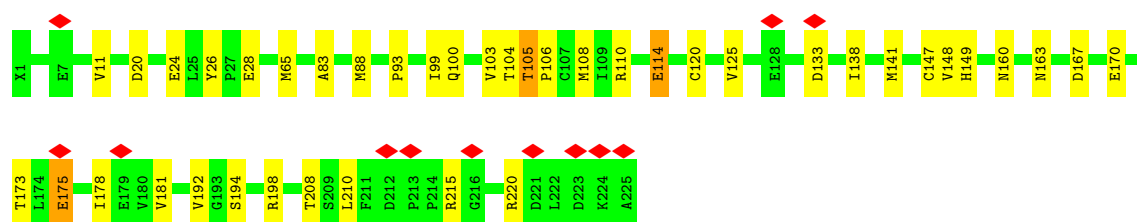
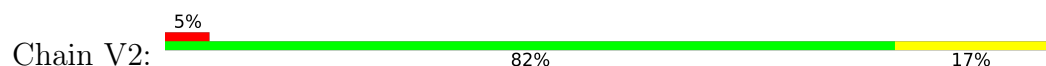


• Molecule 56: NDUFV1

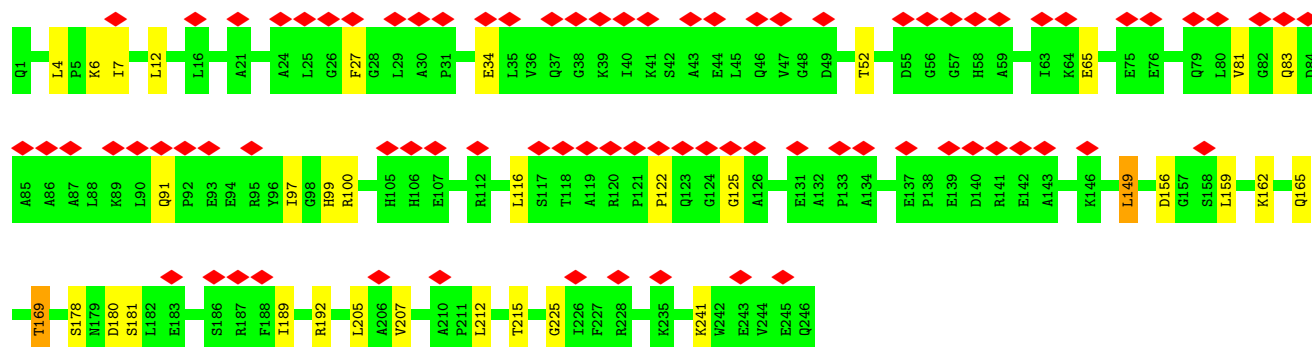
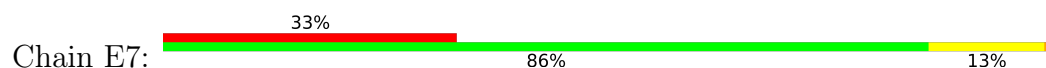




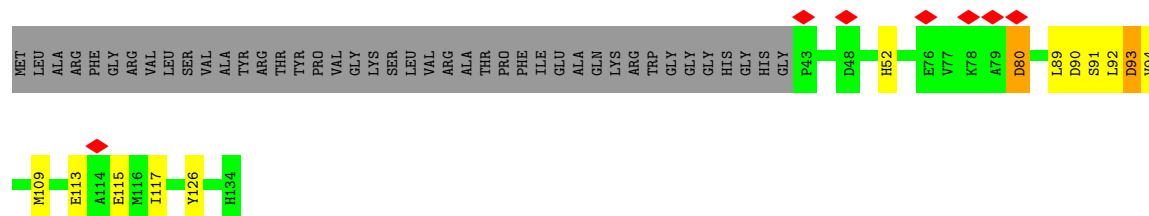
• Molecule 57: NDUFV2



• Molecule 58: NDUEG7



• Molecule 59: NDUFAB1-beta



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	68660	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	61.5	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	130000	Depositor
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	41.957	Depositor
Minimum map value	-16.100	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.994	Depositor
Recommended contour level	5	Depositor
Map size (Å)	446.4, 446.4, 446.4	wwPDB
Map dimensions	480, 480, 480	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.93, 0.93, 0.93	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FES, ZN, ZMP, 2MR, CDL, PC1, K, 3PE, U10, SF4, NDP, FMN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	1A	0.20	0/2858	0.33	0/3878
2	1B	0.17	0/4306	0.28	0/5854
3	2B	0.20	0/958	0.34	0/1306
4	4L	0.19	0/924	0.28	0/1261
5	A1	0.14	0/1108	0.27	0/1511
6	A2	0.15	0/1530	0.28	0/2089
7	A3	0.16	0/1079	0.29	0/1453
8	A5	0.16	0/1282	0.25	0/1737
9	A6	0.15	0/3395	0.27	0/4608
10	A7	0.17	0/1194	0.31	0/1619
11	A8	0.15	0/1879	0.26	0/2543
12	A9	0.17	0/3920	0.29	0/5335
13	AB	0.16	0/704	0.24	0/951
14	AL	0.15	0/2317	0.28	0/3136
15	AM	0.16	0/1533	0.24	0/2079
16	AN	0.15	0/2382	0.25	0/3249
17	B2	0.16	0/947	0.24	0/1291
18	B3	0.17	0/326	0.26	0/441
19	B4	0.18	0/1419	0.29	0/1922
20	B5	0.19	0/1111	0.29	0/1505
21	B6	0.17	0/803	0.27	0/1087
22	B7	0.15	0/877	0.25	0/1172
23	B8	0.19	0/1273	0.27	0/1733
24	B9	0.19	0/1274	0.32	0/1728
25	BL	0.19	0/1266	0.30	0/1710
26	BM	0.19	0/876	0.30	0/1192
27	C4	0.17	0/1592	0.28	0/2158
28	E1	0.14	0/3596	0.27	0/4879
29	E2	0.14	0/3658	0.26	0/4983
30	E3	0.14	0/3320	0.27	0/4520
31	E4	0.15	0/2850	0.26	0/3884
32	E5	0.13	0/2004	0.33	0/2721

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
33	E6	0.14	0/2750	0.25	0/3724
34	E8	0.16	0/1747	0.29	0/2367
35	EA	0.19	0/858	0.29	0/1163
36	EB	0.13	0/650	0.24	0/863
37	EC	0.15	0/676	0.27	0/925
38	ED	0.14	0/1176	0.27	0/1590
39	FX	0.20	0/2035	0.33	0/2763
40	G1	0.18	0/3234	0.31	0/4401
41	G2	0.17	0/1832	0.29	0/2476
42	G3	0.17	0/1957	0.30	0/2646
43	N1	0.17	0/2672	0.33	0/3639
44	N2	0.21	0/2582	0.32	0/3530
45	N3	0.19	0/1068	0.35	0/1456
45	N6	0.17	0/1275	0.32	0/1730
46	N4	0.21	0/4105	0.35	0/5594
47	N5	0.20	0/4963	0.35	0/6758
48	S2	0.21	0/3244	0.33	0/4403
49	S3	0.19	0/2112	0.32	0/2874
50	S4	0.17	0/1573	0.29	0/2107
51	S5	0.13	0/960	0.24	0/1291
52	S6	0.17	0/1232	0.29	0/1659
53	S7	0.18	0/1558	0.28	0/2120
54	S8	0.20	0/1485	0.30	0/2010
56	V1	0.16	0/3990	0.29	0/5394
57	V2	0.16	0/1787	0.29	0/2428
58	E7	0.15	0/1931	0.27	0/2618
59	AC	0.17	0/736	0.25	0/1000
All	All	0.17	0/112749	0.29	0/153064

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	1A	2801	2700	2710	53	0
2	1B	4198	4159	4175	76	0
3	2B	1070	989	1008	24	0
4	4L	890	878	880	19	0
5	A1	1071	1026	1030	15	0
6	A2	1493	1474	1478	15	0
7	A3	1050	1039	1041	15	0
8	A5	1261	1248	1251	15	0
9	A6	3328	3280	3293	57	0
10	A7	1154	1118	1123	23	0
11	A8	1822	1726	1736	40	0
12	A9	3829	3850	3857	57	0
13	AB	694	673	677	14	0
14	AL	2237	2172	2180	27	0
15	AM	1487	1448	1452	26	0
16	AN	2306	2267	2275	16	0
17	B2	913	857	858	13	0
18	B3	449	309	311	11	0
19	B4	1377	1358	1364	27	0
20	B5	1112	1069	1075	15	0
21	B6	773	747	751	19	0
22	B7	857	835	841	13	0
23	B8	1224	1127	1136	25	0
24	B9	1236	1207	1212	34	0
25	BL	1227	1179	1185	13	0
26	BM	910	827	830	25	0
27	C4	1545	1517	1519	30	0
28	E1	3512	3496	3510	48	0
29	E2	3563	3540	3554	44	0
30	E3	3255	3263	3279	45	0
31	E4	2770	2732	2742	51	0
32	E5	1977	2069	2075	45	0
33	E6	2674	2554	2562	25	0
34	E8	1691	1663	1668	37	0
35	EA	961	832	837	11	0
36	EB	774	631	636	11	0
37	EC	660	663	666	10	0
38	ED	1142	1131	1134	16	0
39	FX	1967	1849	1858	58	0
40	G1	3147	2999	3015	58	0
41	G2	1804	1846	1850	31	0
42	G3	1961	1944	1950	35	0
43	N1	2605	2726	2729	49	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
44	N2	2512	2589	2592	67	0
45	N3	1037	1057	1057	22	0
45	N6	1257	1385	1385	38	0
46	N4	4001	4214	4224	110	0
47	N5	4837	5032	5046	152	0
48	S2	3173	3101	3114	76	0
49	S3	2050	1928	1936	61	0
50	S4	1536	1502	1505	33	0
51	S5	991	895	898	7	0
52	S6	1200	1192	1198	23	0
53	S7	1545	1500	1503	27	0
54	S8	1451	1392	1397	38	0
55	U1	60	16	19	1	0
55	U2	60	16	17	1	0
56	V1	3897	3827	3837	76	0
57	V2	1759	1701	1711	28	0
58	E7	1888	1892	1903	24	0
59	AC	721	697	702	15	0
60	1A	4	0	0	0	0
60	V2	4	0	0	0	0
61	1A	16	0	0	0	0
61	S7	8	0	0	0	0
61	S8	16	0	0	5	0
61	V1	8	0	0	1	0
62	1A	1	0	0	0	0
63	A1	80	111	111	3	0
63	A9	66	80	80	1	0
63	AL	50	77	77	0	0
63	AM	97	148	148	1	0
63	AN	48	73	73	0	0
63	B5	108	176	176	2	0
63	C4	38	50	50	1	0
63	E4	51	79	79	0	0
63	E8	171	250	250	3	0
63	ED	54	88	88	1	0
63	N1	89	129	129	0	0
63	N2	37	48	48	1	0
63	N3	42	61	61	0	0
63	N4	72	92	92	2	0
63	N5	131	190	190	1	0
64	A3	58	60	60	1	0
64	AL	202	236	236	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
64	AM	216	273	273	10	0
64	B3	65	74	74	1	0
64	B5	58	60	60	1	0
64	C4	163	223	223	1	0
64	E6	64	72	72	0	0
64	E7	68	80	80	0	0
64	EA	114	116	116	1	0
64	N4	98	149	149	4	0
64	N5	163	223	223	4	0
65	A9	48	26	26	3	0
66	AB	36	0	47	6	0
66	AC	36	0	47	15	0
67	AN	51	81	82	0	0
67	G1	40	56	57	2	0
67	N4	41	55	56	0	0
67	N5	51	81	82	1	0
68	N4	43	55	55	5	0
69	E7	1	0	0	0	0
69	S6	1	0	0	0	0
70	V1	31	19	19	0	0
All	All	113591	112544	113046	1675	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

The worst 5 of 1675 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:E3:222:TYR:HH	37:EC:60:HIS:HE2	1.06	0.99
32:E5:287:LEU:O	32:E5:289:LEU:N	1.97	0.97
10:A7:105:SER:O	49:S3:130:ARG:NH1	2.02	0.93
46:N4:293:THR:HG21	46:N4:365:ILE:HD11	1.50	0.91
17:B2:120:ASP:OD2	17:B2:130:LYS:NZ	2.04	0.90

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	1A	350/385 (91%)	335 (96%)	15 (4%)	0	100	100
2	1B	523/527 (99%)	501 (96%)	22 (4%)	0	100	100
3	2B	112/142 (79%)	103 (92%)	9 (8%)	0	100	100
4	4L	106/171 (62%)	104 (98%)	2 (2%)	0	100	100
5	A1	135/141 (96%)	124 (92%)	11 (8%)	0	100	100
6	A2	190/193 (98%)	186 (98%)	4 (2%)	0	100	100
7	A3	122/125 (98%)	117 (96%)	5 (4%)	0	100	100
8	A5	152/184 (83%)	149 (98%)	3 (2%)	0	100	100
9	A6	421/437 (96%)	405 (96%)	16 (4%)	0	100	100
10	A7	134/136 (98%)	123 (92%)	11 (8%)	0	100	100
11	A8	221/223 (99%)	212 (96%)	9 (4%)	0	100	100
12	A9	482/489 (99%)	463 (96%)	19 (4%)	0	100	100
13	AB	86/134 (64%)	83 (96%)	3 (4%)	0	100	100
14	AL	263/281 (94%)	248 (94%)	15 (6%)	0	100	100
15	AM	182/198 (92%)	175 (96%)	7 (4%)	0	100	100
16	AN	285/287 (99%)	281 (99%)	4 (1%)	0	100	100
17	B2	103/145 (71%)	103 (100%)	0	0	100	100
18	B3	32/62 (52%)	31 (97%)	1 (3%)	0	100	100
19	B4	169/171 (99%)	154 (91%)	15 (9%)	0	100	100
20	B5	132/140 (94%)	130 (98%)	2 (2%)	0	100	100
21	B6	89/91 (98%)	89 (100%)	0	0	100	100
22	B7	95/97 (98%)	94 (99%)	1 (1%)	0	100	100
23	B8	145/176 (82%)	133 (92%)	12 (8%)	0	100	100
24	B9	149/158 (94%)	140 (94%)	9 (6%)	0	100	100
25	BL	142/144 (99%)	139 (98%)	3 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
26	BM	99/112 (88%)	98 (99%)	1 (1%)	0	100	100
27	C4	181/185 (98%)	174 (96%)	7 (4%)	0	100	100
28	E1	448/483 (93%)	434 (97%)	14 (3%)	0	100	100
29	E2	464/467 (99%)	445 (96%)	19 (4%)	0	100	100
30	E3	430/434 (99%)	420 (98%)	10 (2%)	0	100	100
31	E4	349/368 (95%)	338 (97%)	11 (3%)	0	100	100
32	E5	266/290 (92%)	245 (92%)	20 (8%)	1 (0%)	30	60
33	E6	314/371 (85%)	310 (99%)	4 (1%)	0	100	100
34	E8	203/205 (99%)	191 (94%)	12 (6%)	0	100	100
35	EA	96/126 (76%)	92 (96%)	4 (4%)	0	100	100
36	EB	73/101 (72%)	72 (99%)	1 (1%)	0	100	100
37	EC	83/101 (82%)	77 (93%)	6 (7%)	0	100	100
38	ED	136/151 (90%)	130 (96%)	6 (4%)	0	100	100
39	FX	235/325 (72%)	224 (95%)	10 (4%)	1 (0%)	30	60
40	G1	401/436 (92%)	380 (95%)	21 (5%)	0	100	100
41	G2	234/267 (88%)	218 (93%)	16 (7%)	0	100	100
42	G3	253/261 (97%)	235 (93%)	18 (7%)	0	100	100
43	N1	308/670 (46%)	289 (94%)	19 (6%)	0	100	100
44	N2	294/300 (98%)	279 (95%)	15 (5%)	0	100	100
45	N3	119/293 (41%)	115 (97%)	4 (3%)	0	100	100
45	N6	152/293 (52%)	146 (96%)	6 (4%)	0	100	100
46	N4	476/478 (100%)	461 (97%)	15 (3%)	0	100	100
47	N5	582/584 (100%)	552 (95%)	30 (5%)	0	100	100
48	S2	391/395 (99%)	369 (94%)	21 (5%)	1 (0%)	36	65
49	S3	246/277 (89%)	235 (96%)	11 (4%)	0	100	100
50	S4	188/208 (90%)	174 (93%)	14 (7%)	0	100	100
51	S5	110/122 (90%)	108 (98%)	2 (2%)	0	100	100
52	S6	145/147 (99%)	137 (94%)	8 (6%)	0	100	100
53	S7	195/207 (94%)	182 (93%)	13 (7%)	0	100	100
54	S8	180/212 (85%)	176 (98%)	4 (2%)	0	100	100
56	V1	502/526 (95%)	478 (95%)	24 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
57	V2	220/225 (98%)	211 (96%)	9 (4%)	0	100	100
58	E7	244/246 (99%)	238 (98%)	6 (2%)	0	100	100
59	AC	90/134 (67%)	88 (98%)	2 (2%)	0	100	100
All	All	13527/15237 (89%)	12943 (96%)	581 (4%)	3 (0%)	100	100

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
32	E5	288	THR
48	S2	36	HIS
39	FX	276	VAL

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	1A	310/340 (91%)	306 (99%)	4 (1%)	61	75
2	1B	453/454 (100%)	450 (99%)	3 (1%)	76	80
3	2B	109/111 (98%)	105 (96%)	4 (4%)	30	59
4	4L	96/151 (64%)	95 (99%)	1 (1%)	68	77
5	A1	115/118 (98%)	112 (97%)	3 (3%)	40	66
6	A2	159/160 (99%)	158 (99%)	1 (1%)	78	81
7	A3	104/104 (100%)	101 (97%)	3 (3%)	37	63
8	A5	134/152 (88%)	131 (98%)	3 (2%)	45	68
9	A6	346/358 (97%)	344 (99%)	2 (1%)	78	81
10	A7	119/119 (100%)	118 (99%)	1 (1%)	73	80
11	A8	196/196 (100%)	192 (98%)	4 (2%)	48	70
12	A9	420/424 (99%)	417 (99%)	3 (1%)	76	80
13	AB	79/114 (69%)	77 (98%)	2 (2%)	42	66
14	AL	228/242 (94%)	227 (100%)	1 (0%)	84	85

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
15	AM	156/168 (93%)	155 (99%)	1 (1%)	78	81
16	AN	241/241 (100%)	239 (99%)	2 (1%)	73	80
17	B2	97/131 (74%)	97 (100%)	0	100	100
18	B3	30/31 (97%)	30 (100%)	0	100	100
19	B4	144/144 (100%)	141 (98%)	3 (2%)	47	69
20	B5	108/108 (100%)	108 (100%)	0	100	100
21	B6	82/82 (100%)	82 (100%)	0	100	100
22	B7	93/93 (100%)	91 (98%)	2 (2%)	45	68
23	B8	127/148 (86%)	123 (97%)	4 (3%)	35	62
24	B9	132/139 (95%)	131 (99%)	1 (1%)	73	80
25	BL	132/132 (100%)	129 (98%)	3 (2%)	44	68
26	BM	93/93 (100%)	91 (98%)	2 (2%)	45	68
27	C4	166/167 (99%)	164 (99%)	2 (1%)	63	76
28	E1	381/404 (94%)	379 (100%)	2 (0%)	81	83
29	E2	379/380 (100%)	372 (98%)	7 (2%)	51	71
30	E3	339/341 (99%)	333 (98%)	6 (2%)	51	71
31	E4	302/317 (95%)	298 (99%)	4 (1%)	61	75
32	E5	200/205 (98%)	195 (98%)	5 (2%)	42	66
33	E6	272/314 (87%)	266 (98%)	6 (2%)	45	68
34	E8	179/179 (100%)	176 (98%)	3 (2%)	53	72
35	EA	84/86 (98%)	83 (99%)	1 (1%)	63	76
36	EB	70/70 (100%)	70 (100%)	0	100	100
37	EC	73/86 (85%)	73 (100%)	0	100	100
38	ED	121/133 (91%)	119 (98%)	2 (2%)	53	72
39	FX	212/276 (77%)	208 (98%)	4 (2%)	50	71
40	G1	333/365 (91%)	324 (97%)	9 (3%)	39	65
41	G2	192/214 (90%)	188 (98%)	4 (2%)	47	69
42	G3	202/202 (100%)	202 (100%)	0	100	100
43	N1	295/639 (46%)	291 (99%)	4 (1%)	59	74
44	N2	285/289 (99%)	278 (98%)	7 (2%)	42	66
45	N3	116/281 (41%)	111 (96%)	5 (4%)	26	55

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
45	N6	147/281 (52%)	143 (97%)	4 (3%)	39	65
46	N4	455/455 (100%)	447 (98%)	8 (2%)	51	71
47	N5	546/546 (100%)	533 (98%)	13 (2%)	43	67
48	S2	335/336 (100%)	330 (98%)	5 (2%)	57	73
49	S3	224/250 (90%)	216 (96%)	8 (4%)	31	59
50	S4	159/172 (92%)	154 (97%)	5 (3%)	35	62
51	S5	102/102 (100%)	102 (100%)	0	100	100
52	S6	130/130 (100%)	128 (98%)	2 (2%)	57	73
53	S7	165/171 (96%)	163 (99%)	2 (1%)	63	76
54	S8	160/187 (86%)	157 (98%)	3 (2%)	50	71
56	V1	412/427 (96%)	404 (98%)	8 (2%)	50	71
57	V2	190/190 (100%)	180 (95%)	10 (5%)	20	49
58	E7	192/192 (100%)	189 (98%)	3 (2%)	55	73
59	AC	80/111 (72%)	77 (96%)	3 (4%)	29	58
All	All	11801/13051 (90%)	11603 (98%)	198 (2%)	52	72

5 of 198 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
44	N2	96	ASN
47	N5	531	ASN
45	N3	233	ILE
46	N4	397	GLU
48	S2	133	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 217 such sidechains are listed below:

Mol	Chain	Res	Type
33	E6	57	ASN
44	N2	55	ASN
50	S4	39	ASN
34	E8	49	HIS
40	G1	252	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
48	2MR	S2	154	48	10,12,13	2.50	2 (20%)	5,13,15	0.89	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
48	2MR	S2	154	48	-	2/10/13/15	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
48	S2	154	2MR	CZ-NH2	5.48	1.44	1.33
48	S2	154	2MR	CZ-NE	5.10	1.45	1.34

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
48	S2	154	2MR	CG-CD-NE-CZ
48	S2	154	2MR	N-CA-CB-CG

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 64 ligands modelled in this entry, 3 are monoatomic - leaving 61 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
63	PC1	N4	503	-	32,32,53	0.36	0	38,40,61	0.35	0
61	SF4	S8	297	54	0,12,12	-	-	-		
64	CDL	C4	204	-	68,68,99	0.35	0	74,80,111	0.32	0
63	PC1	N2	301	-	36,36,53	0.35	0	42,44,61	0.36	0
64	CDL	AM	217	-	71,71,99	0.35	0	77,83,111	0.34	0
67	3PE	N4	504	-	40,40,50	0.33	0	43,45,55	0.31	0
63	PC1	E8	303	-	32,32,53	0.36	0	38,40,61	0.37	0
63	PC1	N5	606	-	35,35,53	0.35	0	41,43,61	0.32	0
64	CDL	AM	216	-	71,71,99	0.35	0	77,83,111	0.34	0
60	FES	V2	301	57	0,4,4	-	-	-		
65	NDP	A9	559	-	51,52,52	0.33	0	71,80,80	0.39	0
61	SF4	1A	403	1	0,12,12	-	-	-		
66	ZMP	AC	201	59	30,35,36	0.71	1 (3%)	34,42,45	0.81	1 (2%)
64	CDL	B5	201	-	57,57,99	0.38	0	63,69,111	0.34	0
63	PC1	AL	301	-	49,49,53	0.31	0	55,57,61	0.30	0
64	CDL	AM	215	-	71,71,99	0.35	0	77,83,111	0.31	0
63	PC1	A1	203	-	30,30,53	0.37	0	36,38,61	0.35	0
63	PC1	A9	561	-	32,32,53	0.36	0	38,40,61	0.36	0
61	SF4	S7	301	53	0,12,12	-	-	-		
63	PC1	AN	301	-	47,47,53	0.31	0	53,55,61	0.32	0
63	PC1	AM	218	-	48,48,53	0.31	0	54,56,61	0.30	0
63	PC1	A9	560	-	32,32,53	0.37	0	38,40,61	0.36	0
64	CDL	AL	303	-	63,63,99	0.36	0	69,75,111	0.31	0
64	CDL	C4	202	-	93,93,99	0.32	0	99,105,111	0.31	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
64	CDL	AL	304	-	69,69,99	0.35	0	75,81,111	0.33	0
67	3PE	N5	607	-	50,50,50	0.30	0	53,55,55	0.31	0
63	PC1	B5	203	-	53,53,53	0.30	0	59,61,61	0.33	0
63	PC1	E8	304	-	29,29,53	0.38	0	35,37,61	0.37	0
61	SF4	V1	580	56	0,12,12	-	-	-		
64	CDL	E7	301	-	67,67,99	0.35	0	73,79,111	0.31	0
67	3PE	G1	516	-	39,39,50	0.34	0	42,44,55	0.31	0
63	PC1	N1	701	-	48,48,53	0.30	0	54,56,61	0.30	0
63	PC1	N1	702	-	39,39,53	0.33	0	45,47,61	0.31	0
63	PC1	N5	601	-	53,53,53	0.29	0	59,61,61	0.32	0
63	PC1	N3	301	-	41,41,53	0.33	0	47,49,61	0.35	0
70	FMN	V1	579	-	33,33,33	0.28	0	48,50,50	0.32	0
64	CDL	A3	201	-	57,57,99	0.38	0	63,69,111	0.35	0
61	SF4	S8	298	54	0,12,12	-	-	-		
63	PC1	A1	202	-	48,48,53	0.31	0	54,56,61	0.33	0
63	PC1	C4	203	-	37,37,53	0.34	0	43,45,61	0.32	0
64	CDL	AL	302	-	67,67,99	0.36	0	73,79,111	0.31	0
63	PC1	ED	201	-	53,53,53	0.29	0	59,61,61	0.30	0
64	CDL	E6	431	-	63,63,99	0.37	0	69,75,111	0.32	0
64	CDL	B3	102	-	64,64,99	0.37	0	70,76,111	0.35	0
64	CDL	N5	603	-	69,69,99	0.35	0	75,81,111	0.32	0
61	SF4	1A	402	1	0,12,12	-	-	-		
68	U10	N4	505	-	43,43,63	2.47	20 (46%)	54,55,79	1.67	16 (29%)
63	PC1	B5	202	-	53,53,53	0.29	0	59,61,61	0.33	0
63	PC1	E8	301	-	53,53,53	0.29	0	59,61,61	0.31	0
60	FES	1A	401	1	0,4,4	-	-	-		
63	PC1	N4	502	-	38,38,53	0.34	0	44,46,61	0.34	0
63	PC1	N5	605	-	40,40,53	0.33	0	46,48,61	0.30	0
63	PC1	AM	220	-	47,47,53	0.30	0	53,55,61	0.28	0
67	3PE	AN	302	-	50,50,50	0.30	0	53,55,55	0.30	0
64	CDL	N4	501	-	97,97,99	0.31	0	103,109,111	0.27	0
66	ZMP	AB	150	13	30,35,36	0.72	1 (3%)	34,42,45	0.80	1 (2%)
63	PC1	E8	302	-	53,53,53	0.29	0	59,61,61	0.30	0
64	CDL	N5	608	-	92,92,99	0.31	0	98,104,111	0.29	0
63	PC1	E4	401	-	50,50,53	0.30	0	56,58,61	0.33	0
64	CDL	EA	201	-	58,58,99	0.39	0	64,70,111	0.33	0
64	CDL	EA	202	-	54,54,99	0.39	0	60,66,111	0.34	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
63	PC1	N4	503	-	-	5/36/36/57	-
61	SF4	S8	297	54	-	-	0/6/5/5
64	CDL	C4	204	-	-	12/79/79/110	-
63	PC1	N2	301	-	-	8/40/40/57	-
64	CDL	AM	217	-	-	21/82/82/110	-
67	3PE	N4	504	-	-	11/44/44/54	-
63	PC1	E8	303	-	-	8/36/36/57	-
63	PC1	N5	606	-	-	10/39/39/57	-
64	CDL	AM	216	-	-	16/82/82/110	-
60	FES	V2	301	57	-	-	0/1/1/1
65	NDP	A9	559	-	-	2/34/77/77	0/5/5/5
61	SF4	1A	403	1	-	-	0/6/5/5
66	ZMP	AC	201	59	-	23/40/42/43	-
64	CDL	B5	201	-	-	13/68/68/110	-
63	PC1	AL	301	-	-	9/53/53/57	-
64	CDL	AM	215	-	-	15/82/82/110	-
63	PC1	A1	203	-	-	6/34/34/57	-
63	PC1	A9	561	-	-	9/36/36/57	-
61	SF4	S7	301	53	-	-	0/6/5/5
63	PC1	AN	301	-	-	12/51/51/57	-
63	PC1	AM	218	-	-	19/52/52/57	-
63	PC1	A9	560	-	-	11/36/36/57	-
64	CDL	AL	303	-	-	11/74/74/110	-
64	CDL	C4	202	-	-	20/104/104/110	-
64	CDL	AL	304	-	-	24/80/80/110	-
67	3PE	N5	607	-	-	15/54/54/54	-
63	PC1	B5	203	-	-	16/57/57/57	-
63	PC1	E8	304	-	-	11/33/33/57	-
61	SF4	V1	580	56	-	-	0/6/5/5
64	CDL	E7	301	-	-	17/78/78/110	-
67	3PE	G1	516	-	-	4/43/43/54	-
63	PC1	N1	701	-	-	15/52/52/57	-
63	PC1	N1	702	-	-	11/43/43/57	-
63	PC1	N5	601	-	-	8/57/57/57	-
63	PC1	N3	301	-	-	9/45/45/57	-
70	FMN	V1	579	-	-	2/18/18/18	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
64	CDL	A3	201	-	-	7/68/68/110	-
63	PC1	A1	202	-	-	15/52/52/57	-
63	PC1	C4	203	-	-	11/41/41/57	-
61	SF4	S8	298	54	-	-	0/6/5/5
64	CDL	AL	302	-	-	9/78/78/110	-
63	PC1	ED	201	-	-	8/57/57/57	-
64	CDL	E6	431	-	-	22/74/74/110	-
64	CDL	B3	102	-	-	10/75/75/110	-
64	CDL	N5	603	-	-	10/80/80/110	-
68	U10	N4	505	-	-	8/39/63/87	0/1/1/1
61	SF4	1A	402	1	-	-	0/6/5/5
63	PC1	B5	202	-	-	16/57/57/57	-
63	PC1	E8	301	-	-	11/57/57/57	-
60	FES	1A	401	1	-	-	0/1/1/1
63	PC1	N4	502	-	-	16/42/42/57	-
63	PC1	N5	605	-	-	13/44/44/57	-
63	PC1	AM	220	-	-	9/51/51/57	-
67	3PE	AN	302	-	-	10/54/54/54	-
64	CDL	N4	501	-	-	24/108/108/110	-
66	ZMP	AB	150	13	-	16/40/42/43	-
63	PC1	E8	302	-	-	19/57/57/57	-
64	CDL	N5	608	-	-	16/103/103/110	-
63	PC1	E4	401	-	-	13/54/54/57	-
64	CDL	EA	201	-	-	12/69/69/110	-
64	CDL	EA	202	-	-	12/65/65/110	-

The worst 5 of 22 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
68	N4	505	U10	C6-C1	10.52	1.54	1.35
68	N4	505	U10	C4-C3	4.74	1.53	1.36
68	N4	505	U10	C7-C8	2.92	1.55	1.50
68	N4	505	U10	C26-C24	2.68	1.56	1.51
68	N4	505	U10	C31-C29	2.61	1.56	1.51

The worst 5 of 18 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
68	N4	505	U10	C7-C8-C9	-3.57	120.69	126.83
68	N4	505	U10	C30-C29-C31	3.43	121.19	115.23
68	N4	505	U10	C22-C23-C24	-3.32	120.03	127.62
68	N4	505	U10	C7-C6-C1	-3.11	119.55	124.89
68	N4	505	U10	C15-C14-C16	2.96	120.37	115.23

There are no chirality outliers.

5 of 660 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
63	A1	202	PC1	C11-O13-P-O11
63	A1	203	PC1	C11-O13-P-O12
63	A1	203	PC1	C11-O13-P-O14
63	A1	203	PC1	C1-O11-P-O12
63	A1	203	PC1	C1-O11-P-O13

There are no ring outliers.

33 monomers are involved in 78 short contacts:

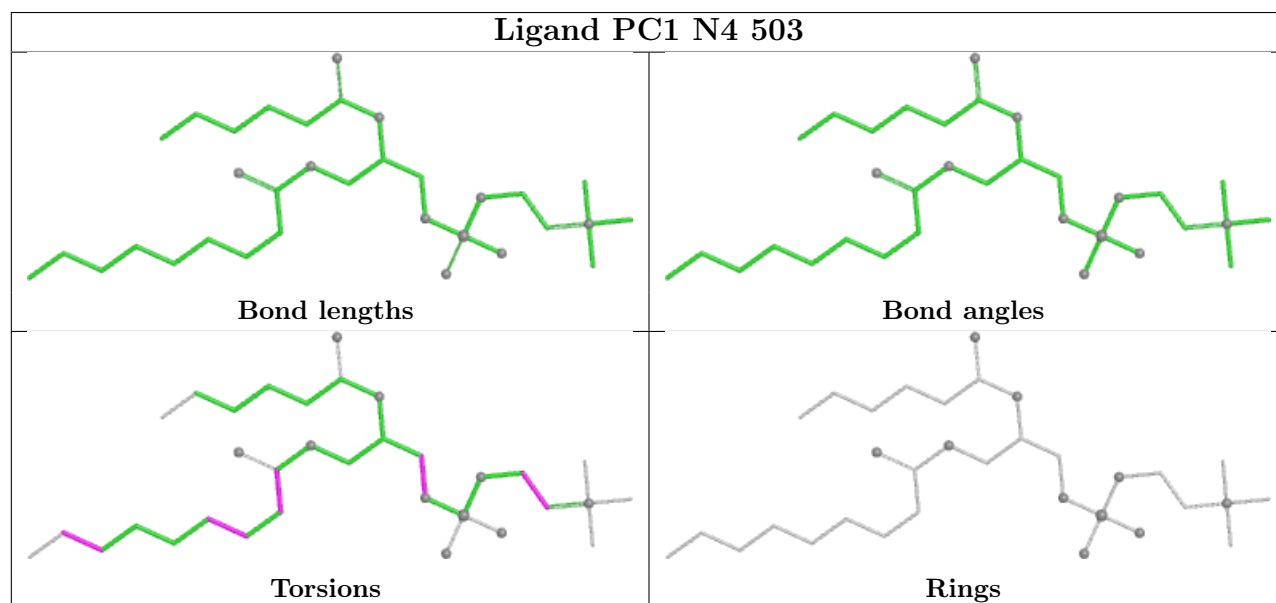
Mol	Chain	Res	Type	Clashes	Symm-Clashes
61	S8	297	SF4	3	0
63	N2	301	PC1	1	0
64	AM	217	CDL	2	0
64	AM	216	CDL	4	0
65	A9	559	NDP	3	0
66	AC	201	ZMP	15	0
64	B5	201	CDL	1	0
64	AM	215	CDL	4	0
63	A1	203	PC1	1	0
63	A9	560	PC1	1	0
64	AL	303	CDL	2	0
64	C4	202	CDL	1	0
67	N5	607	3PE	1	0
63	B5	203	PC1	2	0
61	V1	580	SF4	1	0
67	G1	516	3PE	2	0
63	N5	601	PC1	1	0
64	A3	201	CDL	1	0
61	S8	298	SF4	2	0
63	A1	202	PC1	2	0
63	C4	203	PC1	1	0
63	ED	201	PC1	1	0

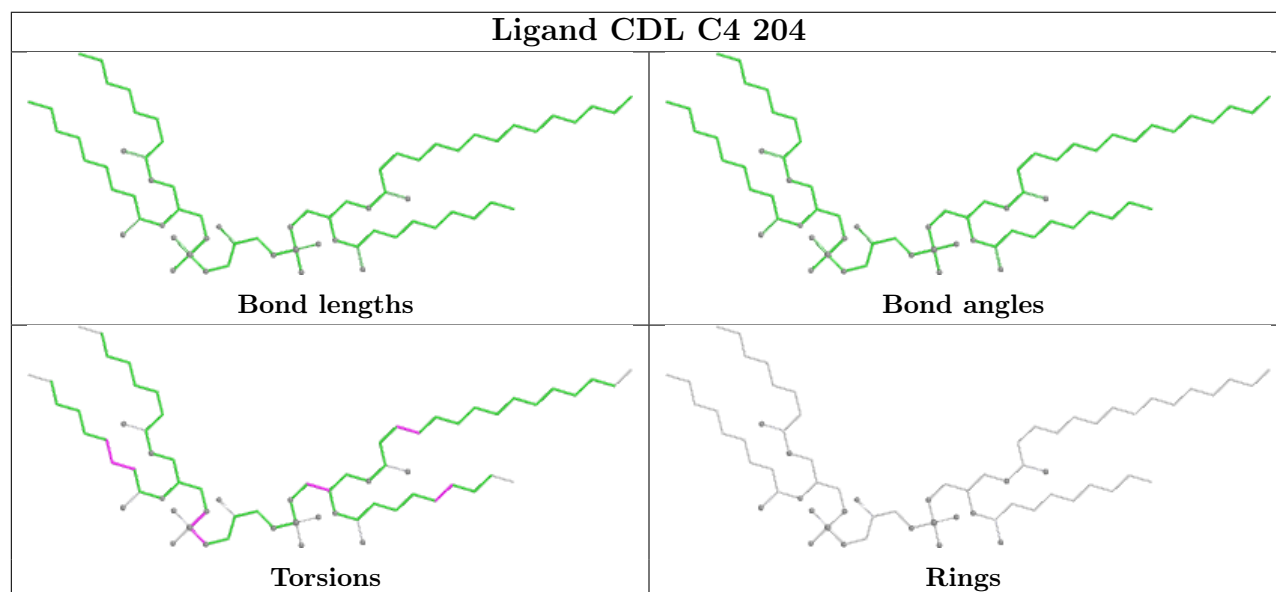
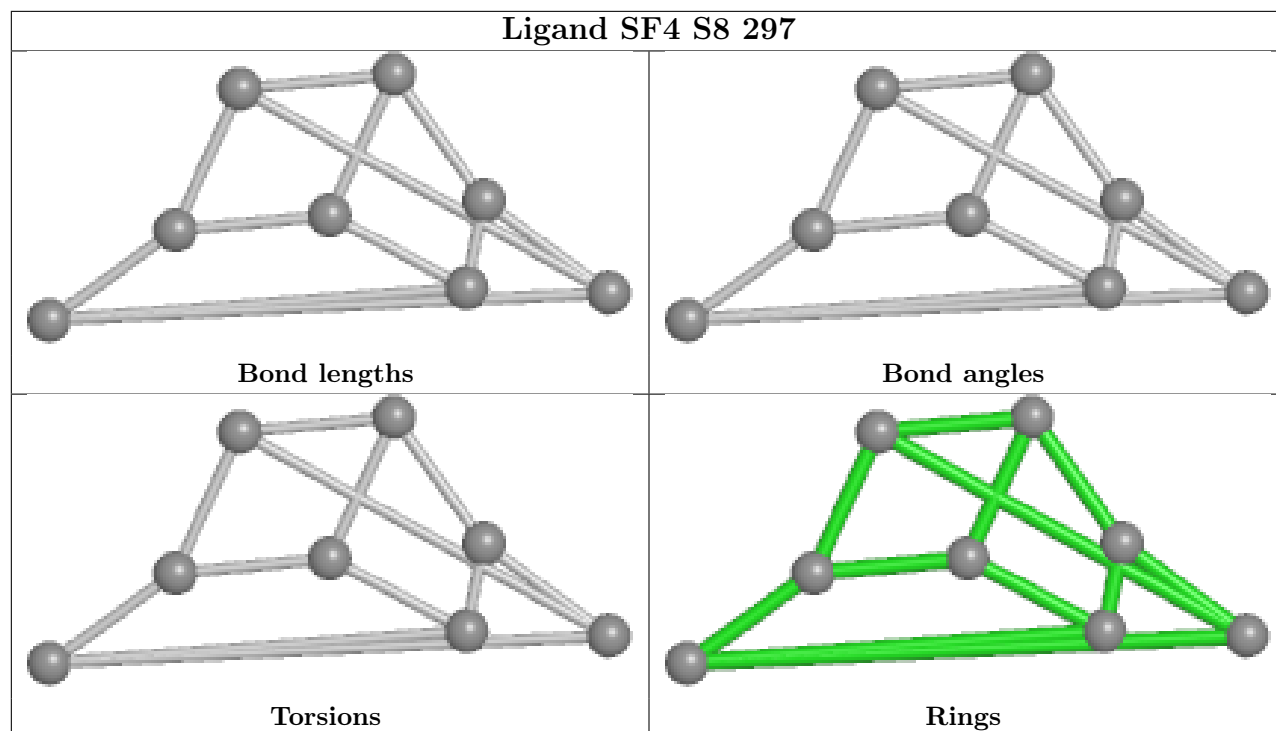
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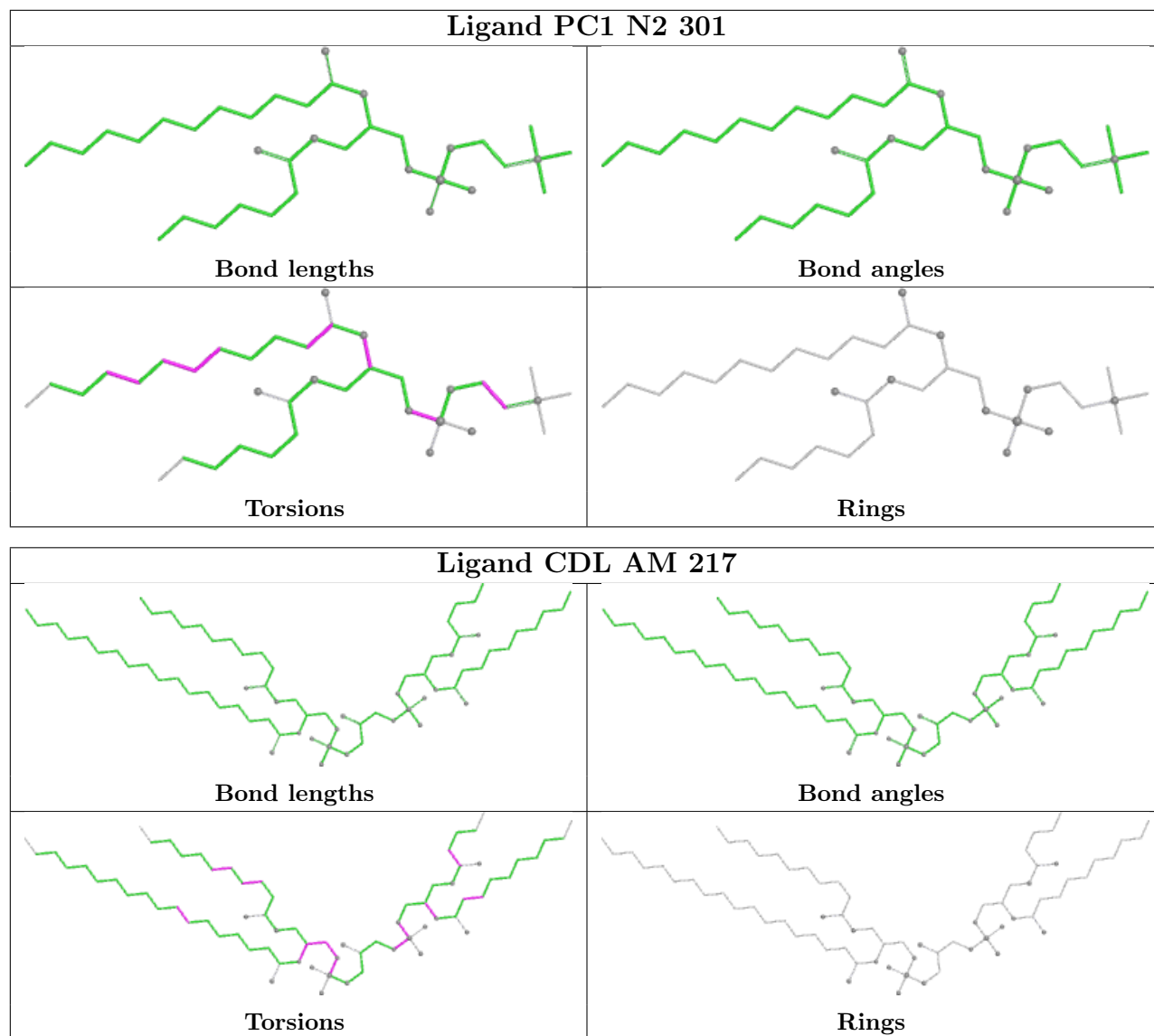
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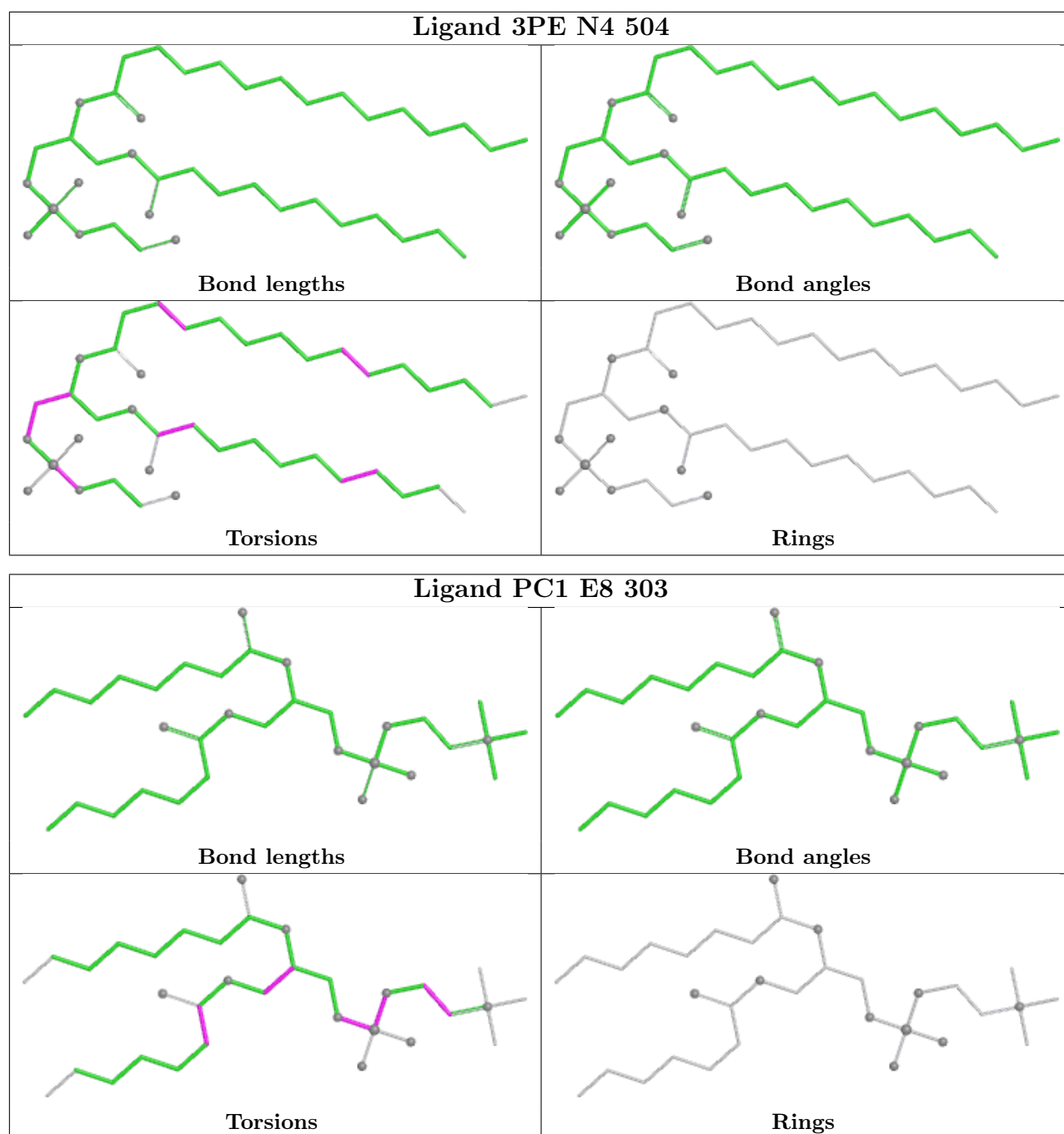
Mol	Chain	Res	Type	Clashes	Symm-Clashes
64	B3	102	CDL	1	0
64	N5	603	CDL	3	0
68	N4	505	U10	5	0
63	E8	301	PC1	2	0
63	N4	502	PC1	2	0
63	AM	220	PC1	1	0
64	N4	501	CDL	4	0
66	AB	150	ZMP	6	0
63	E8	302	PC1	1	0
64	N5	608	CDL	1	0
64	EA	202	CDL	1	0

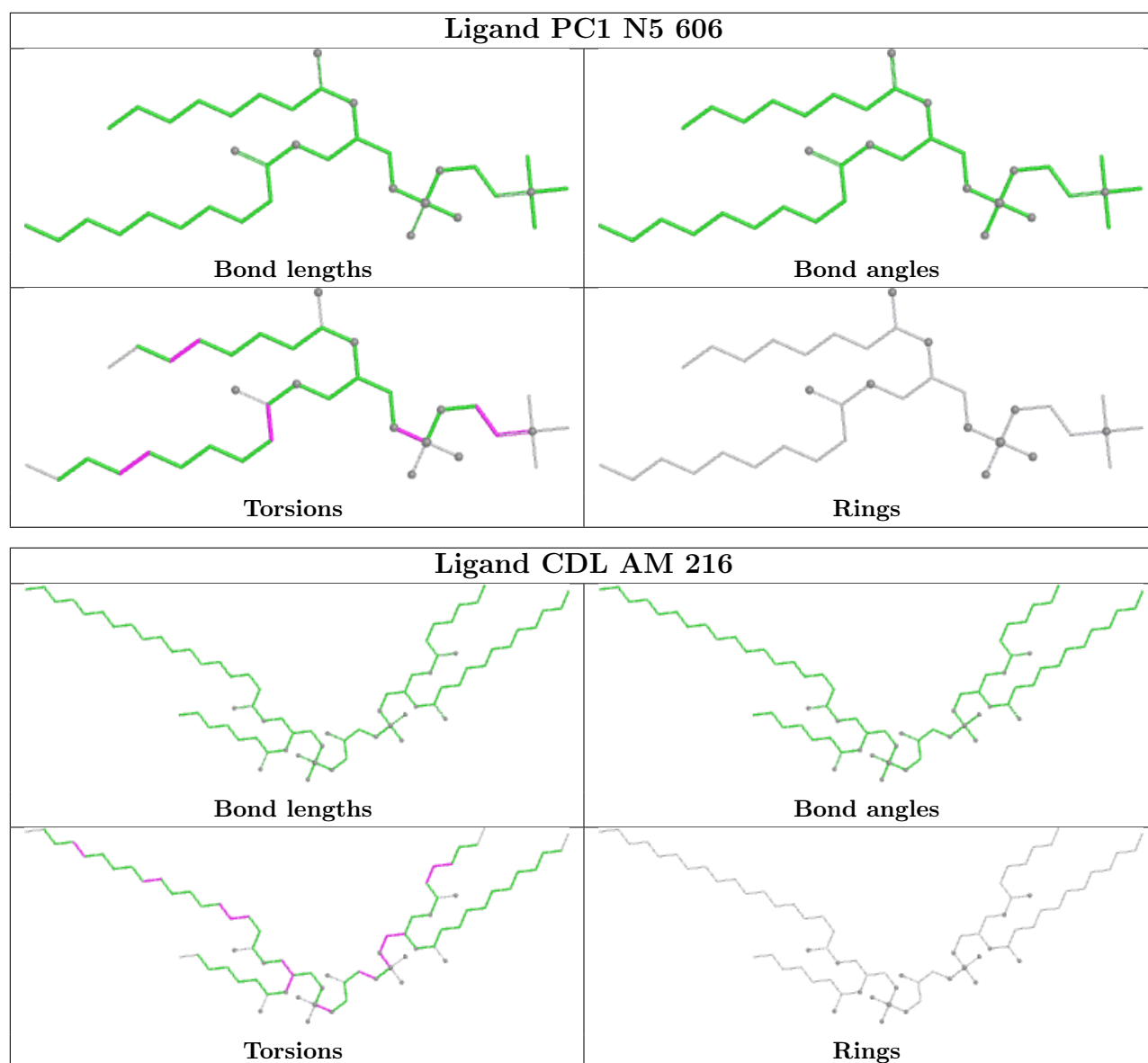
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

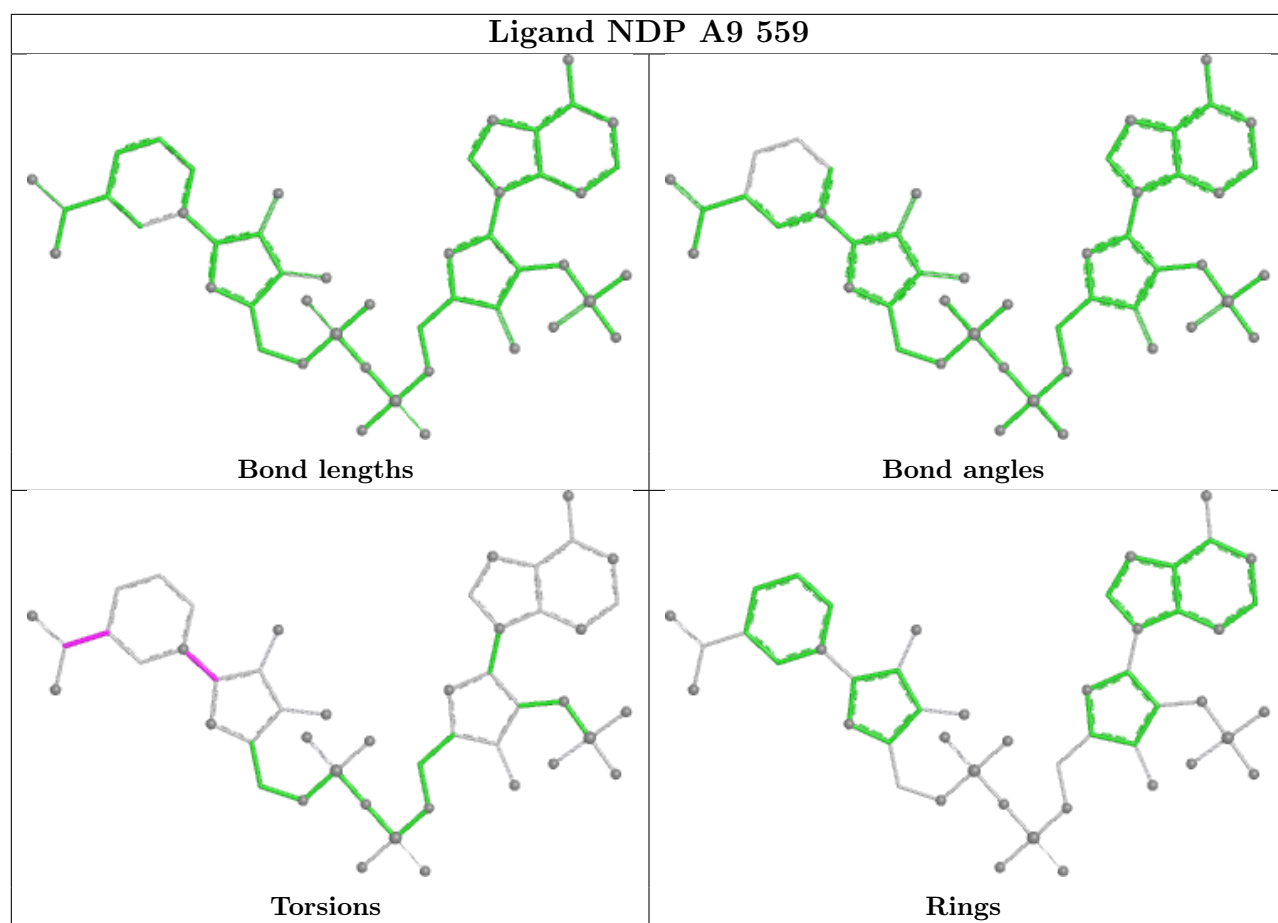


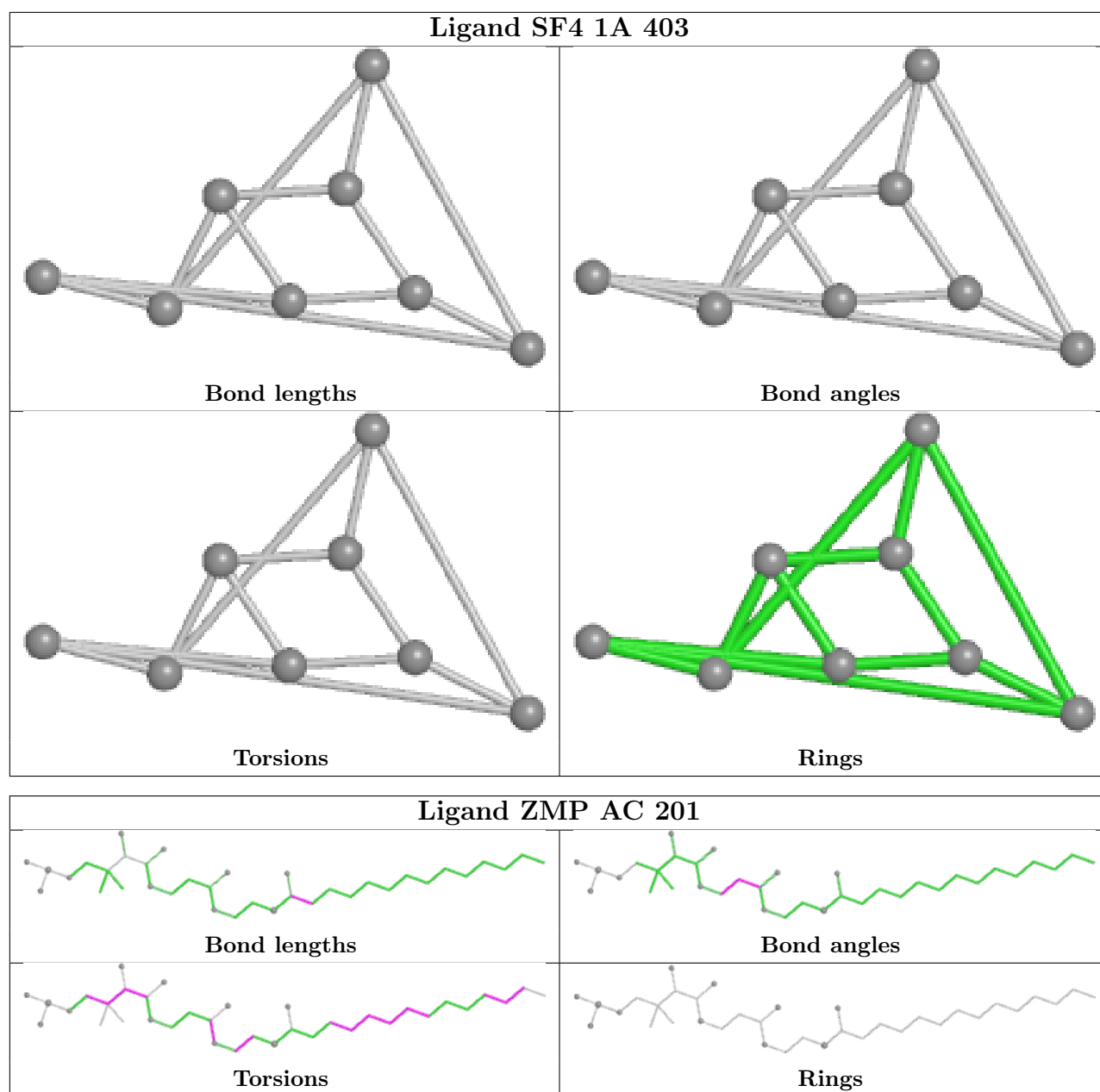


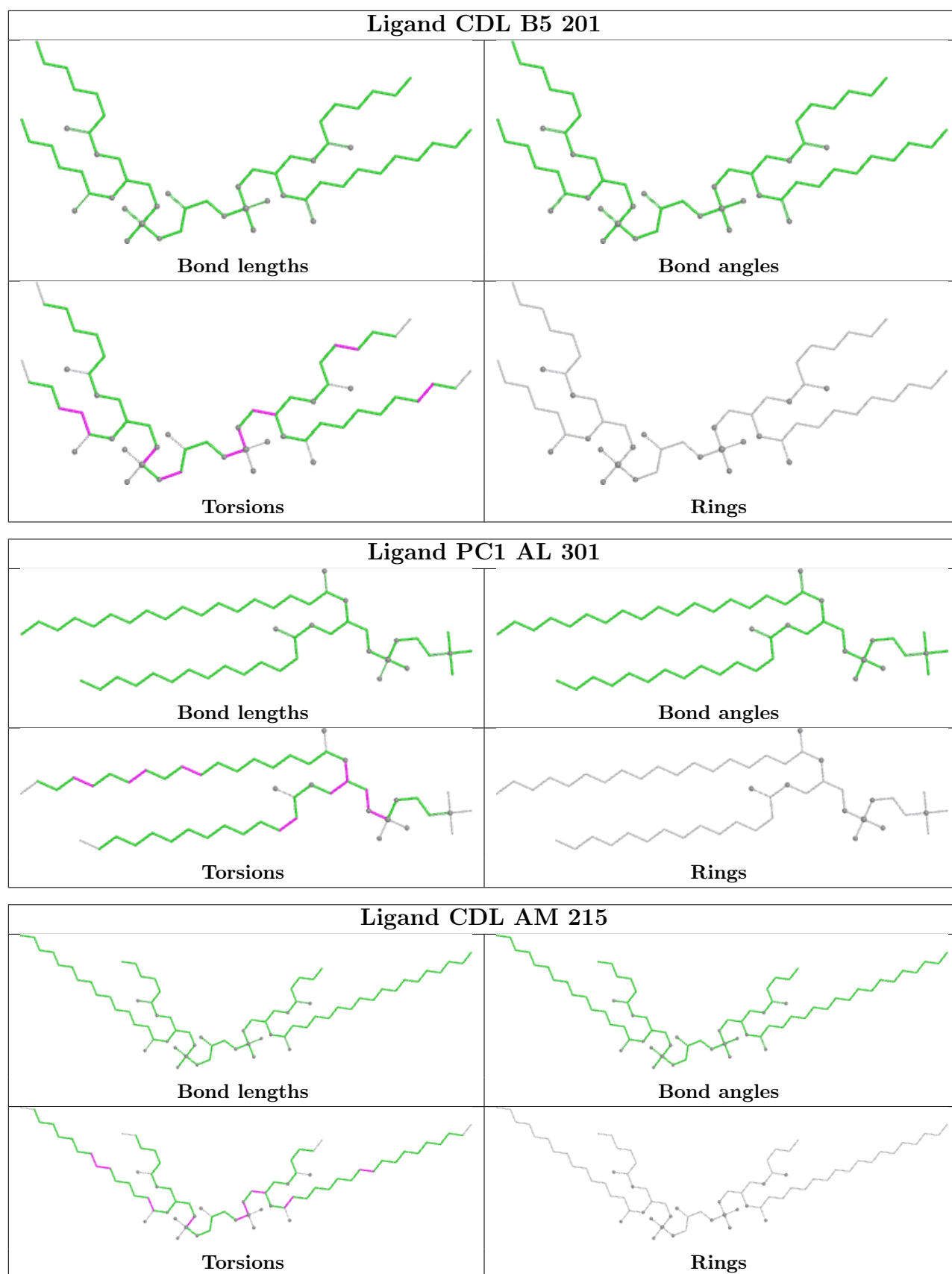


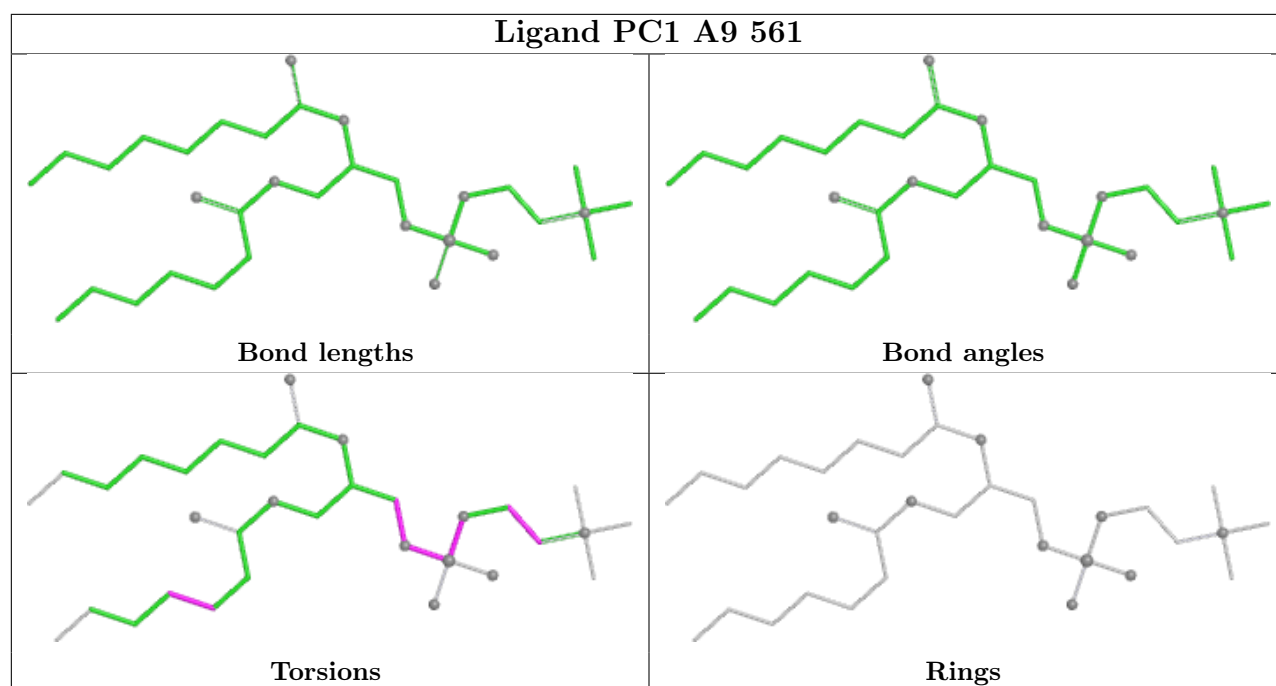
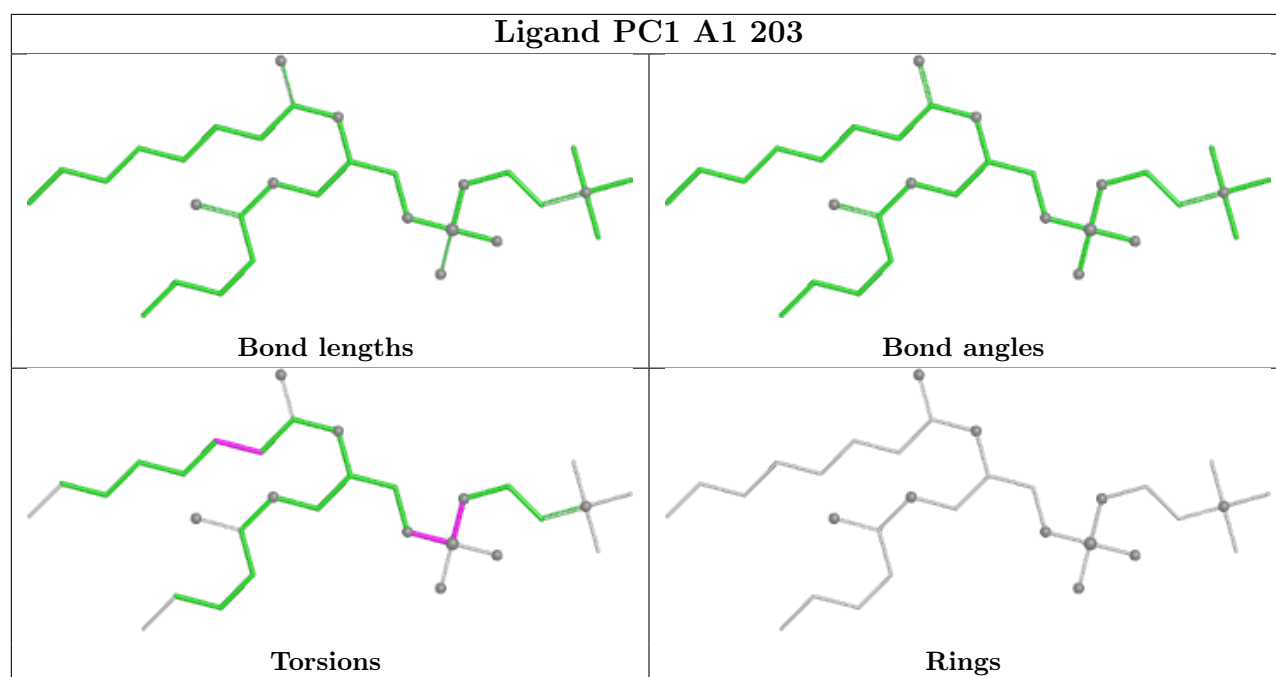


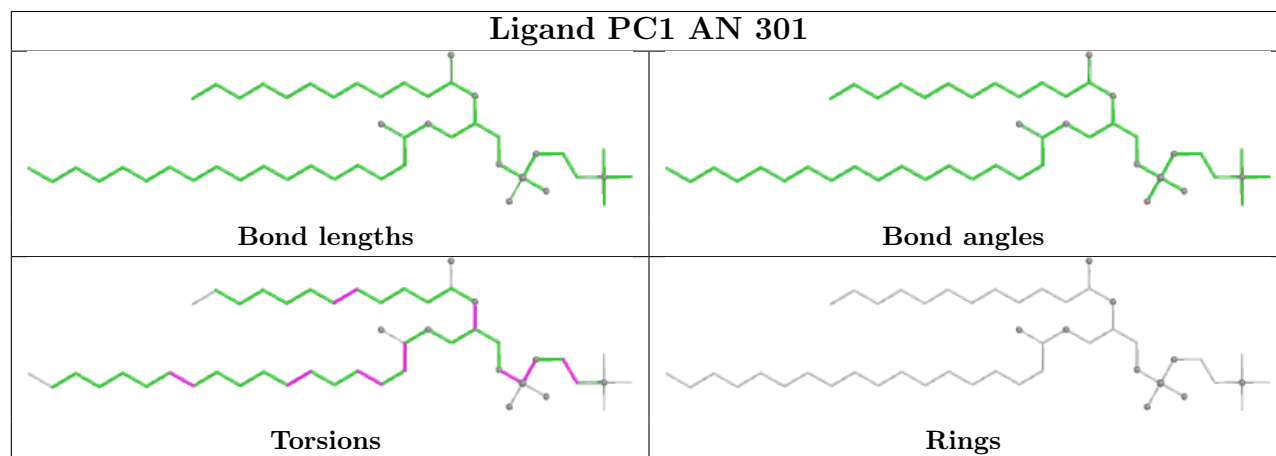
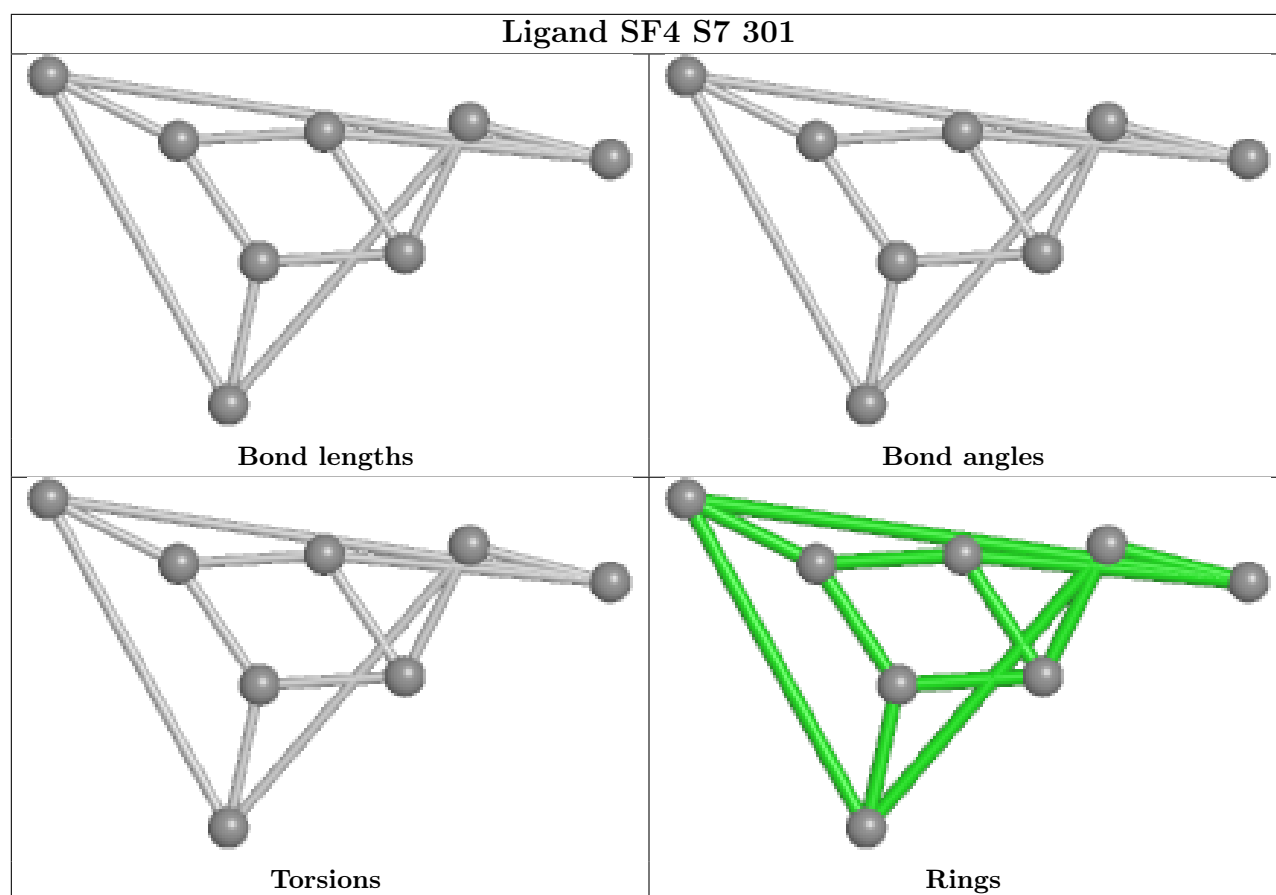


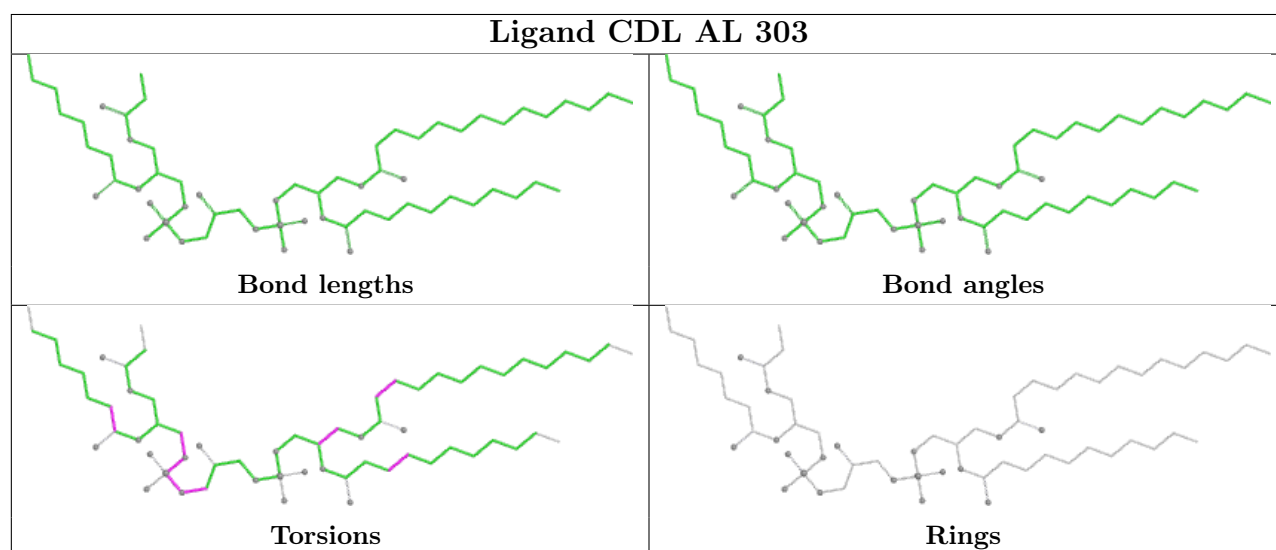
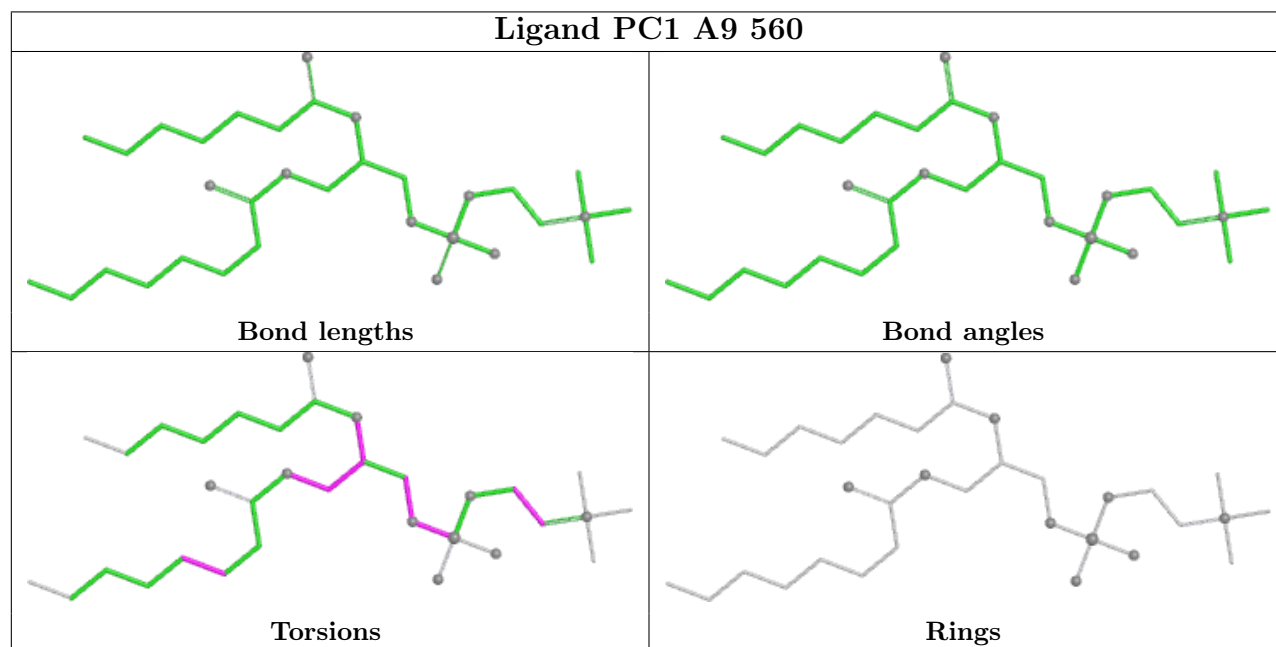
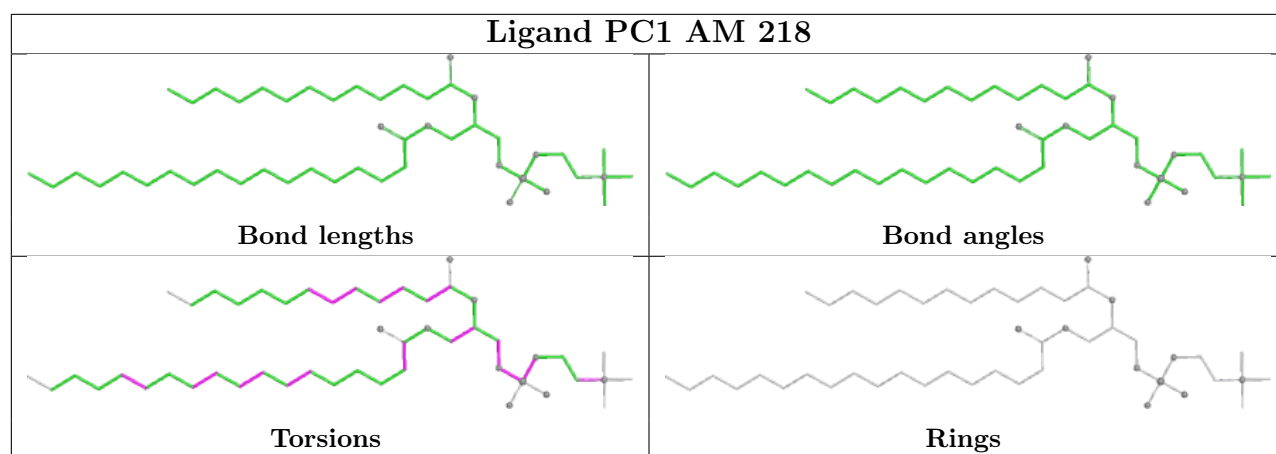


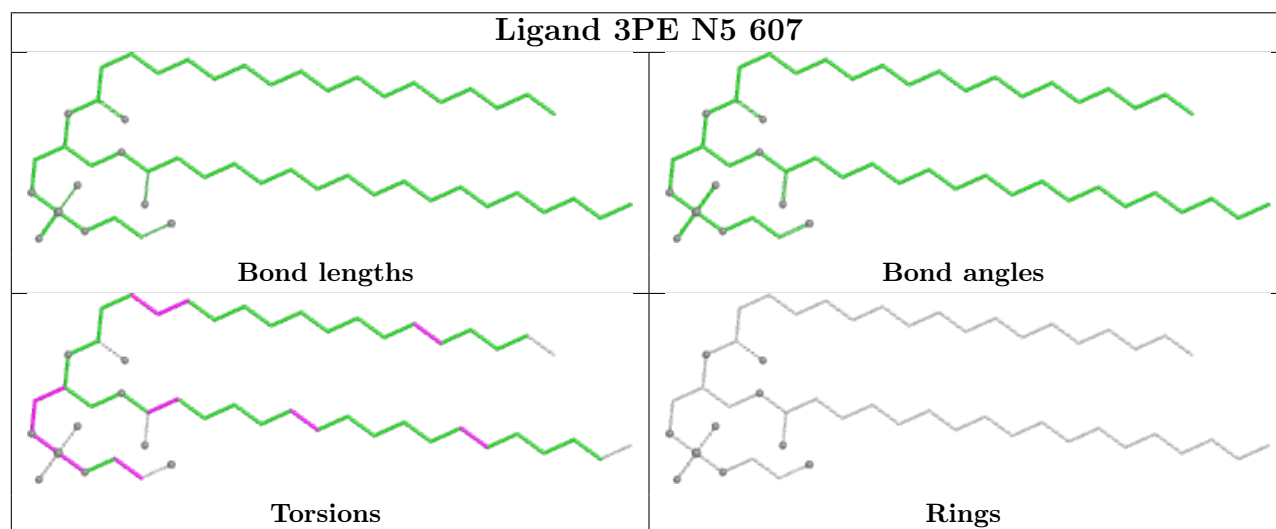
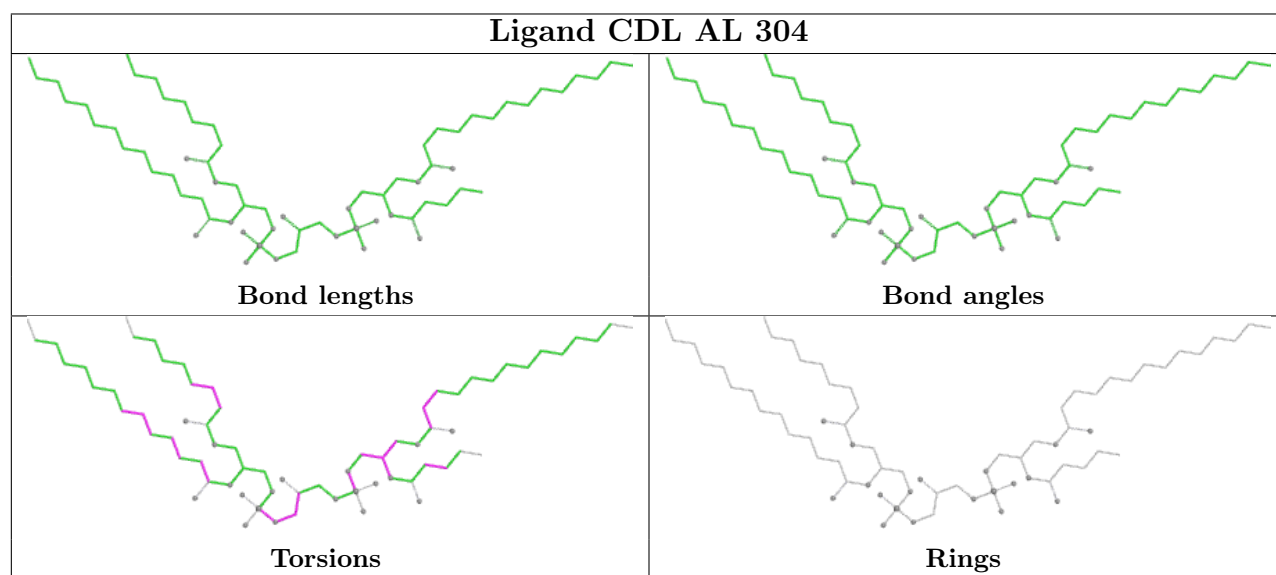
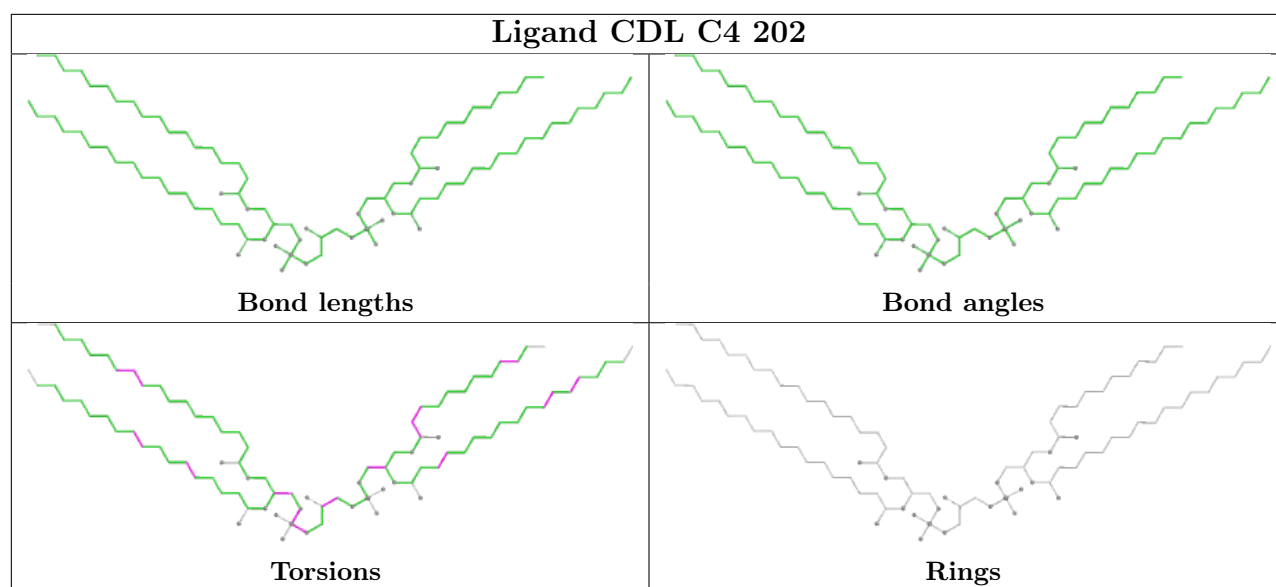


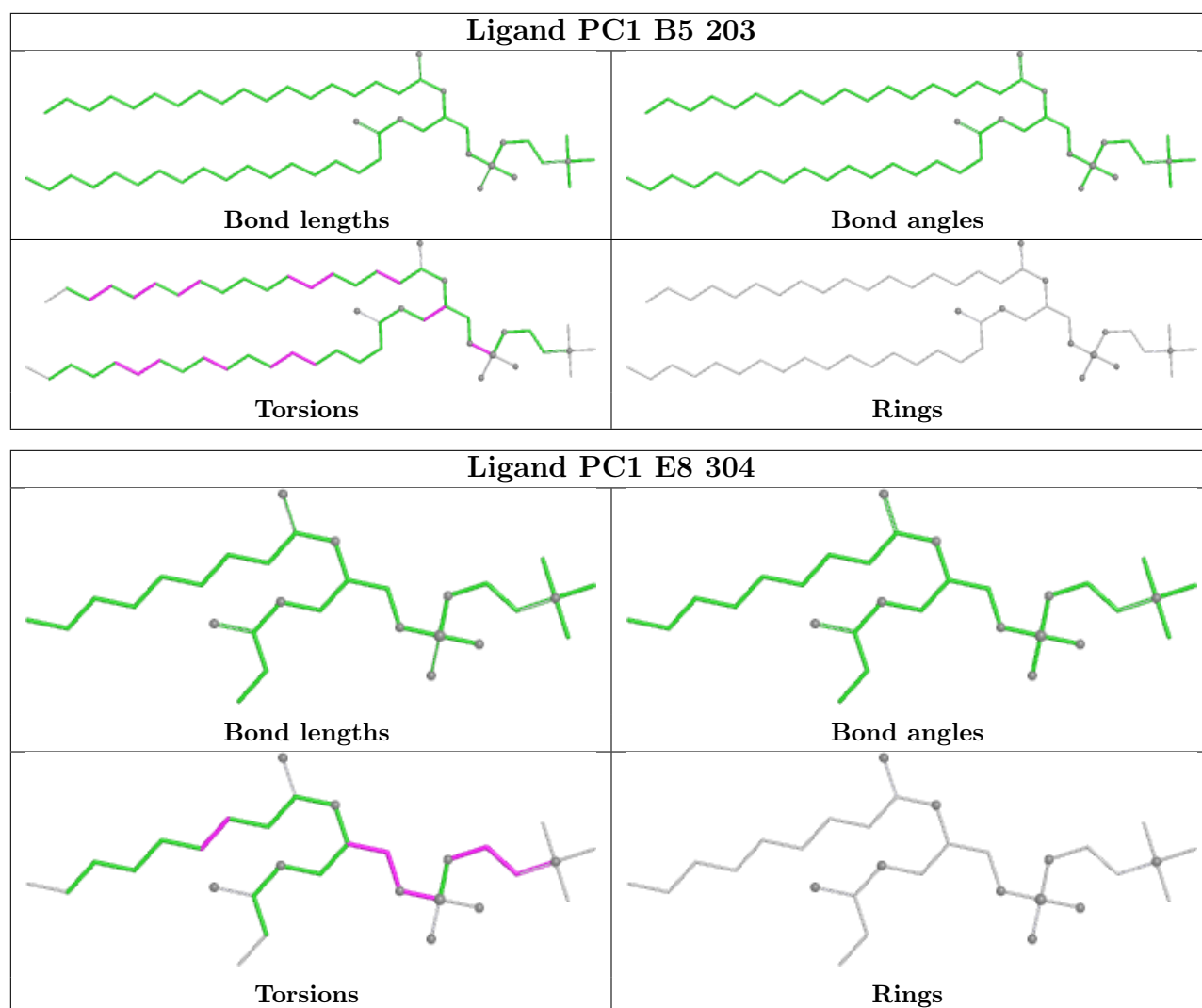


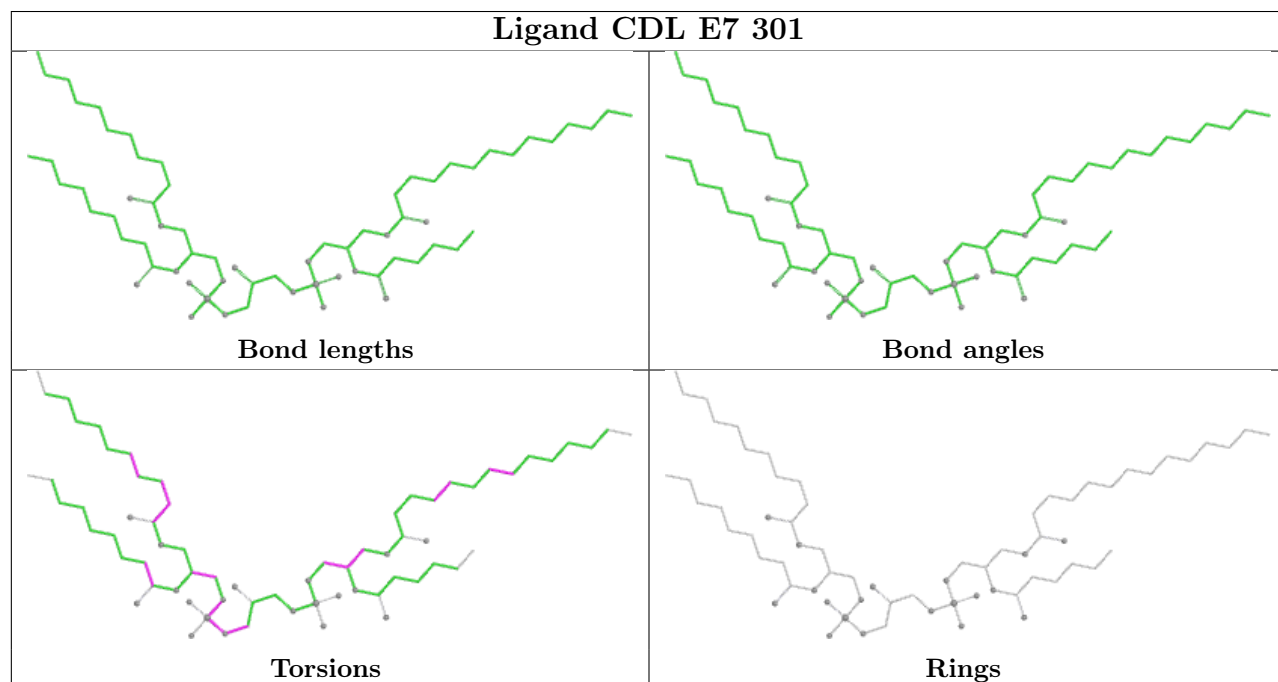
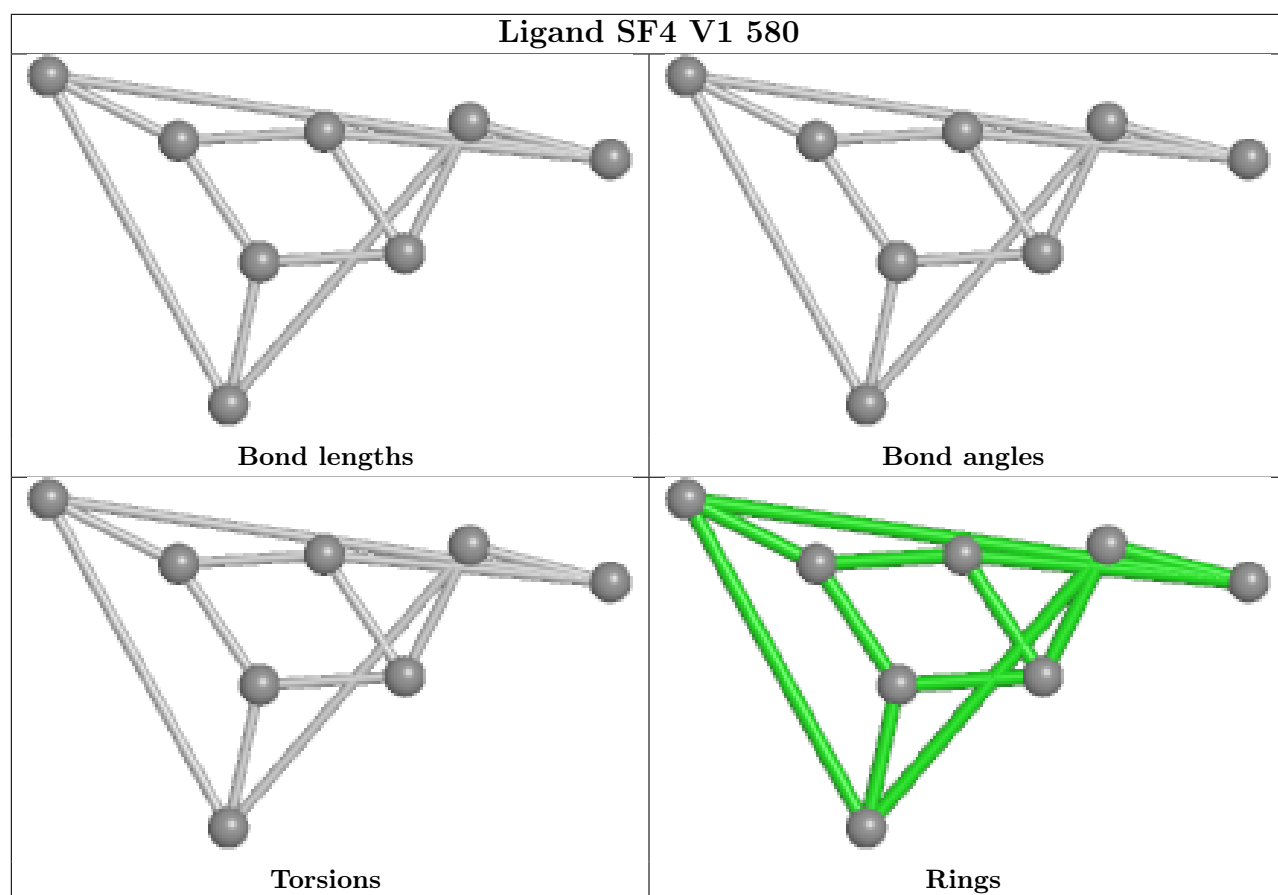


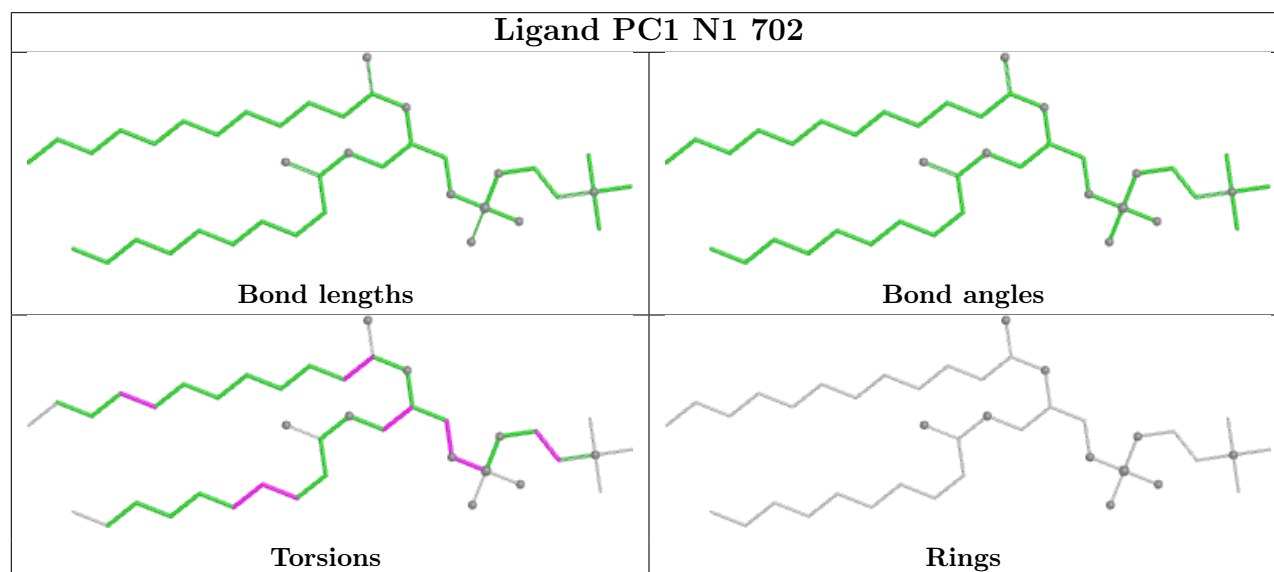
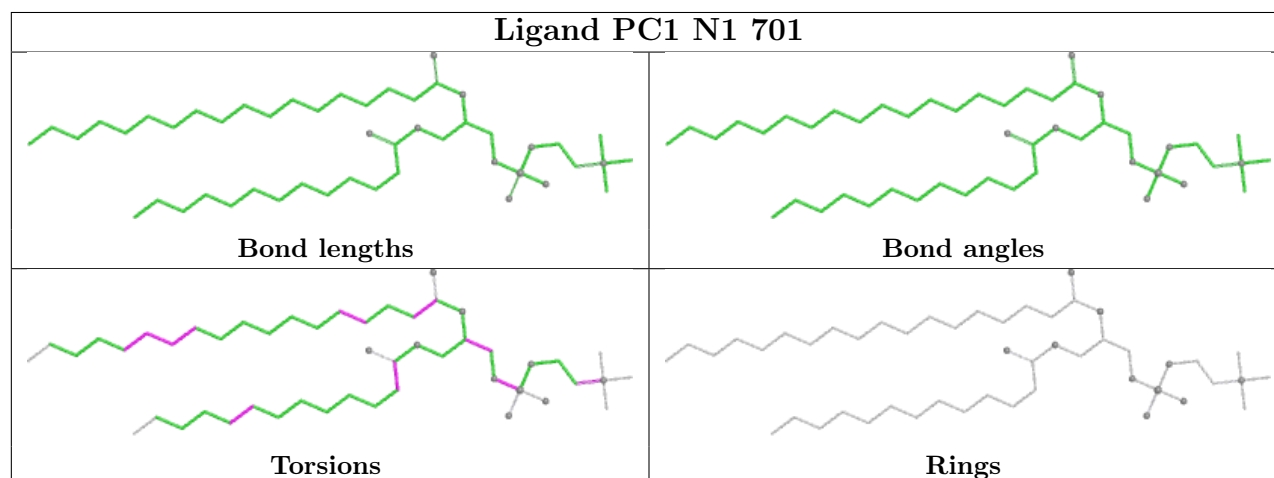
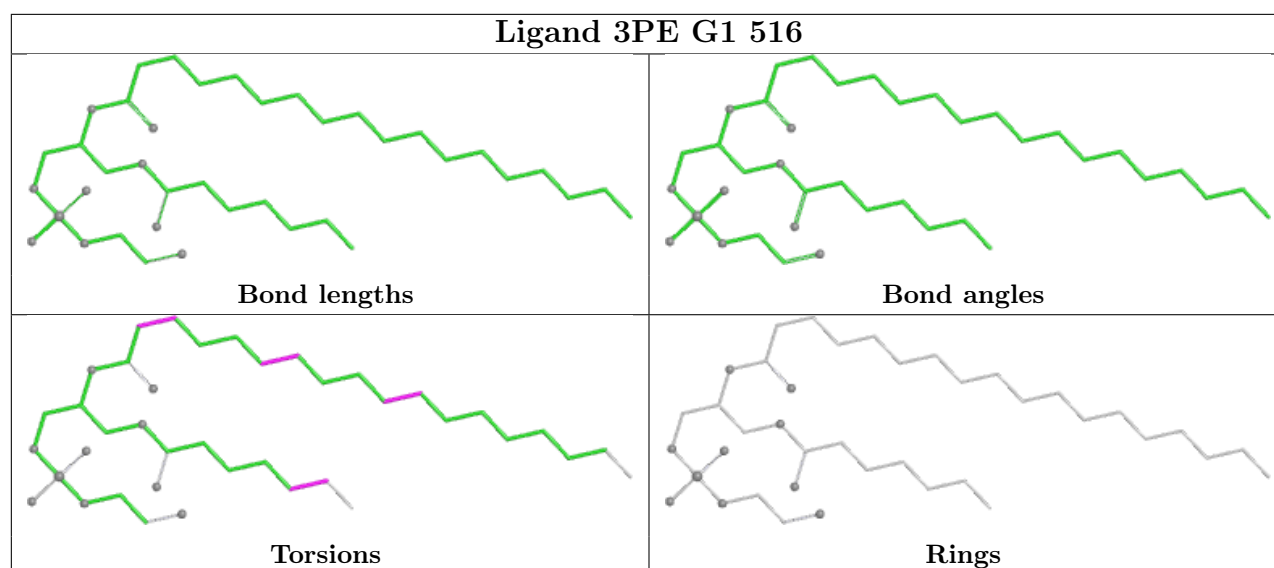


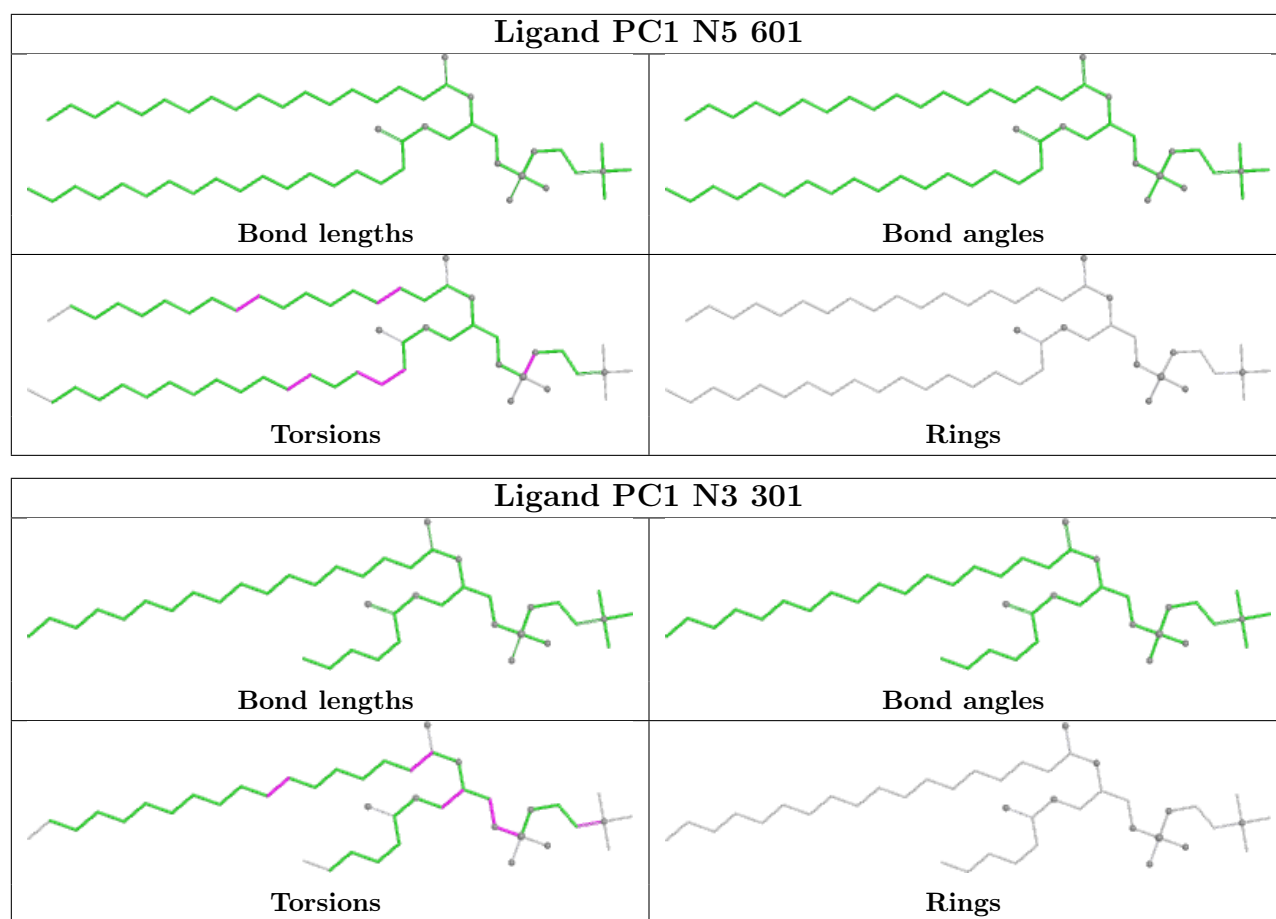


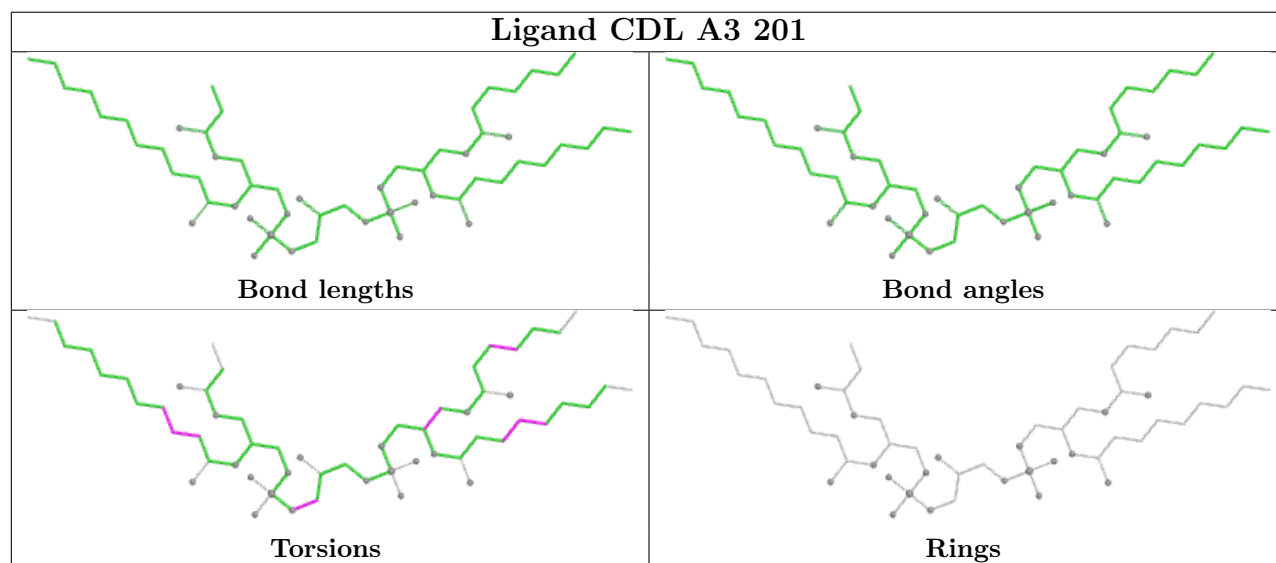
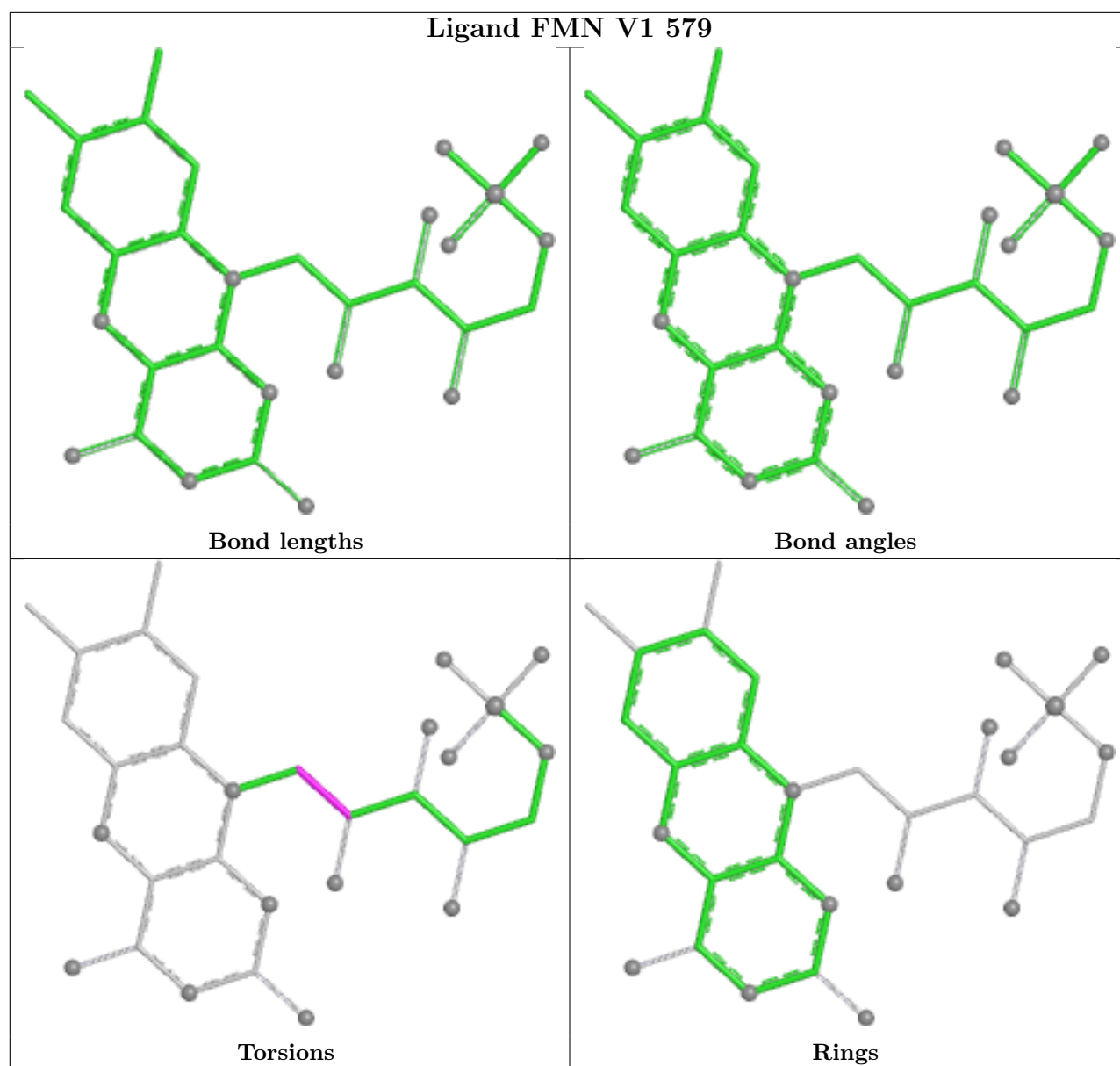


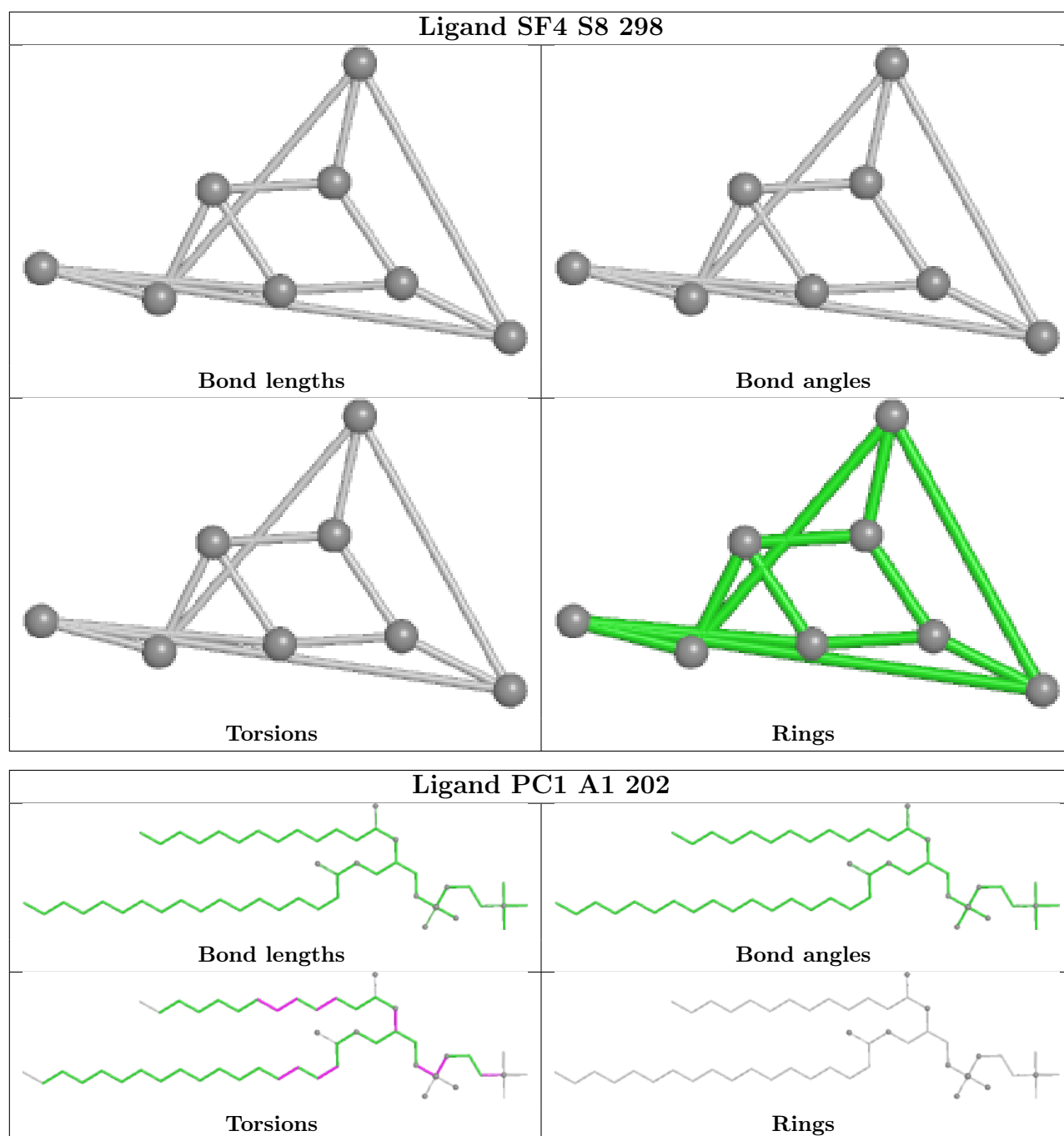


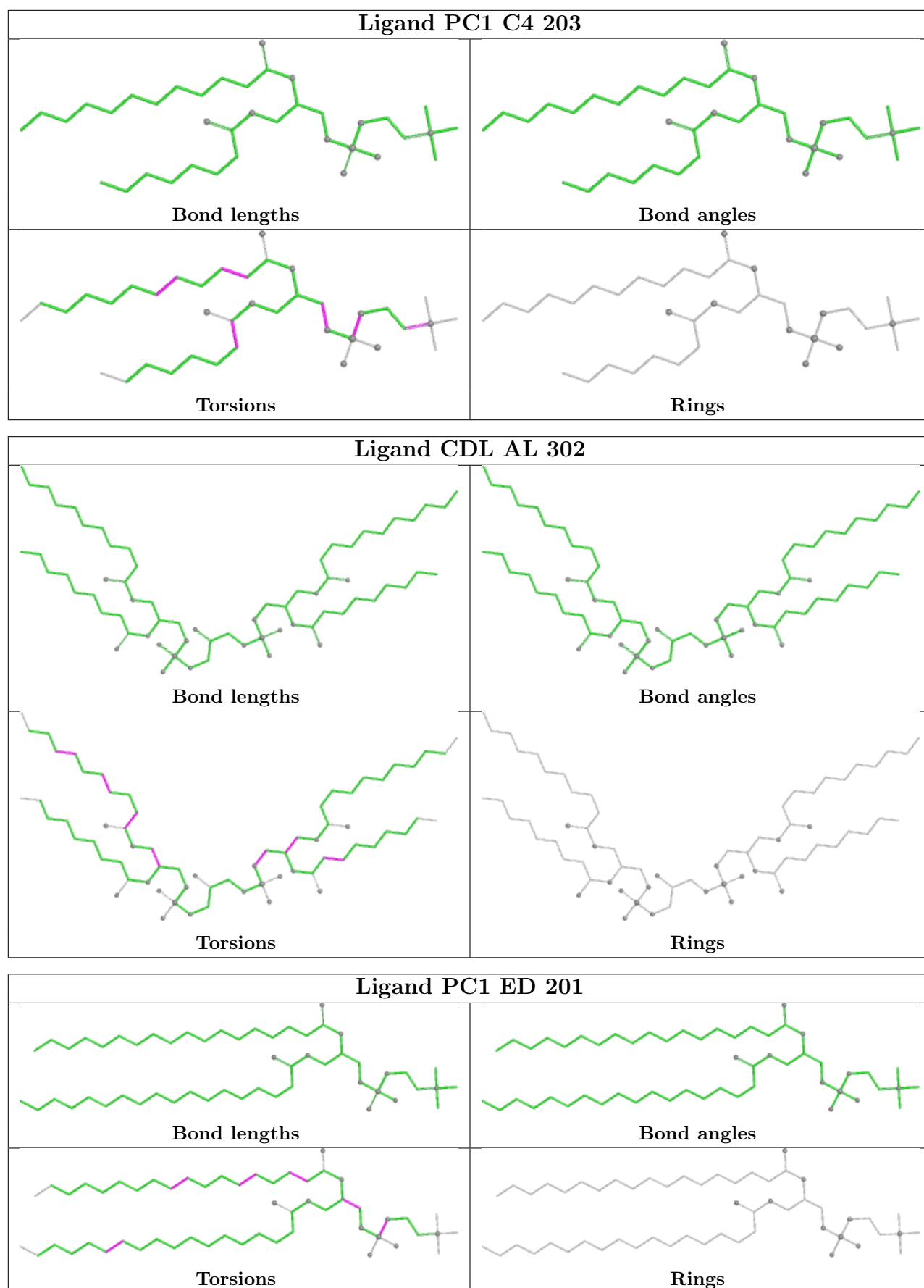


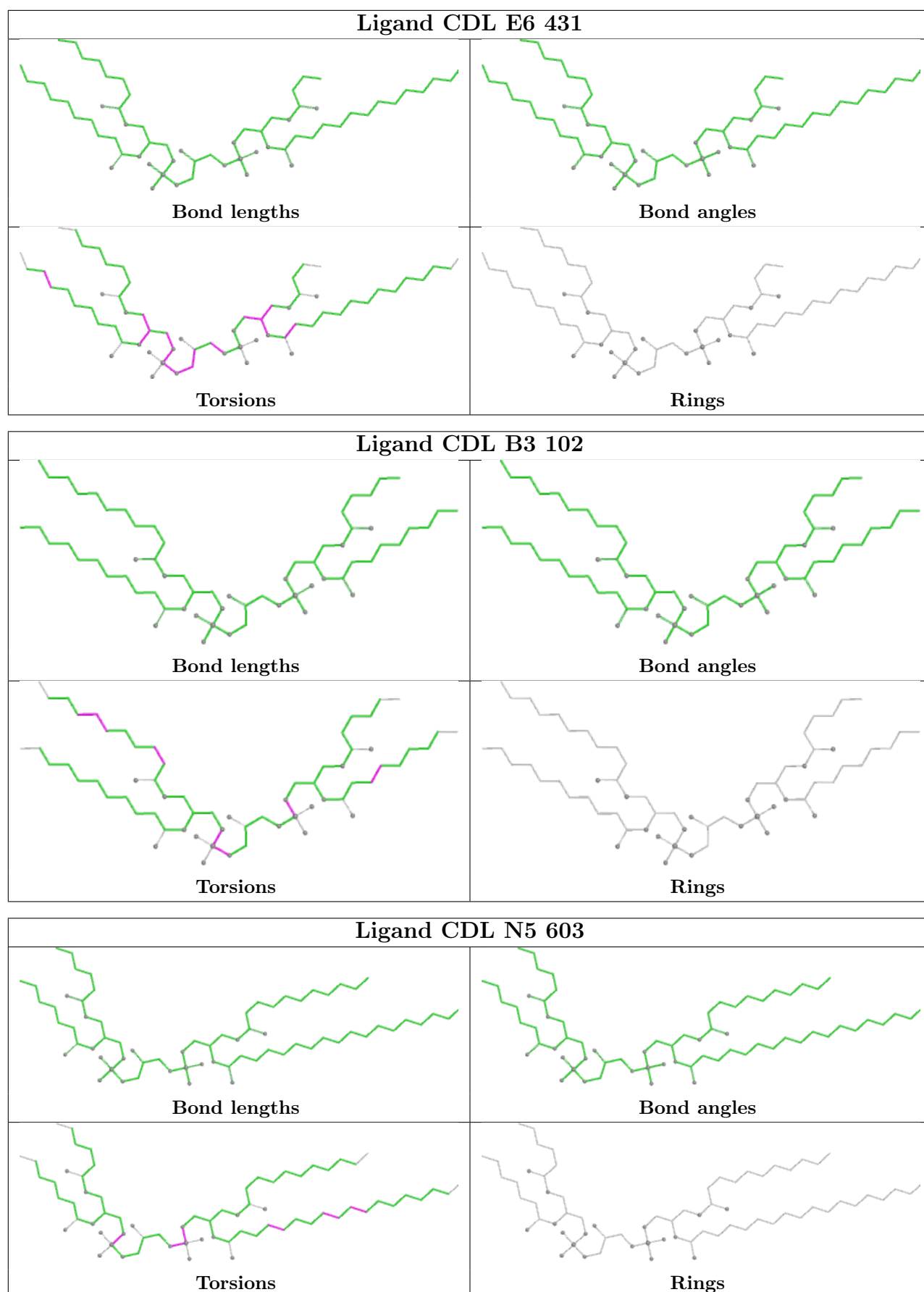


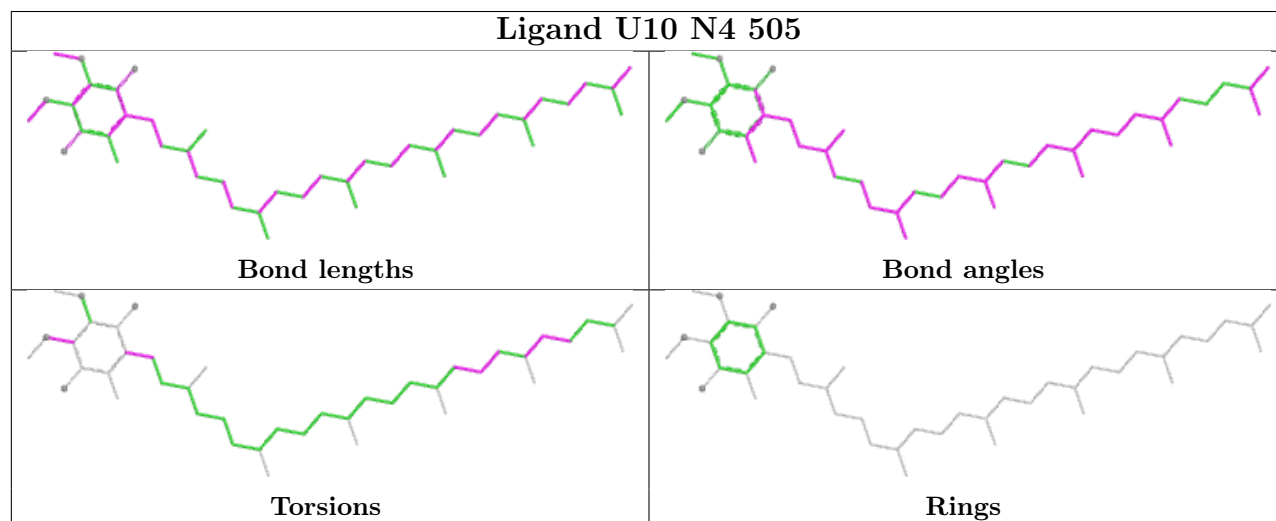
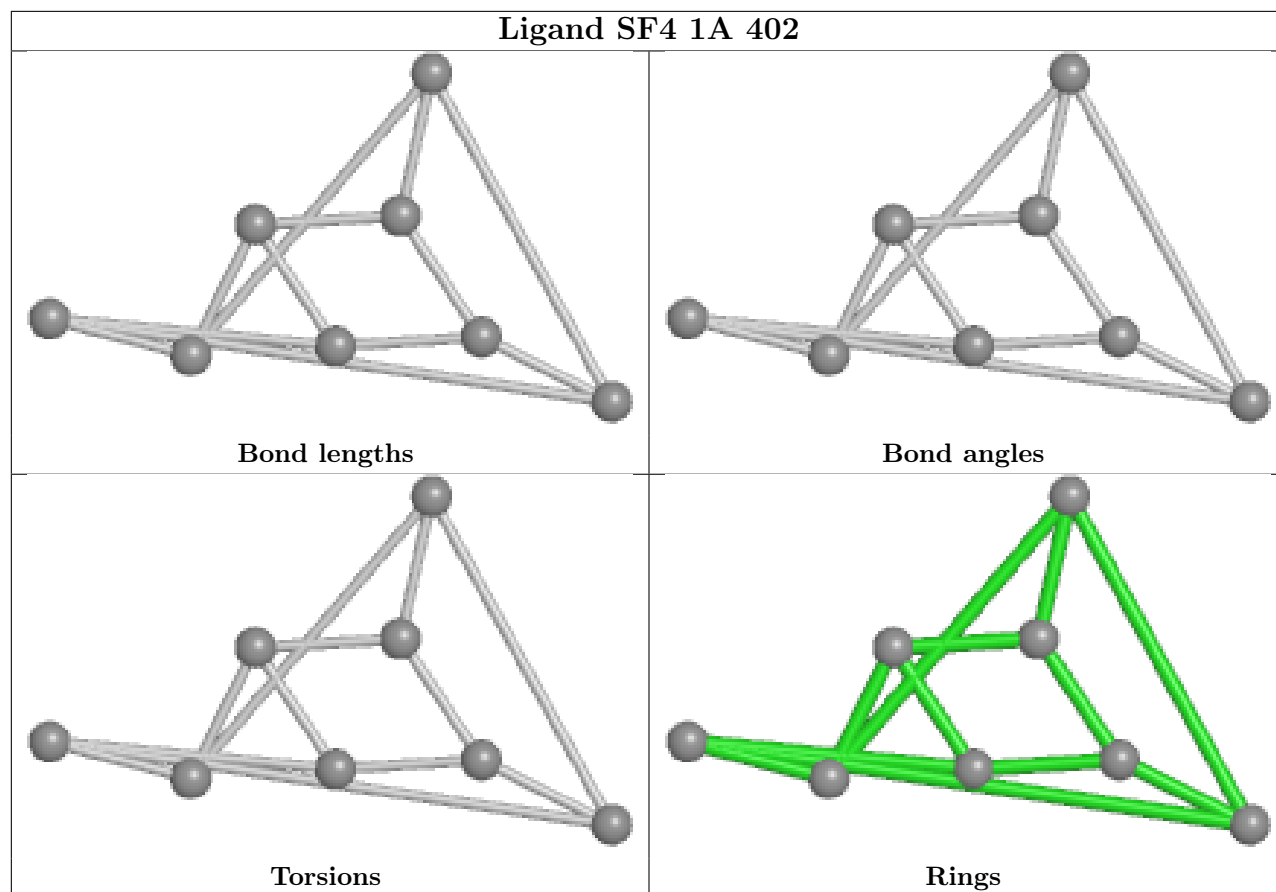


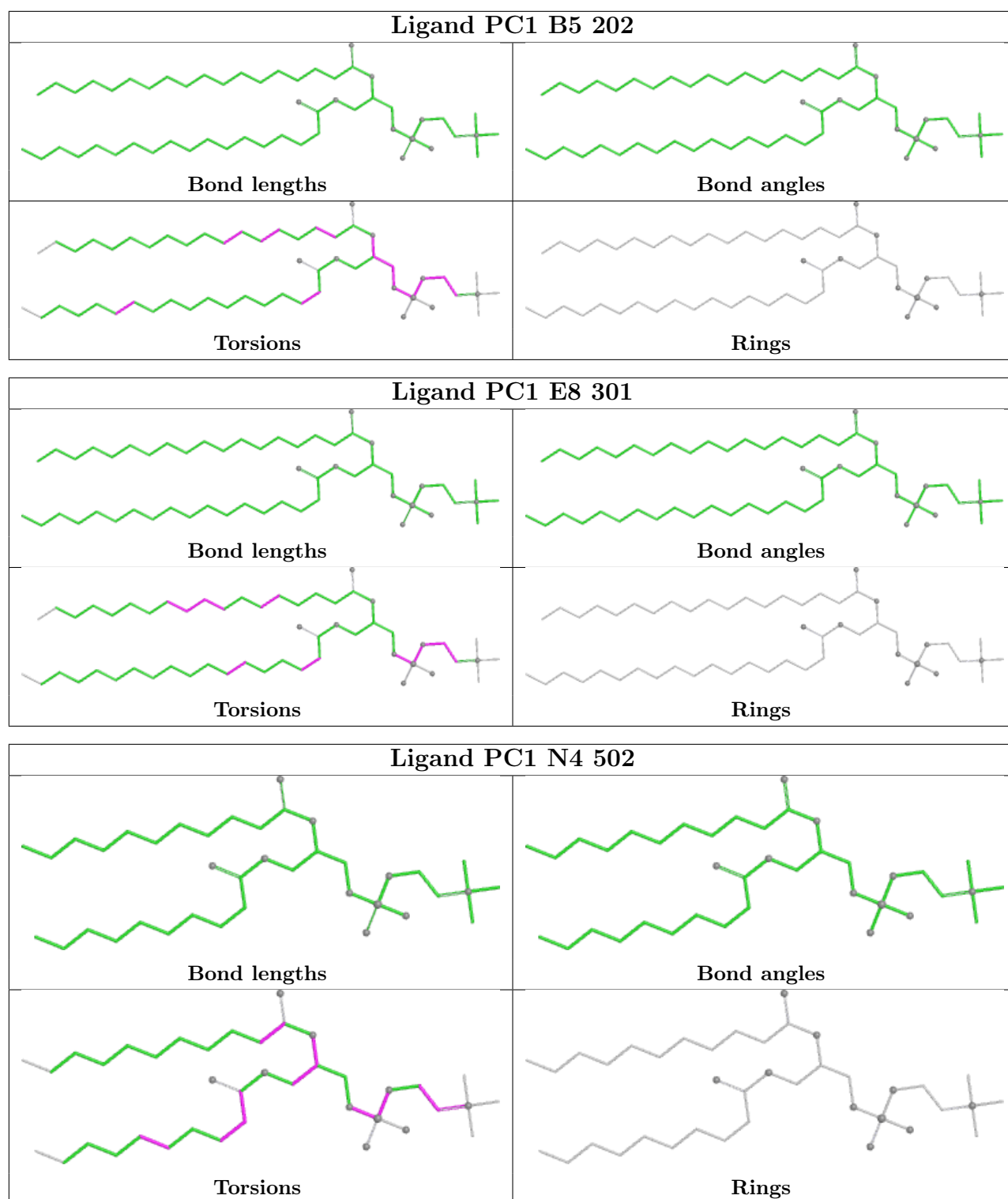


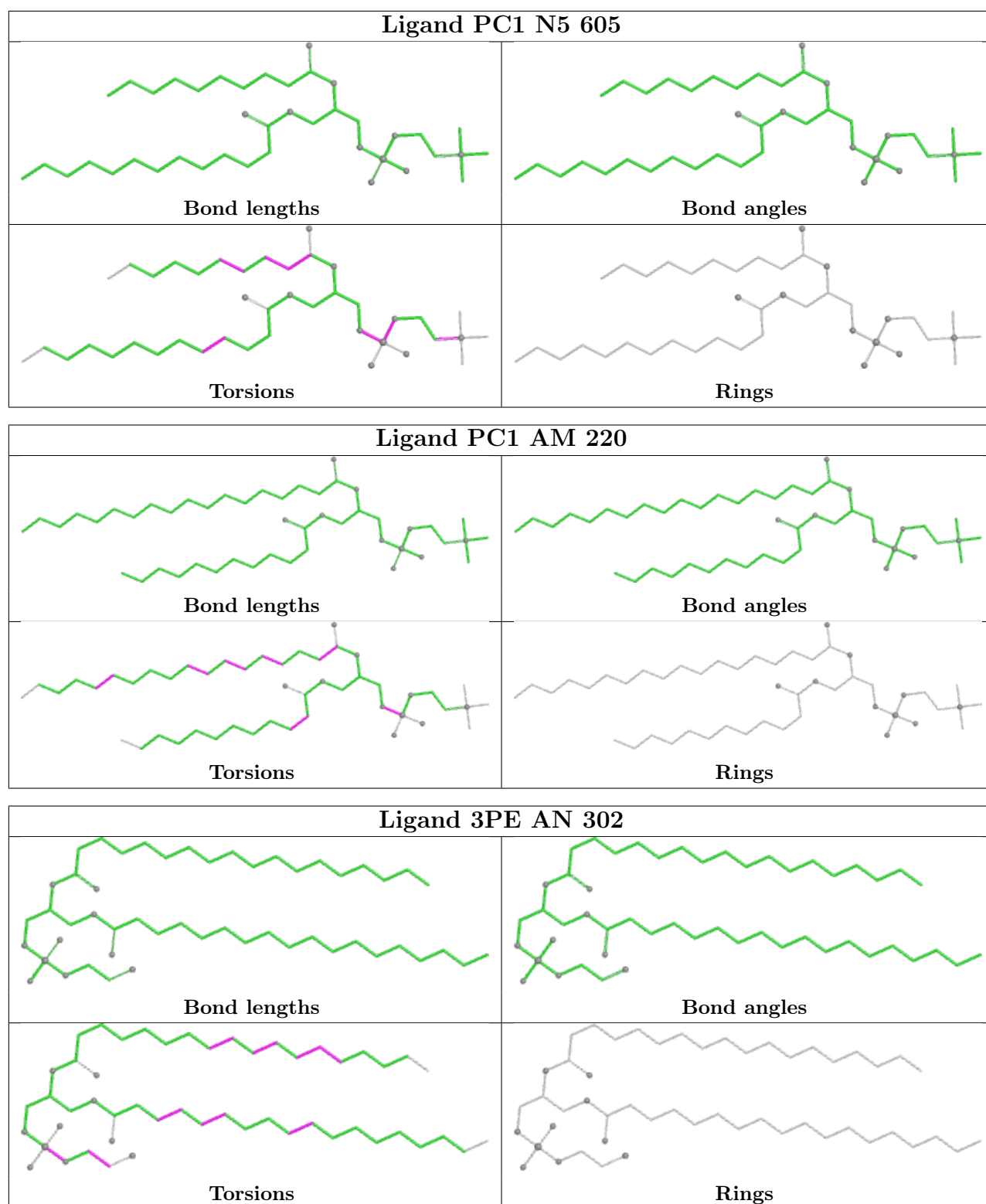


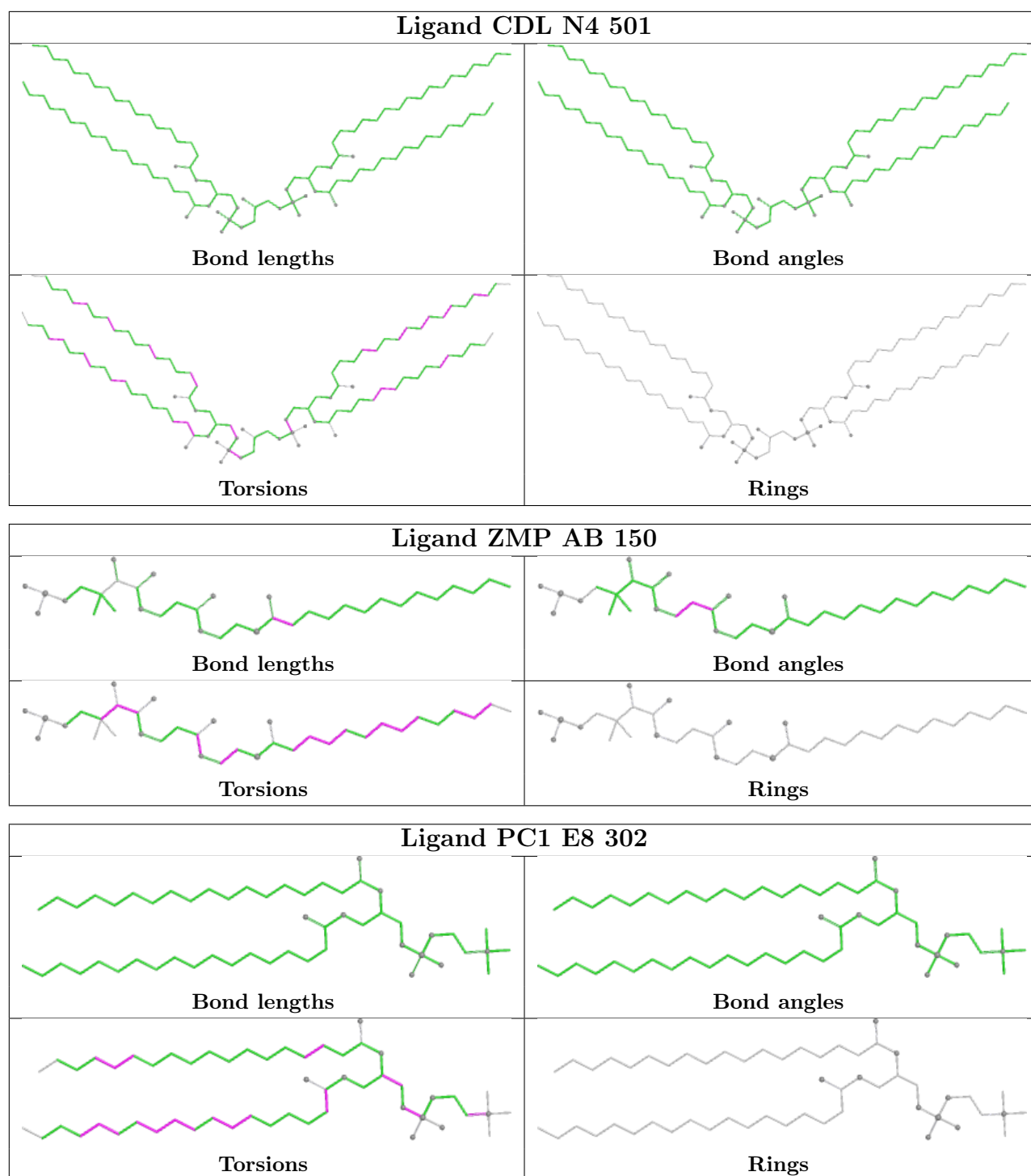


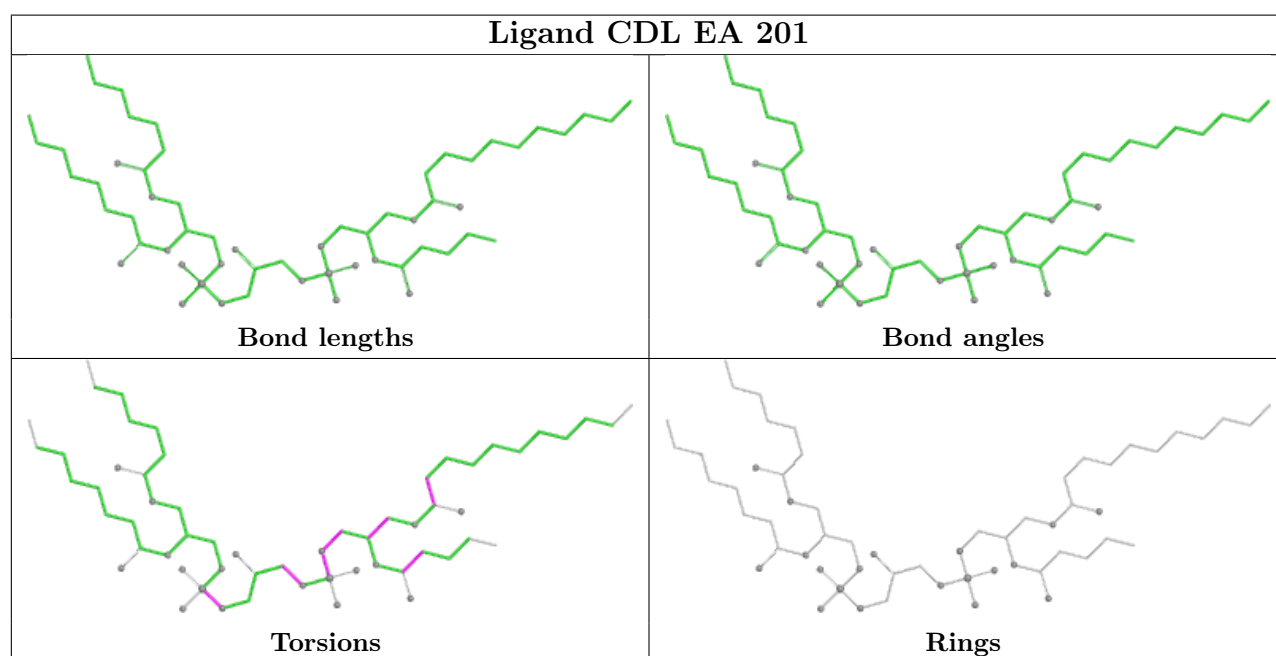
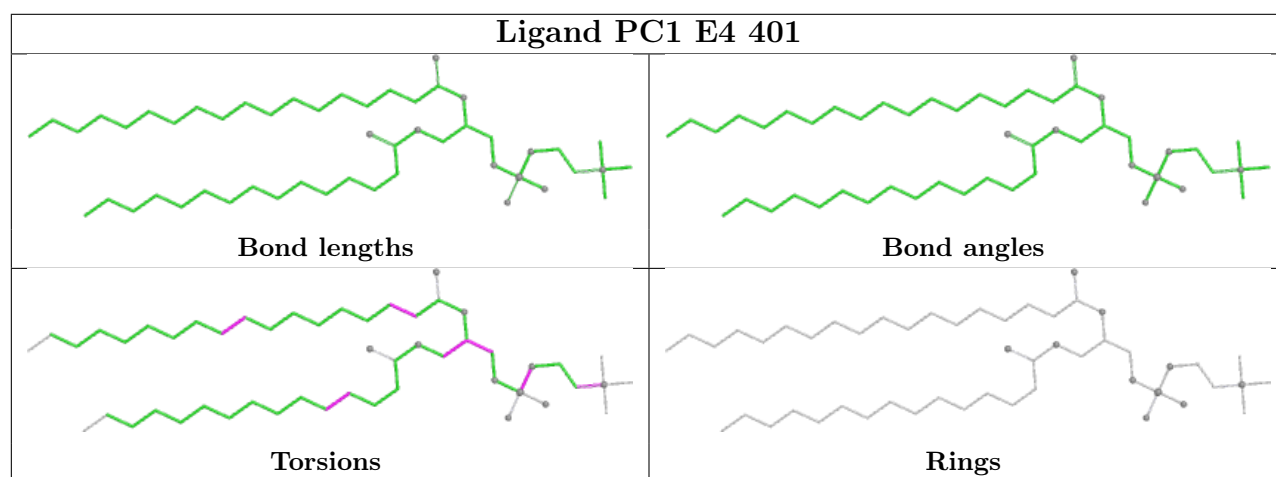
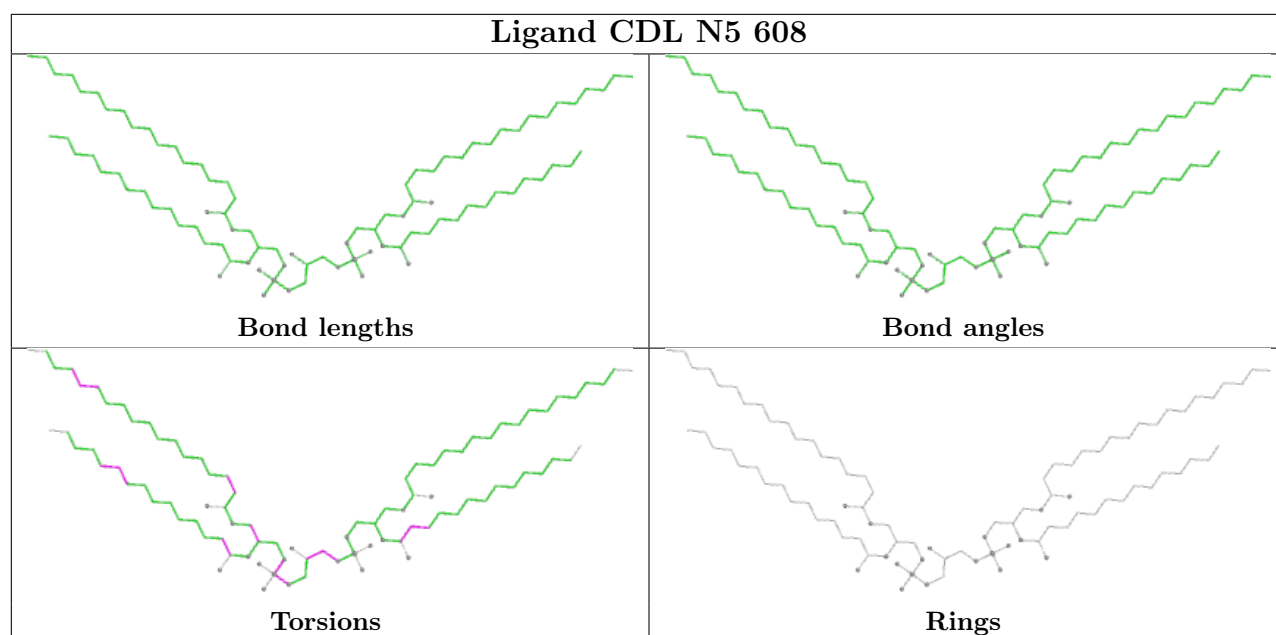


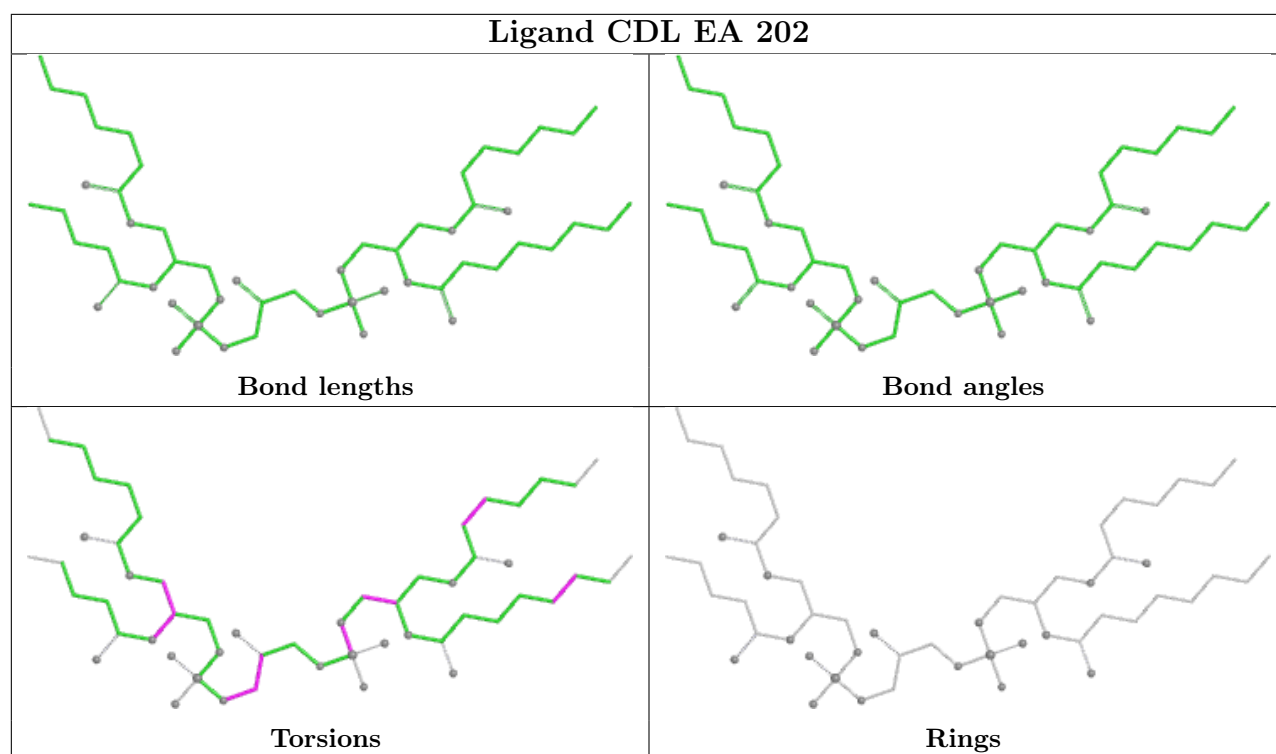












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

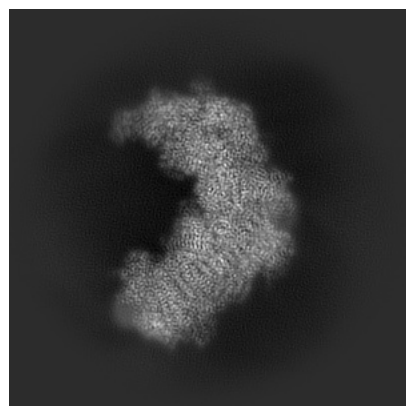
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-36107. These allow visual inspection of the internal detail of the map and identification of artifacts.

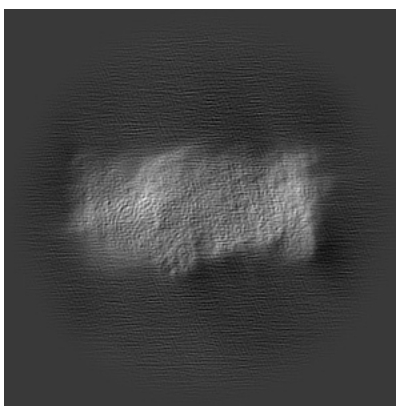
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

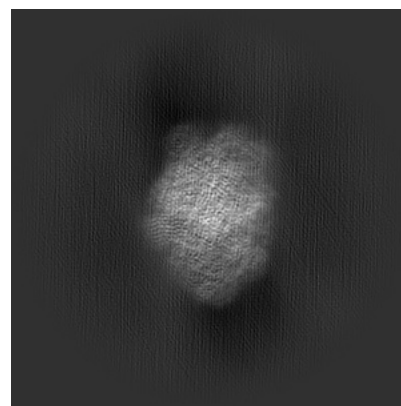
6.1.1 Primary map



X

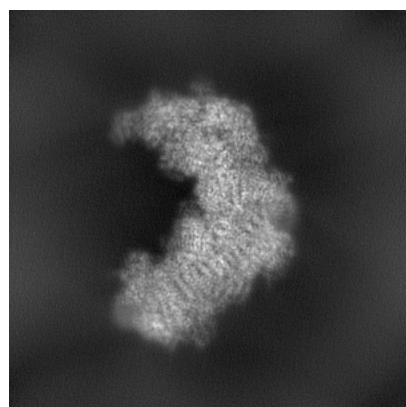


Y

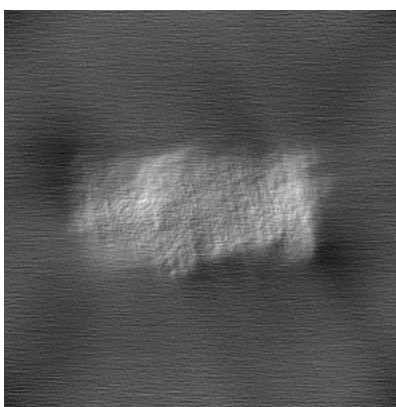


Z

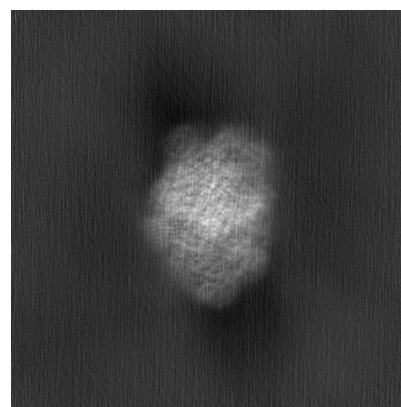
6.1.2 Raw map



X



Y

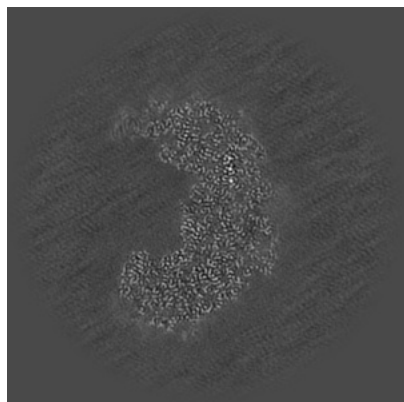


Z

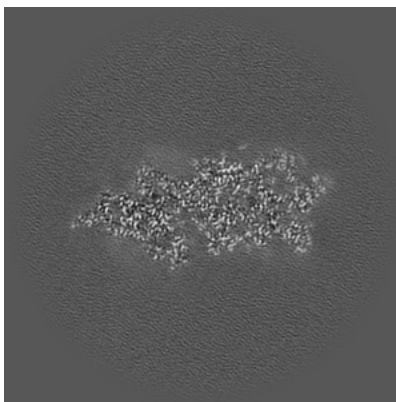
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

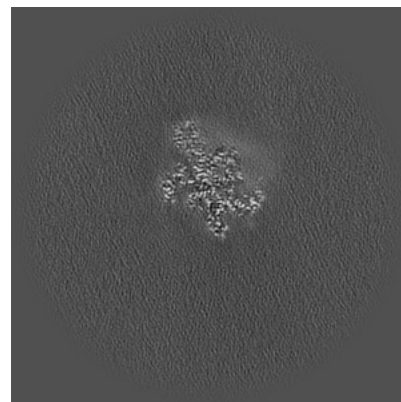
6.2.1 Primary map



X Index: 240



Y Index: 240

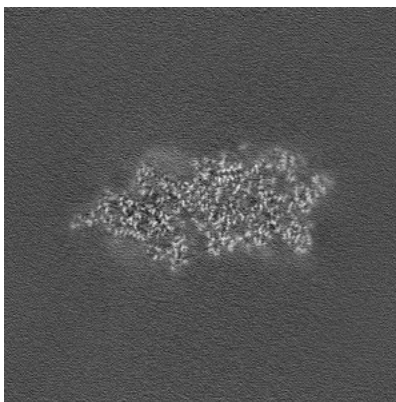


Z Index: 240

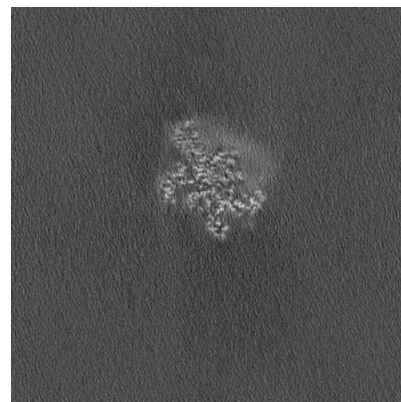
6.2.2 Raw map



X Index: 240



Y Index: 240



Z Index: 240

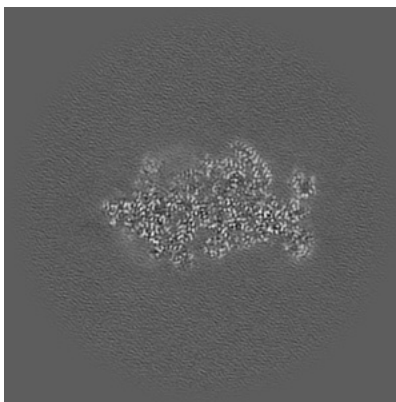
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

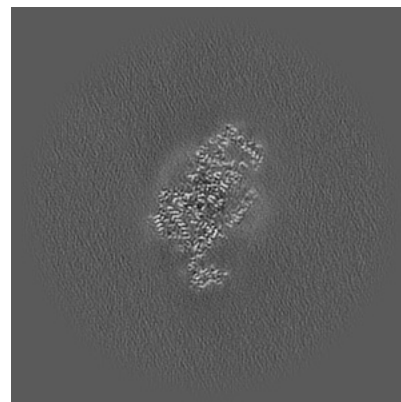
6.3.1 Primary map



X Index: 241



Y Index: 255

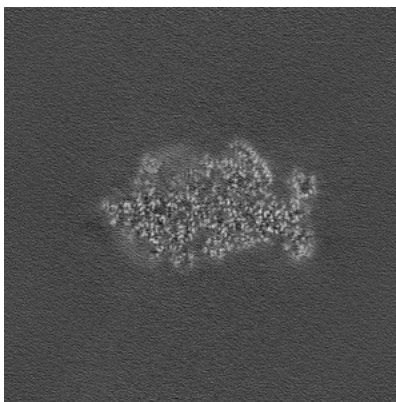


Z Index: 183

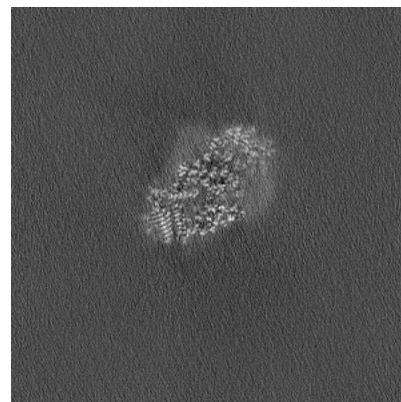
6.3.2 Raw map



X Index: 243



Y Index: 255

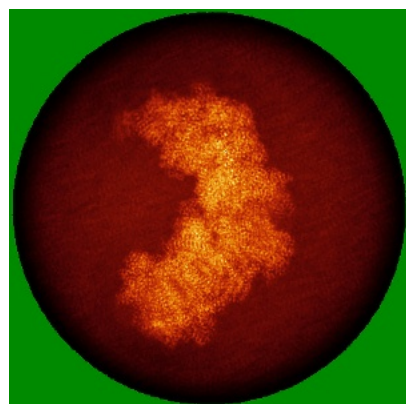


Z Index: 210

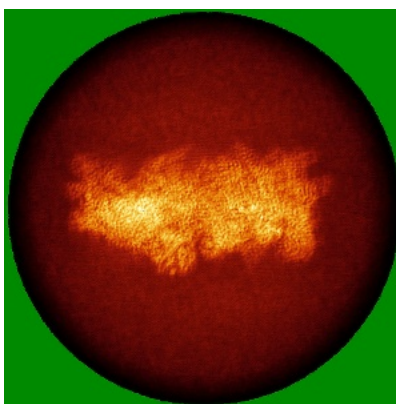
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

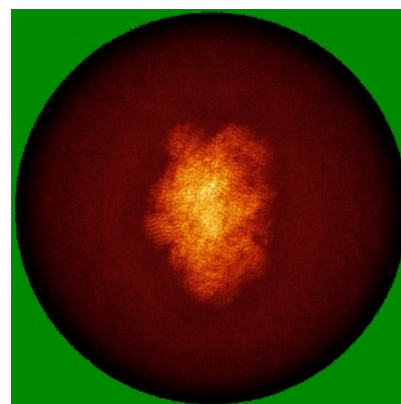
6.4.1 Primary map



X

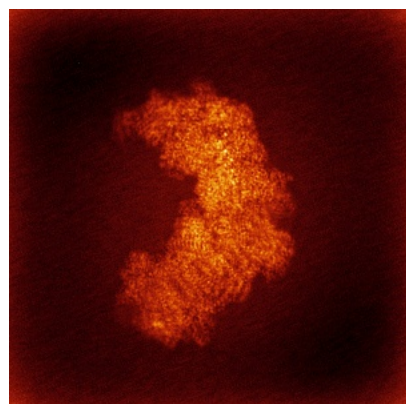


Y

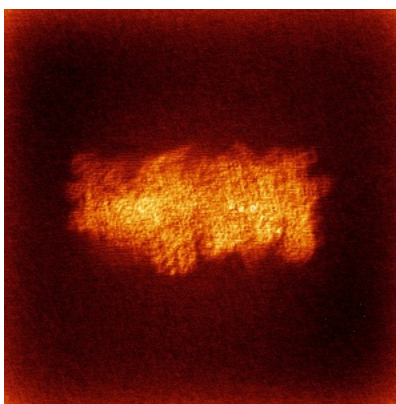


Z

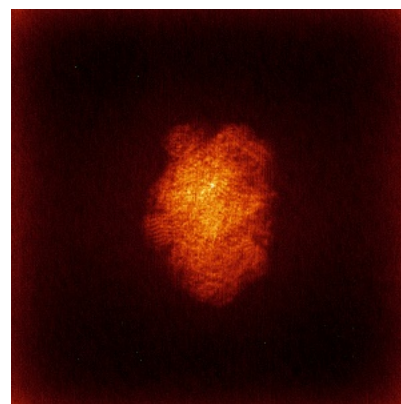
6.4.2 Raw map



X



Y

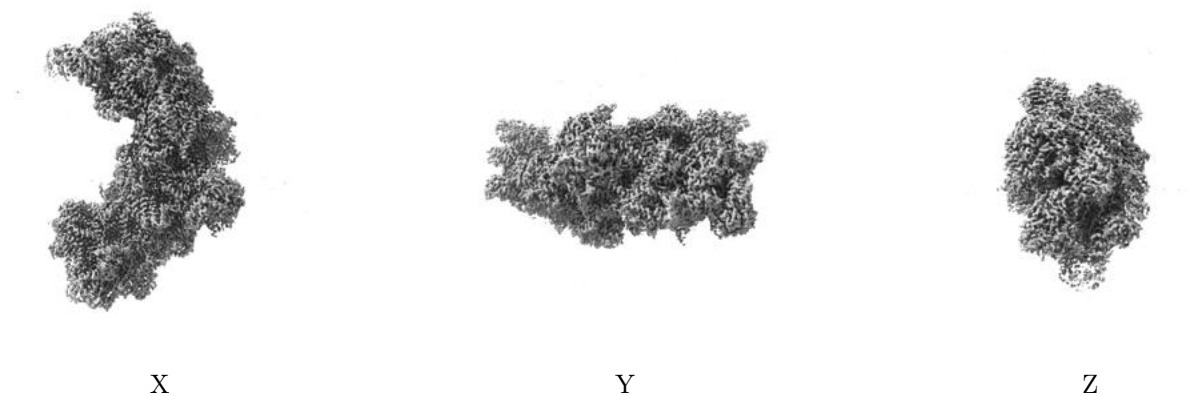


Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 5.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

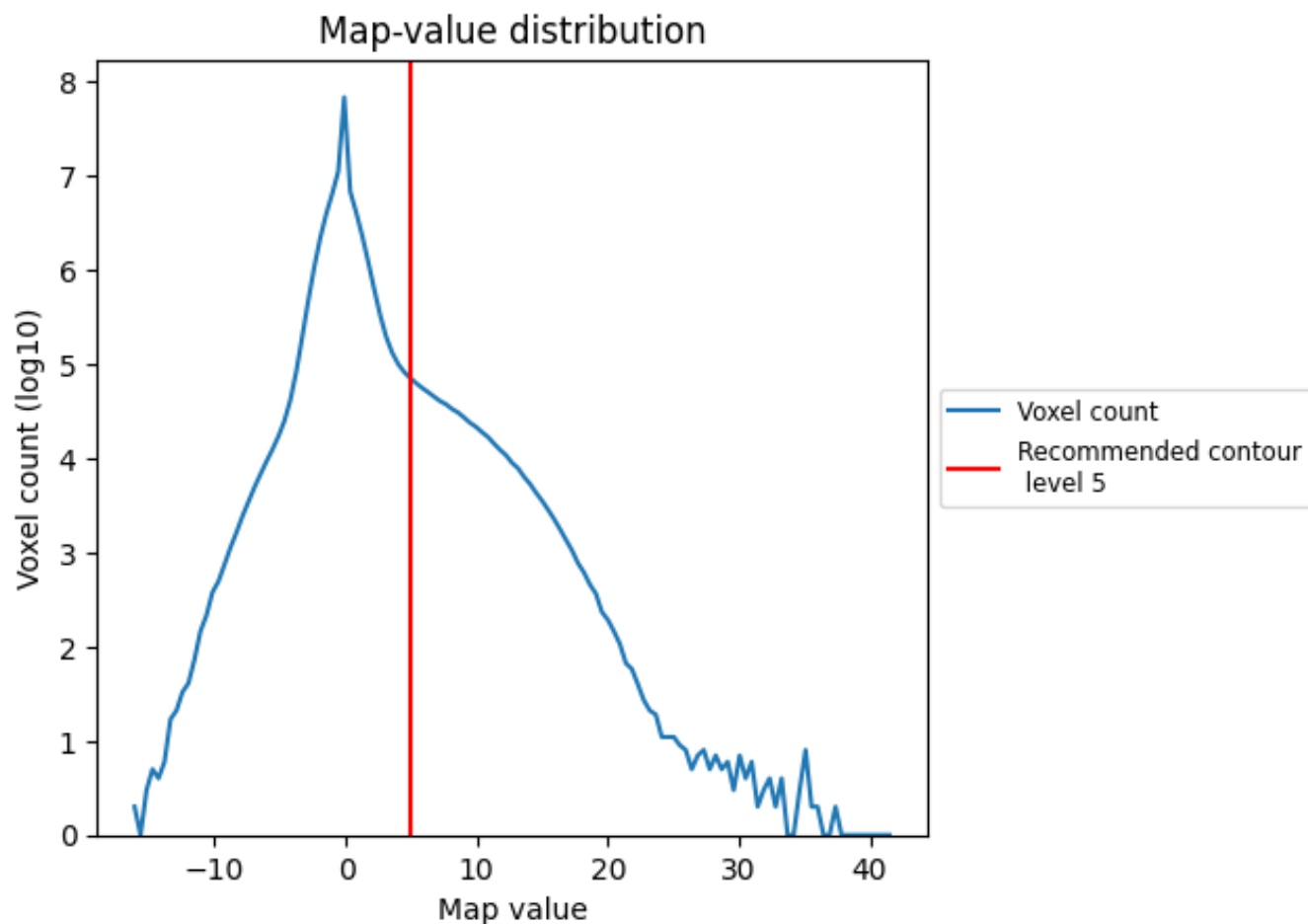
6.6 Mask visualisation [i](#)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

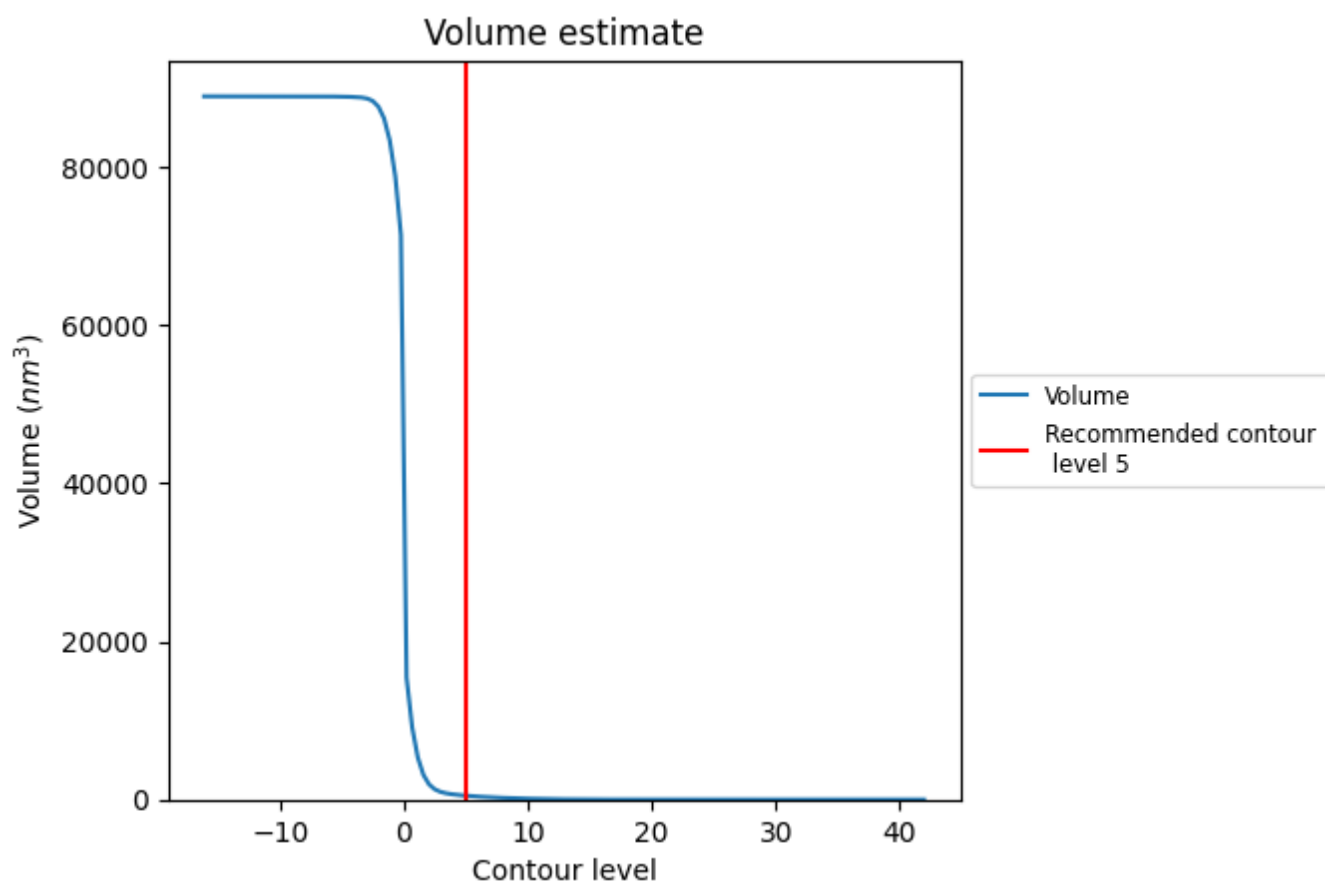
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

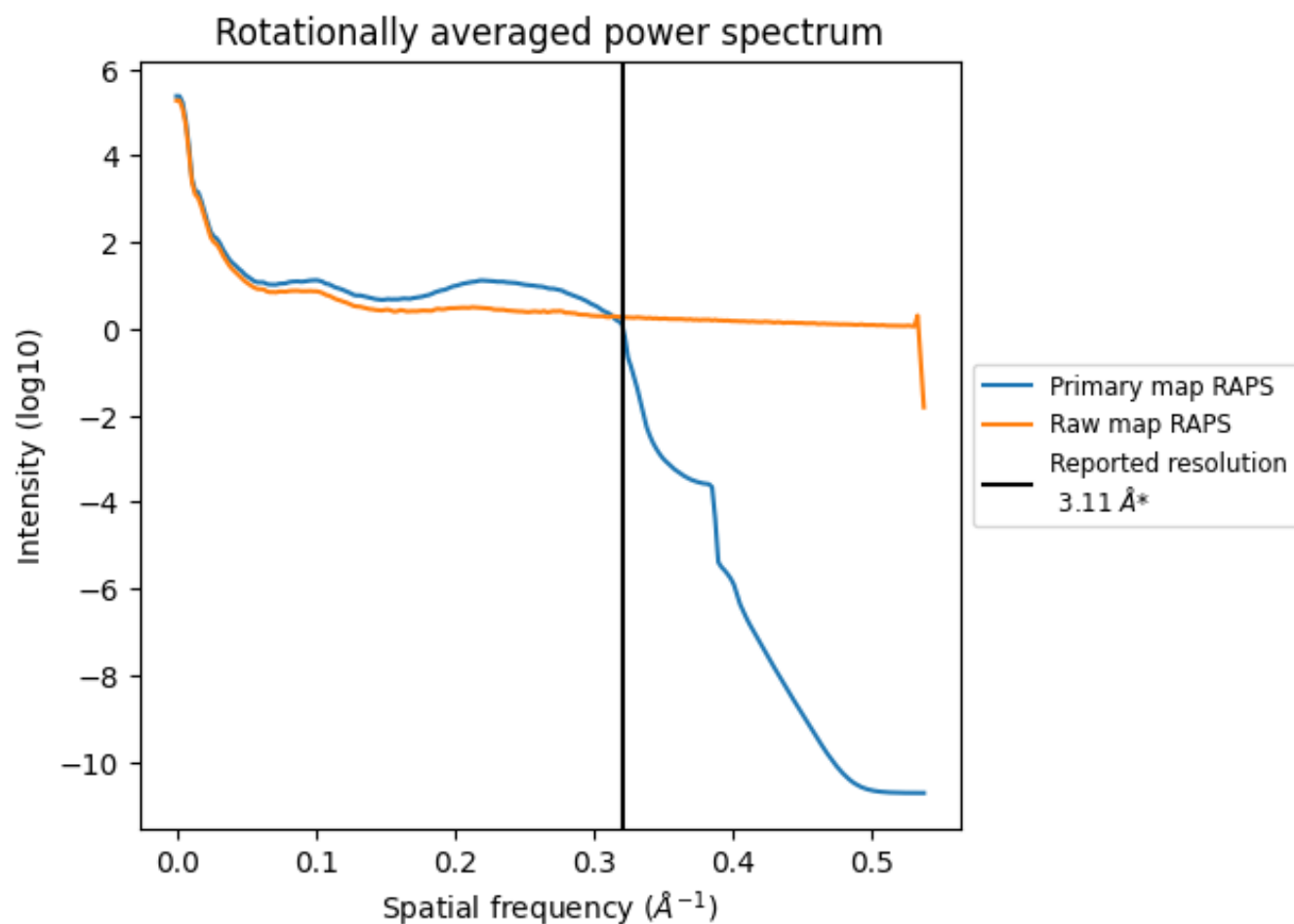
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 492 nm^3 ; this corresponds to an approximate mass of 445 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ

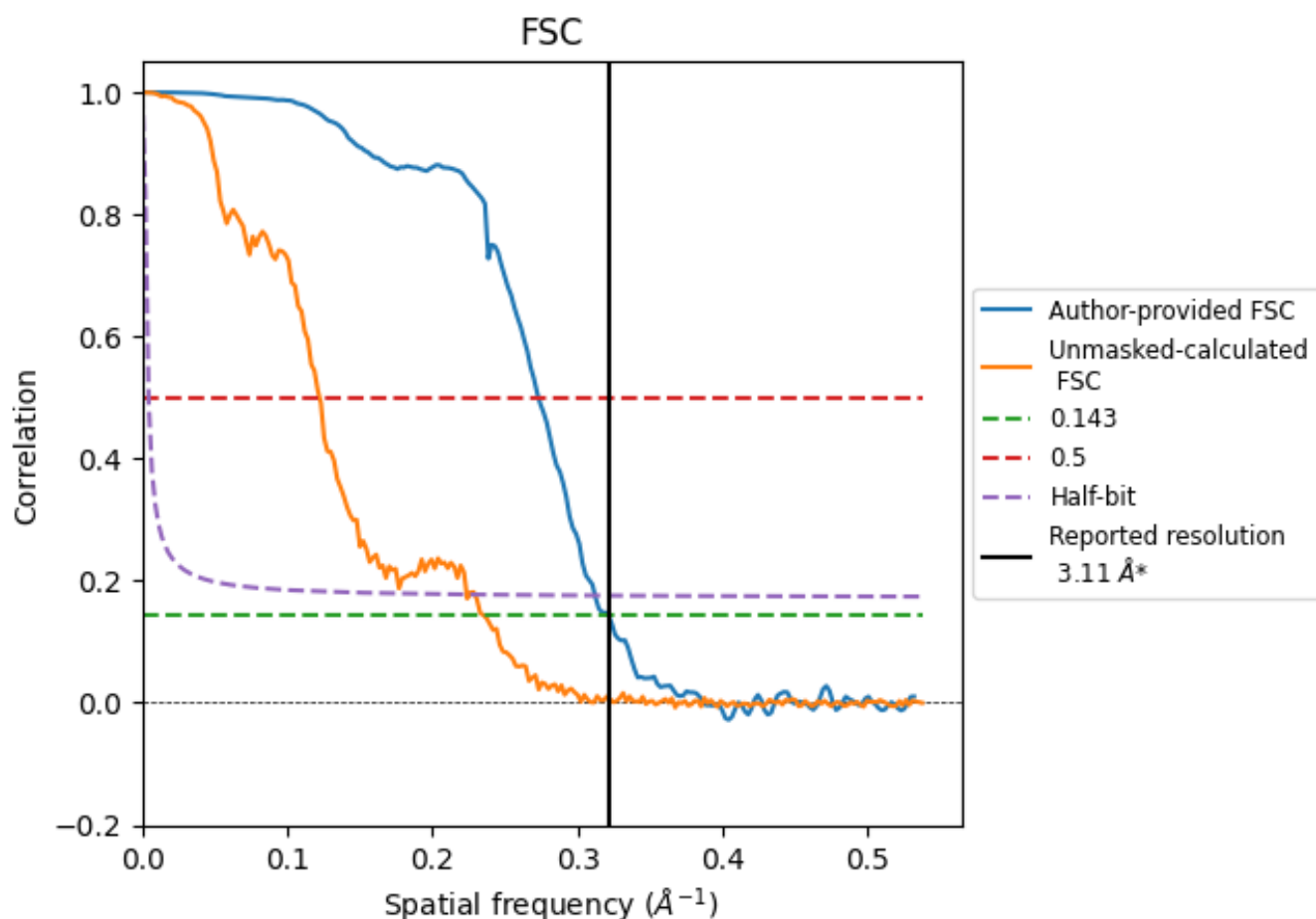


*Reported resolution corresponds to spatial frequency of 0.322 $\text{\AA}^{-1}$

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of $0.322 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

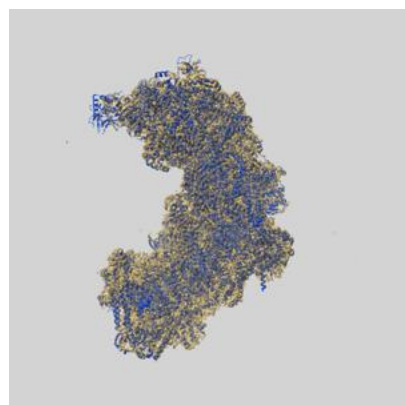
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.11	-	-
Author-provided FSC curve	3.11	3.66	3.20
Unmasked-calculated*	4.25	8.18	4.47

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.25 differs from the reported value 3.11 by more than 10 %

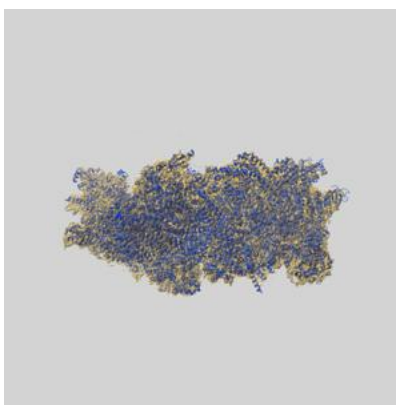
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-36107 and PDB model 8J9H. Per-residue inclusion information can be found in section [3](#) on page [23](#).

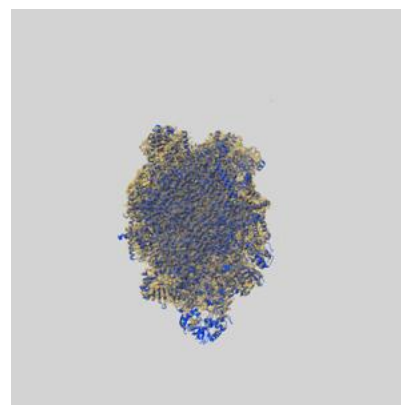
9.1 Map-model overlay [i](#)



X



Y



Z

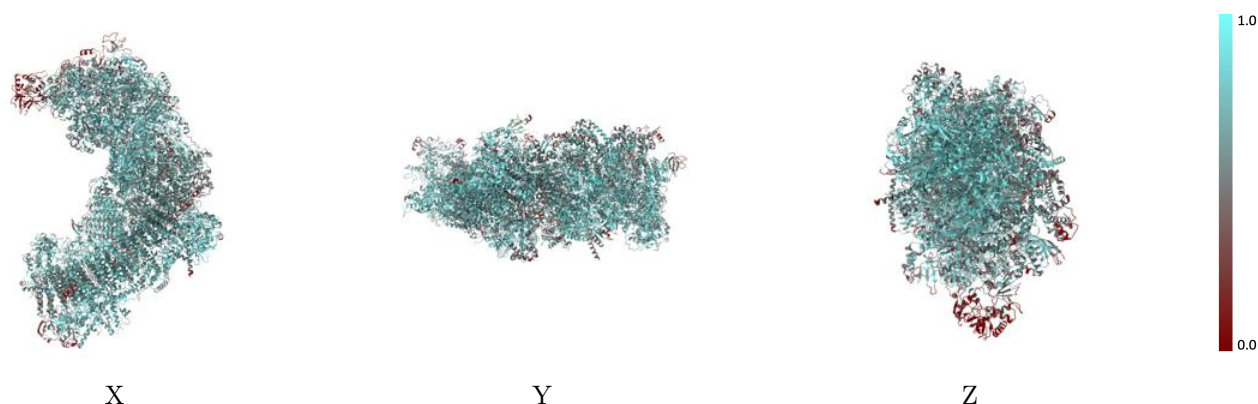
The images above show the 3D surface view of the map at the recommended contour level 5.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



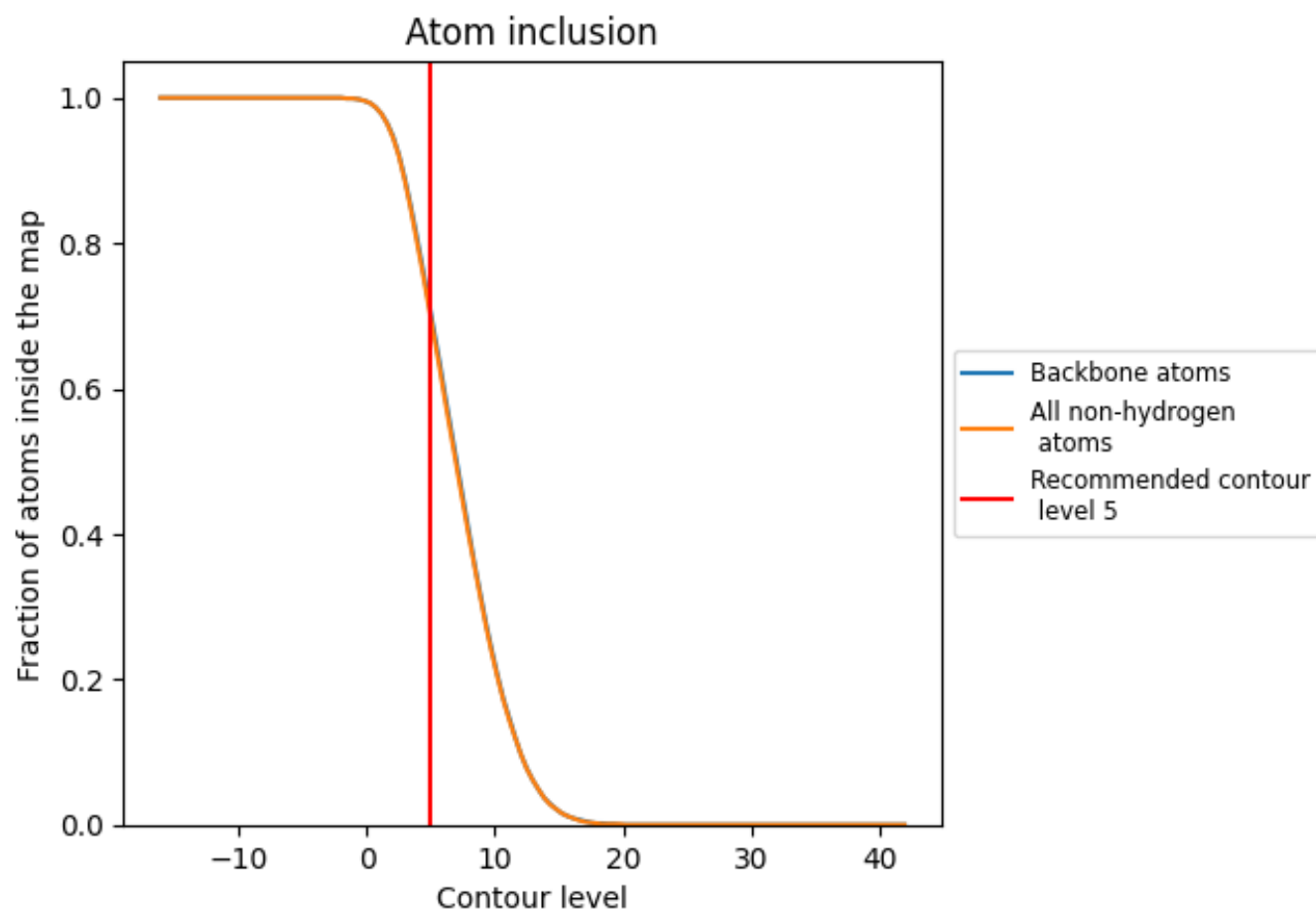
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (5).

9.4 Atom inclusion [i](#)



At the recommended contour level, 70% of all backbone atoms, 70% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (5) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6960	 0.5280
1A	 0.7610	 0.5470
1B	 0.7450	 0.5450
2B	 0.7580	 0.5480
4L	 0.7770	 0.5510
A1	 0.6470	 0.5160
A2	 0.6960	 0.5400
A3	 0.7050	 0.5380
A5	 0.7070	 0.5210
A6	 0.6010	 0.5050
A7	 0.7700	 0.5480
A8	 0.7400	 0.5270
A9	 0.7160	 0.5320
AB	 0.6660	 0.5280
AC	 0.7030	 0.5440
AL	 0.6980	 0.5340
AM	 0.6620	 0.5160
AN	 0.6990	 0.5280
B2	 0.7240	 0.5380
B3	 0.6950	 0.5310
B4	 0.7480	 0.5450
B5	 0.7350	 0.5270
B6	 0.7300	 0.5370
B7	 0.7620	 0.5290
B8	 0.7890	 0.5470
B9	 0.7620	 0.5520
BL	 0.7490	 0.5370
BM	 0.7280	 0.5400
C4	 0.6920	 0.5320
E1	 0.6710	 0.5130
E2	 0.6260	 0.5010
E3	 0.6040	 0.5050
E4	 0.6490	 0.5310
E5	 0.2050	 0.3270
E6	 0.6120	 0.5020



Continued on next page...

Continued from previous page...

Chain	Atom inclusion	Q-score
E7	 0.5100	 0.4620
E8	 0.6260	 0.5000
EA	 0.7720	 0.5550
EB	 0.7180	 0.5230
EC	 0.5870	 0.5080
ED	 0.6320	 0.5100
FX	 0.7710	 0.5550
G1	 0.7410	 0.5510
G2	 0.6850	 0.5280
G3	 0.7170	 0.5330
N1	 0.7090	 0.5360
N2	 0.7710	 0.5510
N3	 0.6980	 0.5420
N4	 0.7570	 0.5410
N5	 0.7310	 0.5320
N6	 0.6690	 0.5210
S2	 0.7950	 0.5550
S3	 0.7990	 0.5530
S4	 0.7350	 0.5510
S5	 0.7240	 0.5220
S6	 0.7780	 0.5520
S7	 0.7600	 0.5480
S8	 0.7980	 0.5540
U1	 0.5170	 0.4690
U2	 0.6170	 0.5230
V1	 0.7360	 0.5310
V2	 0.7490	 0.5310