



wwPDB EM Validation Summary Report ⓘ

Mar 7, 2026 – 03:44 AM UTC

PDB ID : 8J9I / pdb_00008j9i
EMDB ID : EMD-36108
Title : Cryo-EM structure of Euglena gracilis complex I, turnover 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 : 2.87 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

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:

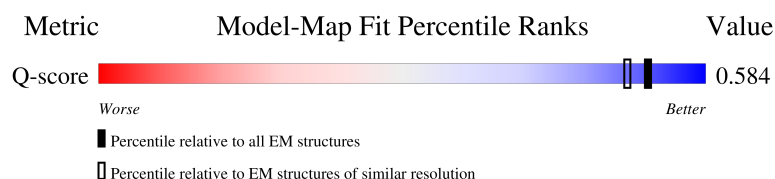
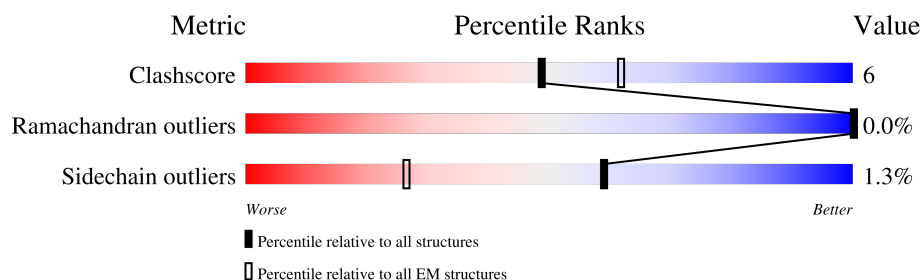
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 2.87 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.



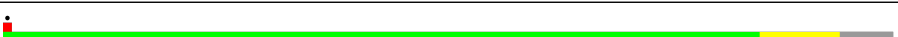
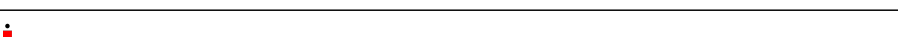
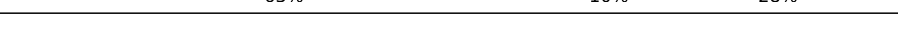
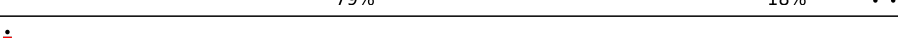
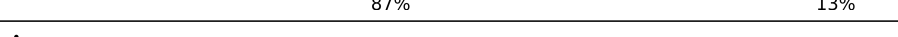
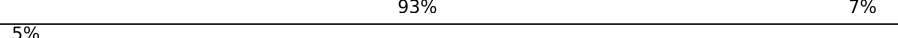
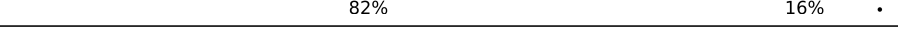
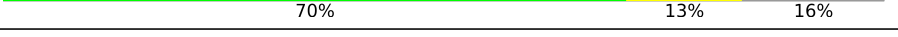
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	12062 (2.37 - 3.37)

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	AC	134	
15	AL	281	
16	AM	198	
17	AN	287	
18	B2	145	
19	B3	62	
20	B4	171	
21	B5	140	
22	B6	91	
23	B7	97	
24	B8	176	
25	B9	158	
26	BL	144	
27	BM	112	
28	C4	185	
29	E1	483	

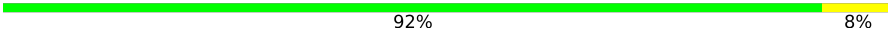
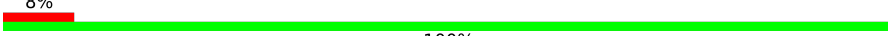
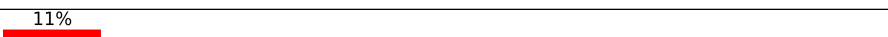
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Mol	Chain	Length	Quality of chain
30	E2	467	
31	E3	434	
32	E4	368	
33	E5	290	
34	E6	371	
35	E8	205	
36	EA	126	
37	EB	101	
38	EC	101	
39	ED	151	
40	FX	325	
41	G1	436	
42	G2	267	
43	G3	261	
44	N1	670	
45	N2	300	
46	N3	293	
46	N6	293	
47	N4	478	
48	N5	584	
49	S2	395	
50	S3	277	
51	S4	208	
52	S5	122	
53	S6	147	

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Mol	Chain	Length	Quality of chain
54	S7	207	 82% 14%
55	S8	212	 69% 16% 14%
56	U1	12	 92% 8%
56	U2	12	 8% 100%
57	V1	526	 77% 19%
58	V2	225	 84% 16%
59	E7	246	 11% 89% 11%

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	-

2 Entry composition

There are 71 unique types of molecules in this entry. The entry contains 226854 atoms, of which 112888 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 NDUFS1A.

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

- Molecule 2 is a protein called NDUFS1B.

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

- Molecule 3 is a protein called ND2B.

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

- Molecule 4 is a protein called ND4L.

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

- Molecule 5 is a protein called NDUFA1.

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

- Molecule 6 is a protein called NDUFA2.

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

- 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 NDUFAB1-beta.

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

- Molecule 15 is a protein called NDUFA12.

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

- Molecule 16 is a protein called NDUFA13.

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

- Molecule 17 is a protein called NDUFA11.

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

- Molecule 18 is a protein called NDUFB2.

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

- Molecule 19 is a protein called NDUFB3.

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

- Molecule 20 is a protein called NDUFB4.

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

- Molecule 21 is a protein called NDUFB5.

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

- Molecule 22 is a protein called NDUFB6.

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

- Molecule 23 is a protein called NDUFB7.

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

- Molecule 24 is a protein called NDUFB8.

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

- Molecule 25 is a protein called NDUFB9.

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

- Molecule 26 is a protein called NDUFB10.

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

- Molecule 27 is a protein called NDUFB11.

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

- Molecule 28 is a protein called NDUFC2.

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

- Molecule 29 is a protein called NDUEG1.

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

- Molecule 30 is a protein called NDUEG2.

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

- Molecule 31 is a protein called NDUEG3.

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

- Molecule 32 is a protein called NDUEG4.

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

- Molecule 33 is a protein called NDUEG5.

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

- Molecule 34 is a protein called NDUEG6.

Mol	Chain	Residues	Atoms						AltConf	Trace
34	E6	342	Total	C	H	N	O	S	0	0
			5629	1839	2758	507	513	12		

- Molecule 35 is a protein called NDUEG8.

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

- Molecule 36 is a protein called NDUEG10.

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

- Molecule 37 is a protein called NDUEG11.

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

- Molecule 38 is a protein called NDUEG12.

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

- Molecule 39 is a protein called NDUEG13.

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

- Molecule 40 is a protein called NDUFEX.

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

- Molecule 41 is a protein called NDUCA1.

Mol	Chain	Residues	Atoms						AltConf	Trace
41	G1	418	Total	C	H	N	O	S	0	0
			6420	2072	3139	581	612	16		

- Molecule 42 is a protein called NDUCA2.

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

- Molecule 43 is a protein called NDUCA3.

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

- Molecule 44 is a protein called ND1.

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

- Molecule 45 is a protein called ND2A.

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

- Molecule 46 is a protein called ND3.

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

- Molecule 47 is a protein called NADH-ubiquinone oxidoreductase chain 4.

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

- Molecule 48 is a protein called ND5.

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

- Molecule 49 is a protein called NDUFS2.

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

- Molecule 50 is a protein called NDUFS3.

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

- Molecule 51 is a protein called NDUFS4.

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

- Molecule 52 is a protein called NDUFS5.

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

- Molecule 53 is a protein called NDUFS6.

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

- Molecule 54 is a protein called NDUFS7.

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

- Molecule 55 is a protein called NDUFS8.

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

- Molecule 56 is a protein called UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK.

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

- Molecule 57 is a protein called NDUFV1.

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

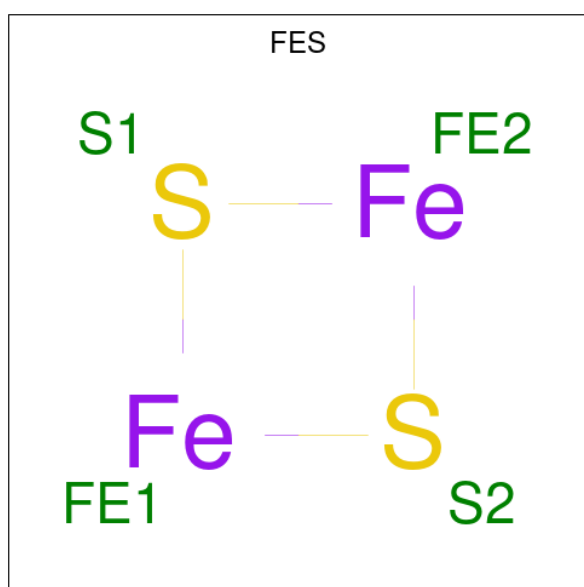
- Molecule 58 is a protein called NDUFV2.

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

- Molecule 59 is a protein called NDUEG7.

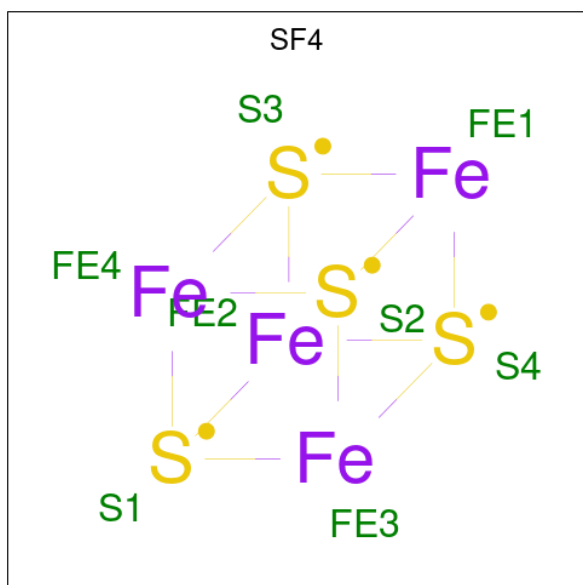
Mol	Chain	Residues	Atoms						AltConf	Trace
59	E7	246	Total	C	H	N	O	S	0	0
			3780	1205	1892	332	344	7		

- Molecule 60 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe₂S₂).



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).

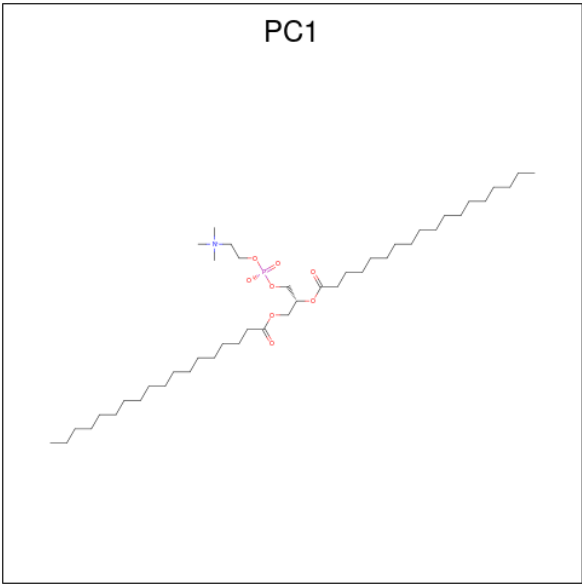


Mol	Chain	Residues	Atoms			AltConf
61	1A	1	Total	Fe	S	0
			8	4	4	
61	1A	1	Total	Fe	S	0
			8	4	4	
61	S7	1	Total	Fe	S	0
			8	4	4	
61	S8	1	Total	Fe	S	0
			8	4	4	
61	S8	1	Total	Fe	S	0
			8	4	4	
61	V1	1	Total	Fe	S	0
			8	4	4	

- 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: C₄₄H₈₈NO₈P).



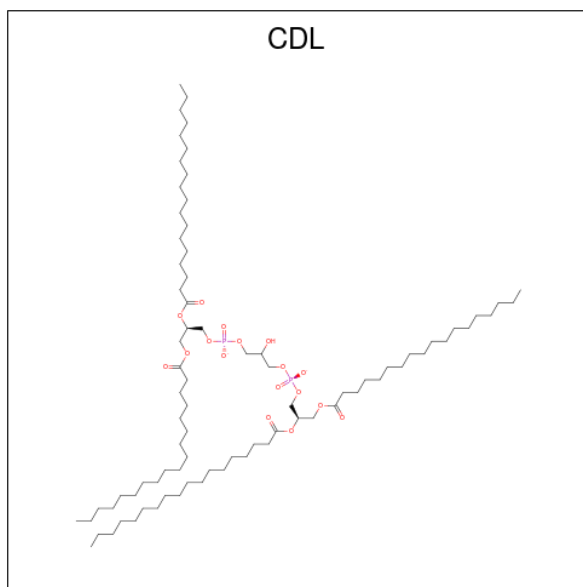
Mol	Chain	Residues	Atoms						AltConf
63	A1	1	Total	C	H	N	O	P	0
			124	39	75	1	8	1	
63	A1	1	Total	C	H	N	O	P	0
			67	21	36	1	8	1	
63	A9	1	Total	C	H	N	O	P	0
			73	23	40	1	8	1	
63	A9	1	Total	C	H	N	O	P	0
			73	23	40	1	8	1	
63	AL	1	Total	C	H	N	O	P	0
			127	40	77	1	8	1	
63	AM	1	Total	C	H	N	O	P	0
			124	39	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 142	C 44	H 88	N 1	O 8	P 1	0
63	E8	1	Total 142	C 44	H 88	N 1	O 8	P 1	0
63	E8	1	Total 73	C 23	H 40	N 1	O 8	P 1	0
63	E8	1	Total 64	C 20	H 34	N 1	O 8	P 1	0
63	ED	1	Total 142	C 44	H 88	N 1	O 8	P 1	0
63	N1	1	Total 124	C 39	H 75	N 1	O 8	P 1	0
63	N1	1	Total 94	C 30	H 54	N 1	O 8	P 1	0
63	N2	1	Total 85	C 27	H 48	N 1	O 8	P 1	0
63	N3	1	Total 103	C 32	H 61	N 1	O 8	P 1	0
63	N4	1	Total 91	C 29	H 52	N 1	O 8	P 1	0
63	N4	1	Total 73	C 23	H 40	N 1	O 8	P 1	0
63	N5	1	Total 142	C 44	H 88	N 1	O 8	P 1	0
63	N5	1	Total 97	C 31	H 56	N 1	O 8	P 1	0
63	N5	1	Total 82	C 26	H 46	N 1	O 8	P 1	0

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



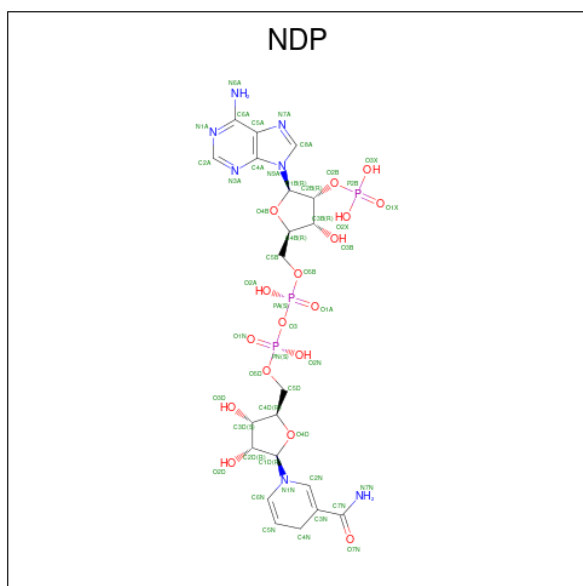
Mol	Chain	Residues	Atoms					AltConf
64	A3	1	Total	C	H	O	P	0
			118	39	60	17	2	
64	AL	1	Total	C	H	O	P	0
			148	49	80	17	2	
64	AL	1	Total	C	H	O	P	0
			136	45	72	17	2	
64	AL	1	Total	C	H	O	P	0
			154	51	84	17	2	
64	AM	1	Total	C	H	O	P	0
			163	53	91	17	2	
64	AM	1	Total	C	H	O	P	0
			163	53	91	17	2	
64	AM	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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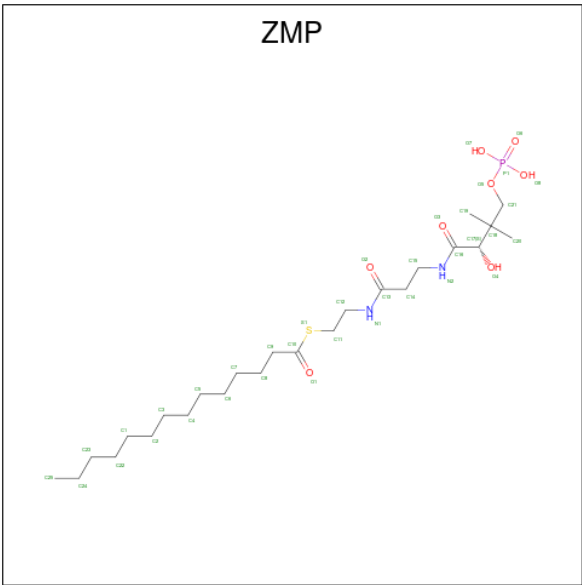
Mol	Chain	Residues	Atoms					AltConf
64	N4	1	Total	C	H	O	P	0
			247	79	149	17	2	
64	N5	1	Total	C	H	O	P	0
			157	51	87	17	2	
64	N5	1	Total	C	H	O	P	0
			229	74	136	17	2	
64	E7	1	Total	C	H	O	P	0
			148	49	80	17	2	

- Molecule 65 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$).



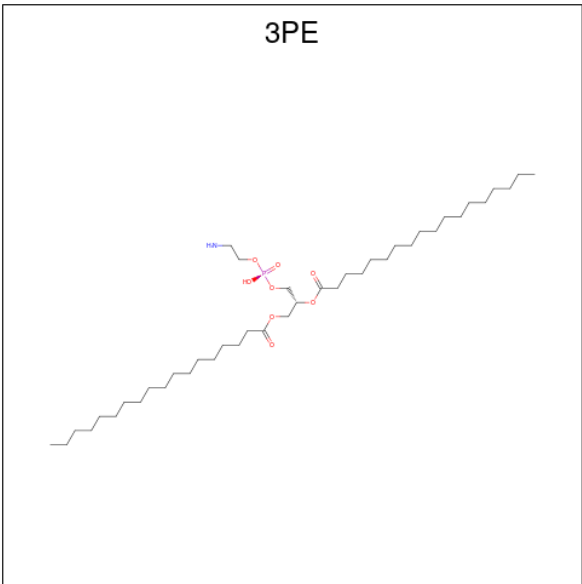
Mol	Chain	Residues	Atoms						AltConf
65	A9	1	Total	C	H	N	O	P	0
			74	21	26	7	17	3	

- Molecule 66 is S-[2-(N-[(2S)-2-hydroxy-3,3-dimethyl-4-(phosphonooxy)butanoyl]-beta-alanyl}amino)ethyl] tetradecanethioate (CCD ID: ZMP) (formula: $C_{25}H_{49}N_2O_8PS$) (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).



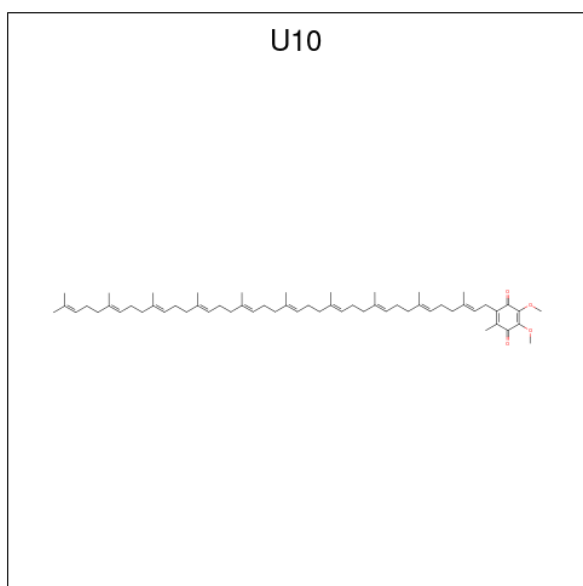
Mol	Chain	Residues	Atoms						AltConf
67	AN	1	Total	C	H	N	O	P	0
			132	41	81	1	8	1	

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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).

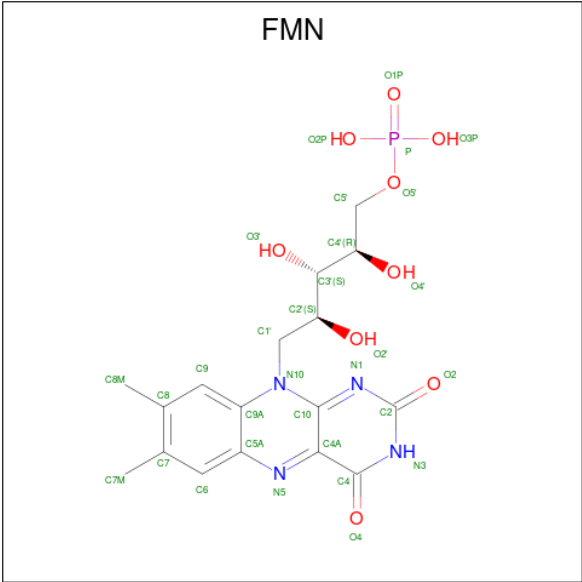


Mol	Chain	Residues	Atoms				AltConf
68	N4	1	Total	C	H	O	
			98	39	55	4	0

- 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

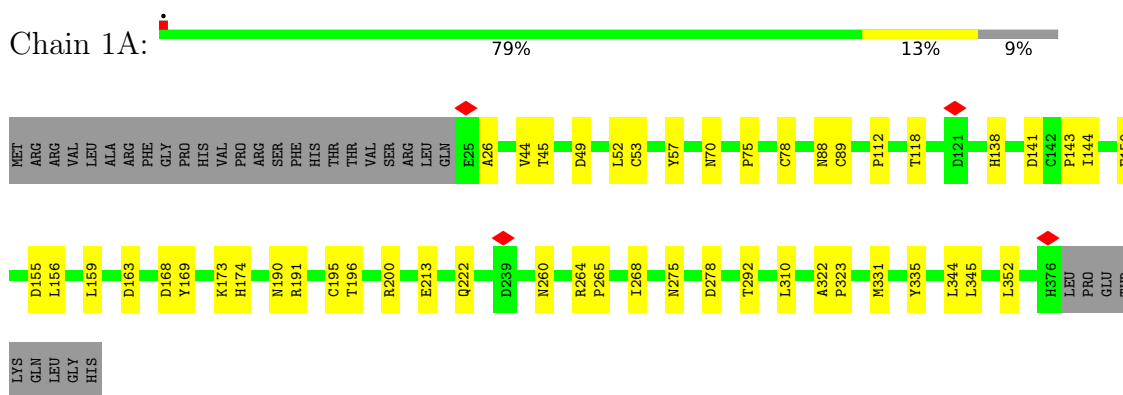
- Molecule 70 is FLAVIN MONONUCLEOTIDE (CCD ID: FMN) (formula: $C_{17}H_{21}N_4O_9P$).



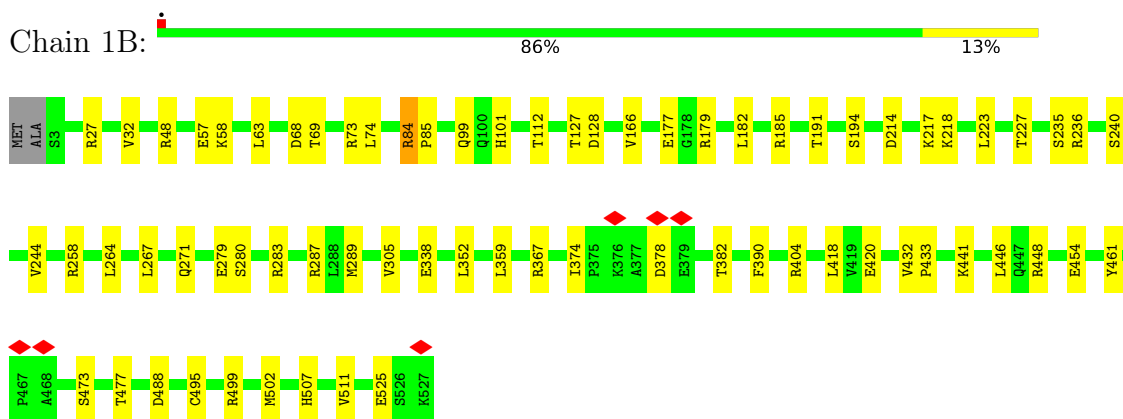
3 Residue-property plots

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.

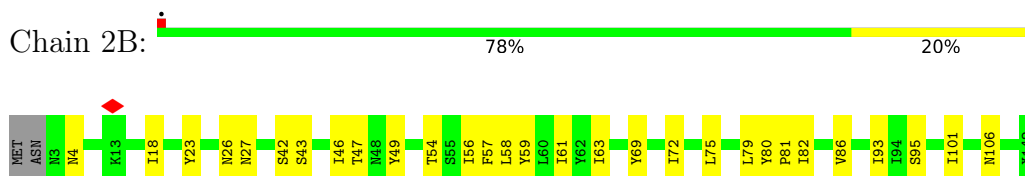
• Molecule 1: NDUFS1A



• Molecule 2: NDUFS1B



• Molecule 3: ND2B



• Molecule 4: ND4L

Chain 4L:  51% 12% 37%

MET LEU LYS ALA ILE GLN PRO MET SER SER LEU ARG SER ARG VAL VAL SER SER ARG ALA LEU SER SER GLY TYR LYS GLN PHE GLY LEU THR PHE LYS GLU VAL LYS GLN ASN PRO THR THR PHE MET GLN ARG PRO LEU ALA ILE ALA ASN GLY LEU PHE

V61 I78 G79 P80 Y81 I84 R91 K95 R103 V107 E108 T111 L114 F118 F119 I120 P126 L127 M135 L141 K142 T143 L144 M145 V163 K168 ARG LEU LYS

• Molecule 5: NDUFA1

Chain A1:  89% 8%

MET P2 C31 H37 K38 D43 Y49 A65 V69 E96 N124 V127 F137 P138 GLY TYR TYR

• Molecule 6: NDUFA2

Chain A2:  91% 8%

MET A2 R8 R14 A23 P26 E27 A28 A29 A30 I31 N32 R58 L64 D118 L147 D157 D163 T164 I165 M193

• Molecule 7: NDUFA3

Chain A3:  90% 10%

H1 R10 H17 N22 R23 E25 Q28 R31 W32 E33 E49 A55 G99 R116 A121 A122 V123 Q124 GLY

• Molecule 8: NDUFA5

Chain A5:  5% 76% 8% 16%

MET LEU ARG ALA ALA GLU MET LEU ARG CYS ALA VAL ALA VAL ILE GLY ARG ARG GLY TYR A21 K38 V39 R40 T50 D60 Y63 R64 R65 S66 V67 A81 E82 L107 E123 R124 V125 L126 D127 Y146 E155 K161 E169 K170 A171 R172 V173 A174 ALA

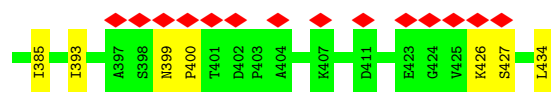
LYS ALA ALA GLU MET LEU ARG CYS ALA VAL ALA VAL ILE GLY ARG ARG GLY TYR

• Molecule 9: NDUFA6

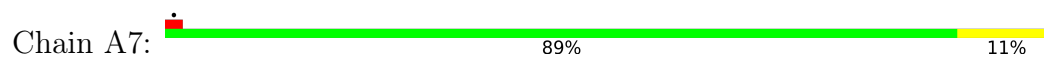
Chain A6:  7% 85% 11%

MET PRO GLN MET MET VAL ALA THR LYS LYS VAL VAL ARG ARG PRO PHE S12 V13 R14 D24 V31 V32 R33 R36 E39 F40 E41 H45 P48 R59 Y62 L66 K67 R86 R92 S92 L99 L116 E117 K132 M135 M136 D137 R158 G163

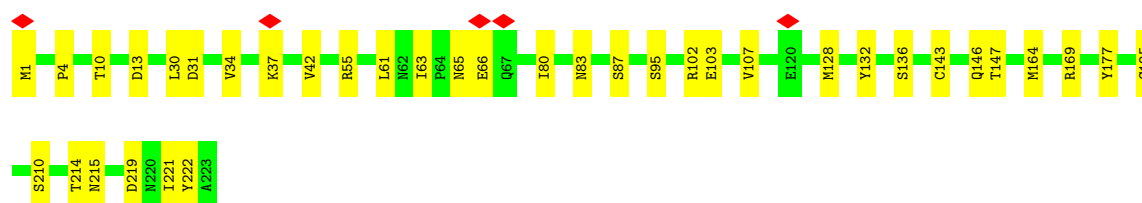
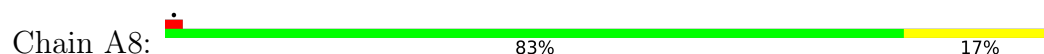
E169 E175 D176 V177 T178 A189 L194 G199 D200 R204 H205 L206 L207 V210 Q215 L221 E280 L294 D297 L303 L304 D309 R312 A316 G317 L318 L319 S320 T321 T322 E323 P324 D325 Y339 L344 L360 V363 D364 R365 R366 L367



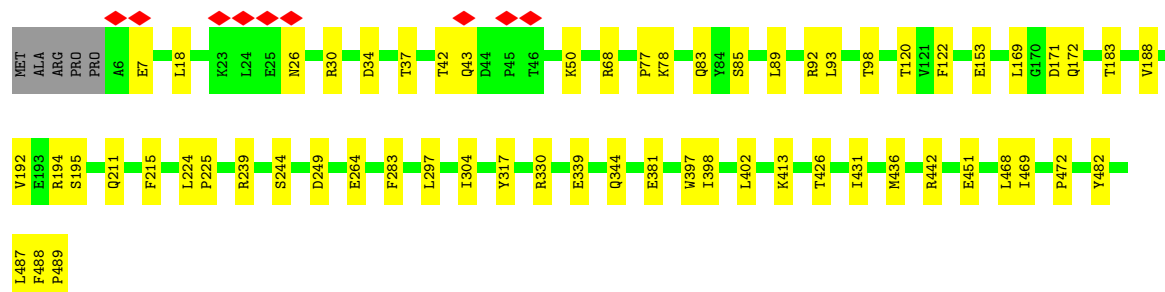
• Molecule 10: NDUFA7



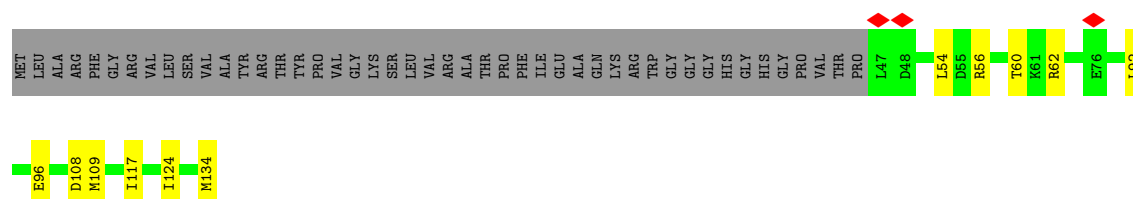
• Molecule 11: NDUFA8



• Molecule 12: NDUFA9

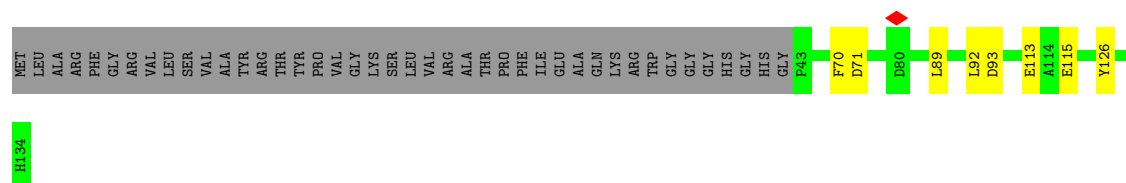


• Molecule 13: NDUFAB1-alpha

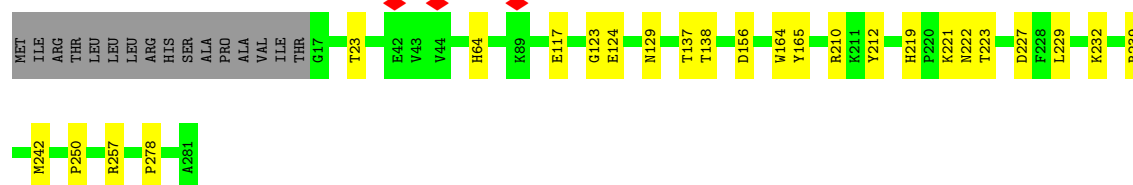
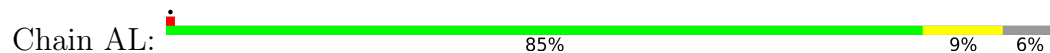


• Molecule 14: NDUFAB1-beta

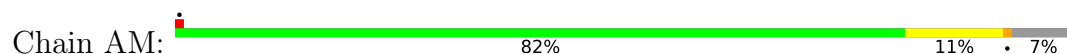




• Molecule 15: NDUFA12



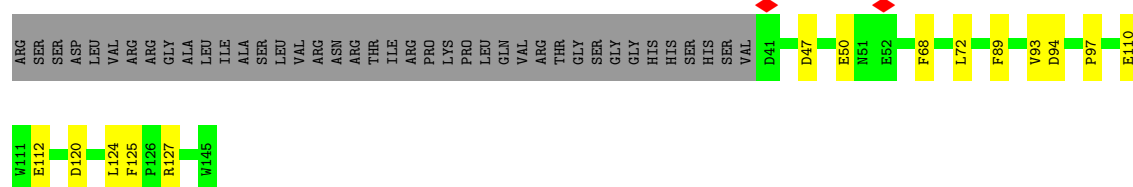
• Molecule 16: NDUFA13



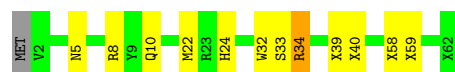
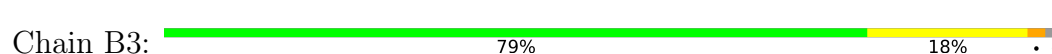
• Molecule 17: NDUFA11



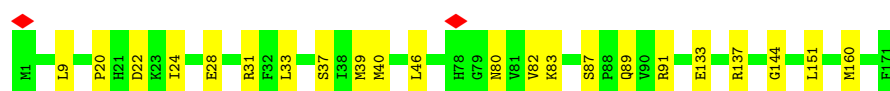
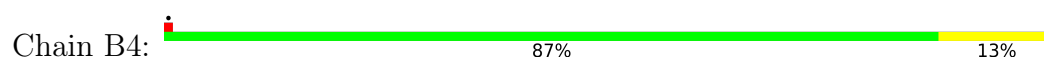
• Molecule 18: NDUFB2



• Molecule 19: NDUFB3



• Molecule 20: NDUFB4



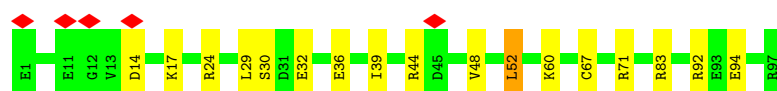
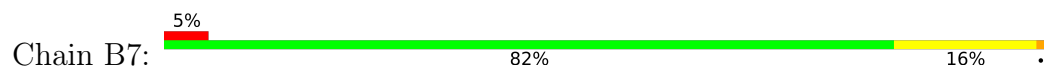
- Molecule 21: NDUFB5



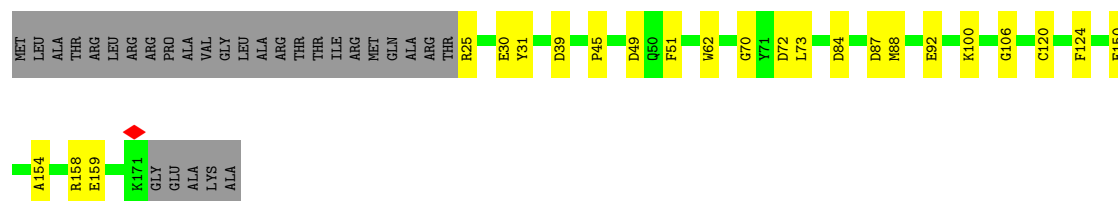
- Molecule 22: NDUFB6



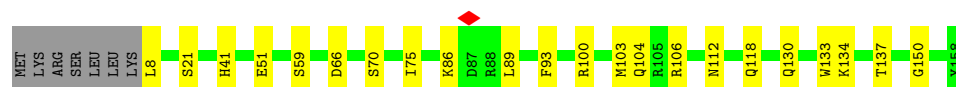
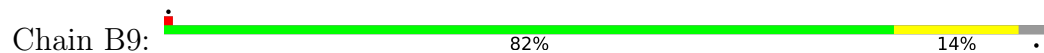
- Molecule 23: NDUFB7



- Molecule 24: NDUFB8



- Molecule 25: NDUFB9



- Molecule 26: NDUFB10

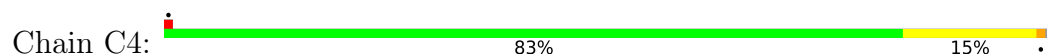




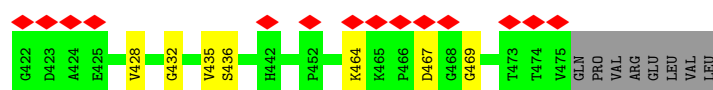
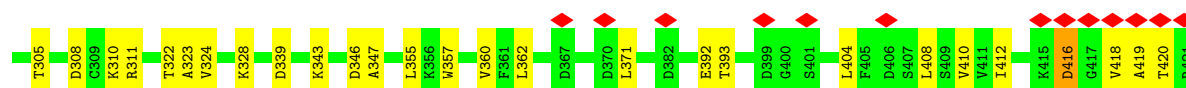
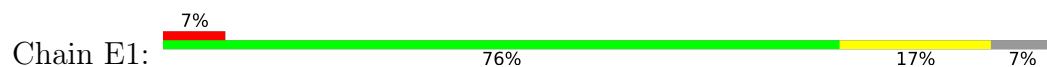
• Molecule 27: NDUFB1



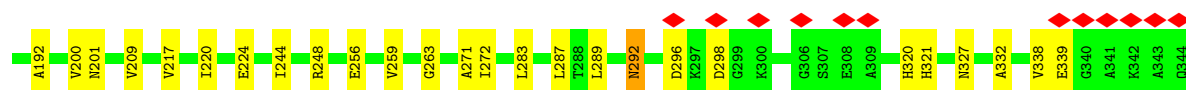
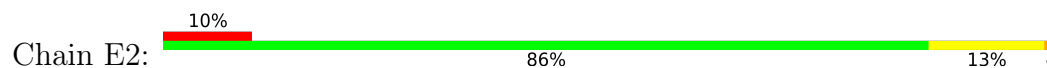
• Molecule 28: NDUF2

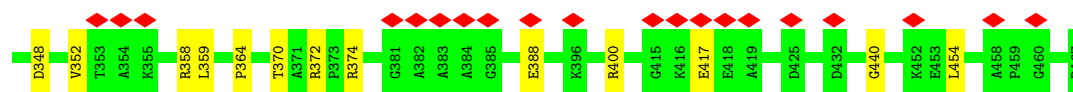


• Molecule 29: NDUEG1

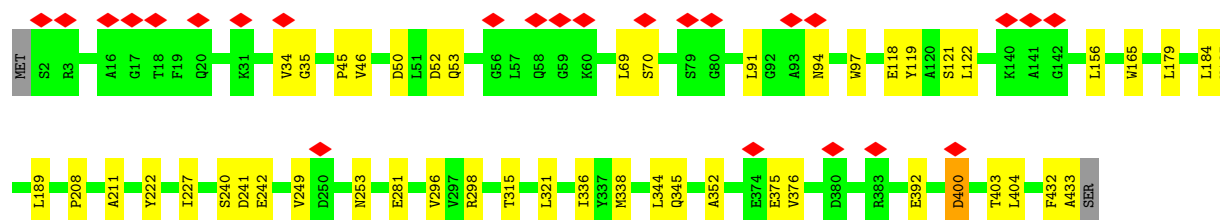


• Molecule 30: NDUEG2

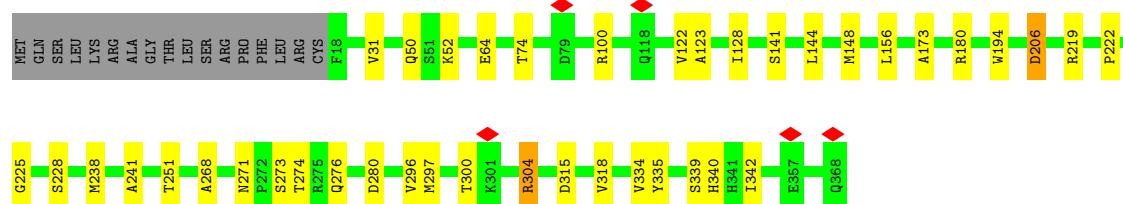
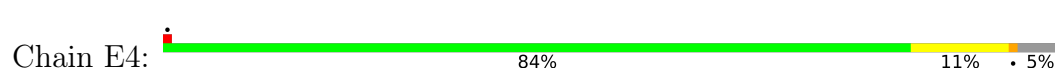




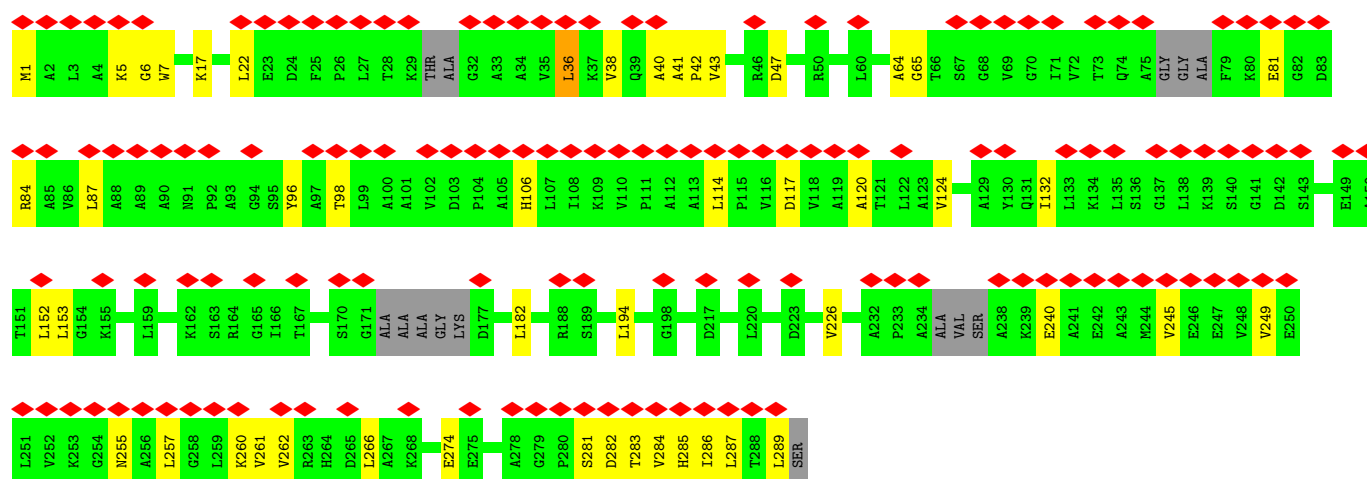
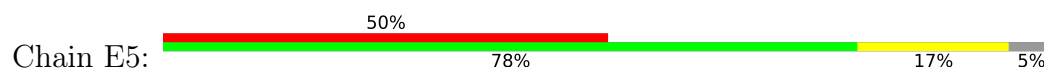
• Molecule 31: NDUEG3



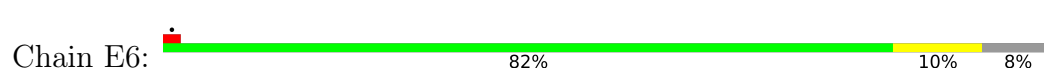
• Molecule 32: NDUEG4

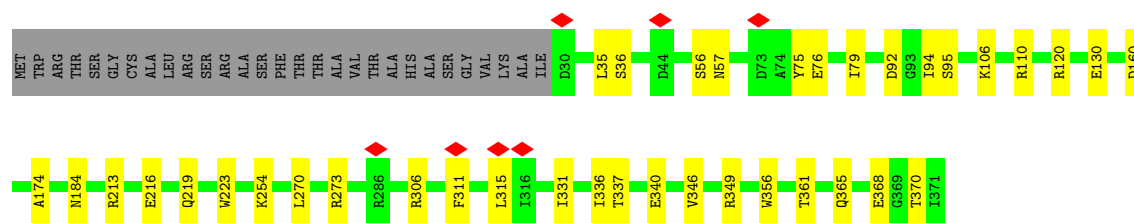


• Molecule 33: NDUEG5

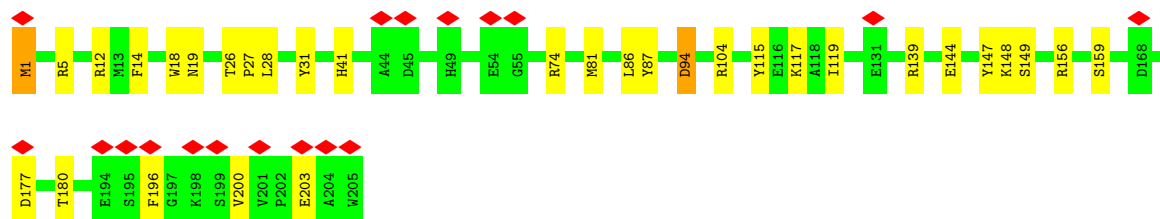
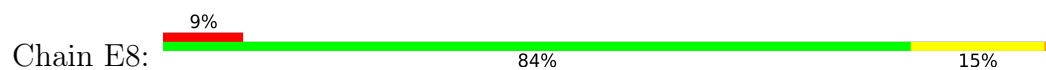


• Molecule 34: NDUEG6

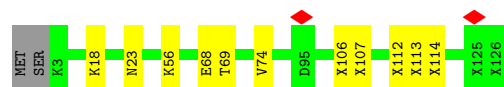




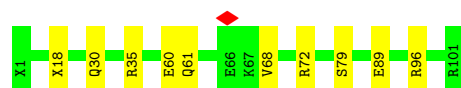
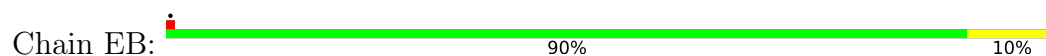
• Molecule 35: NDUEG8



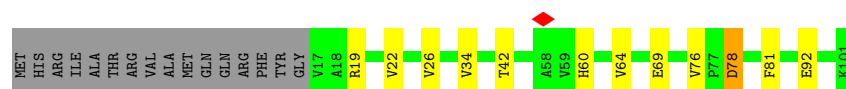
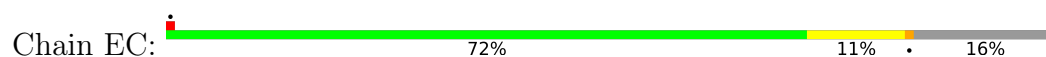
• Molecule 36: NDUEG10



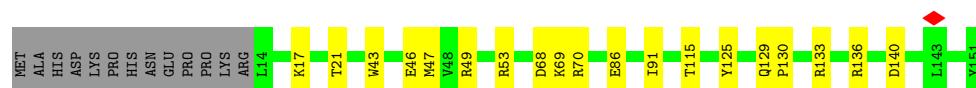
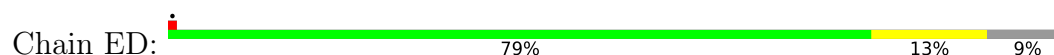
• Molecule 37: NDUEG11



• Molecule 38: NDUEG12

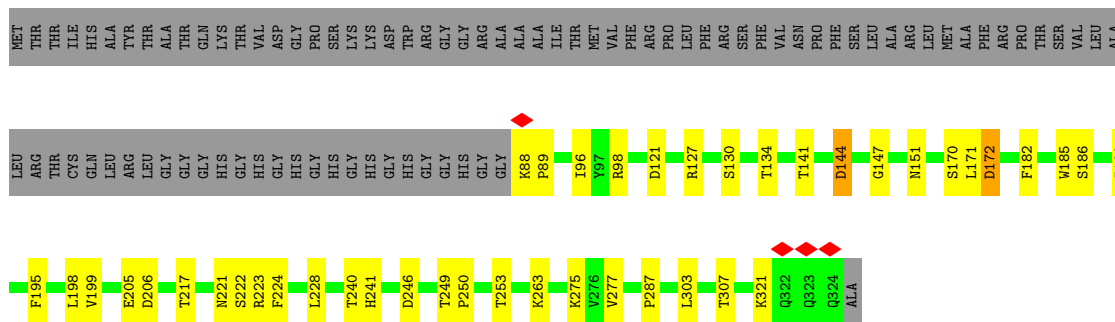


• Molecule 39: NDUEG13



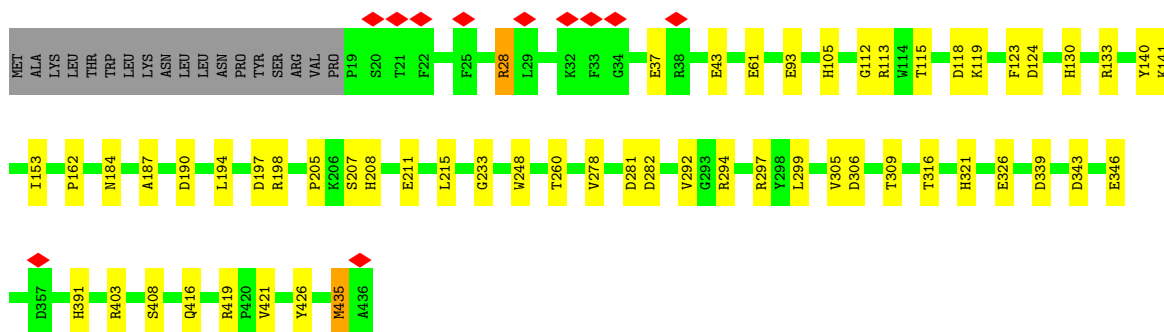
• Molecule 40: NDUFX

Chain FX:  60% 13% 27%



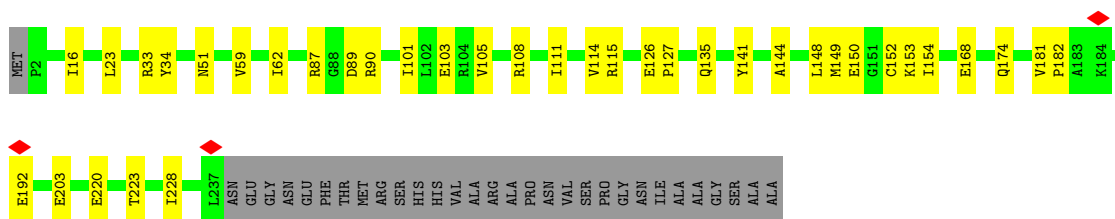
• Molecule 41: NDUCA1

Chain G1:  83% 13%



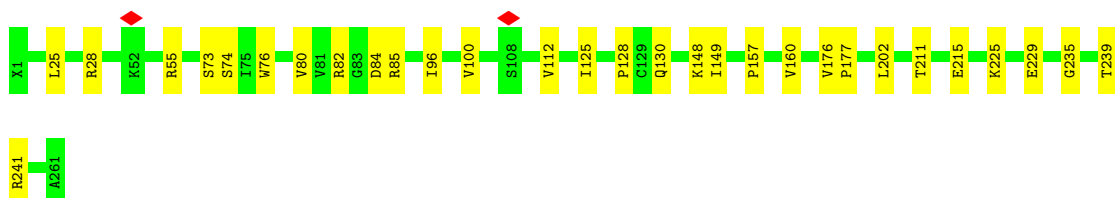
• Molecule 42: NDUCA2

Chain G2:  75% 14% 12%



• Molecule 43: NDUCA3

Chain G3:  88% 12%



• Molecule 44: ND1

Y506	S361	PHE	LEU	LEU	ASN	ILE	MET
L514	Q367	ILE	PHE	VAL	TYR	ILE	LEU
Y516	L370	PRO	LEU	VAL	HIS	ILE	ASN
L525	L373	PHE	LEU	HIS	ASN	THR	TYR
Y526		ALA	ILE	THR	SER	ILE	ILE
L527	L376	SER	LEU	PHE	LEU	LEU	ILE
L528	S377	ILE	LEU	LEU	ASN	HIS	ILE
L534	L378	ASN	PHE	GLY	ASP	ILE	TYR
Y535	S379	THR	VAL	PHE	GLN	ASN	GLY
L536	S380	THR	VAL	PHE	GLN	ASN	GLY
	L381	TYR	ASN	GLY	ILE	ILE	ILE
Y539	L382	TYR	ILE	LEU	TYR	ILE	TYR
L540	V384	HIS	TYR	GLY	ILE	ILE	MET
E541	L385	TYR	HIS	THR	GLY	ILE	SER
	E386	LYS	TYR	VAL	THR	ASN	SER
L562		TYR	ASN	ILE	ASN	ASN	ILE
L565	Q394	TYR	ASN	ILE	TYR	ILE	ILE
E566	R395	ASN	ASN	THR	TYR	ILE	ILE
	R396	TYR	LYS	ASP	ASN	GLU	ILE
	I397	ASN	SER	ASN	ASN	TYR	ILE
Y579		ASN	ILE	ASN	THR	ASN	ILE
	G403	TYR	PHE	LEU	MET	ILE	MET
L582		ASN	ASN	ILE	ILE	CYS	PHE
	D413	ILE	THR	PHE	PHE	TYR	LEU
S586	L419	ASN	THR	VAL	ILE	LEU	ASN
L587	Y419	ASN	ASN	PHE	LYS	ILE	PHE
E588		ASN	LYS	TYR	ASP	GLY	SER
Y602	Y422	THR	SER	ILE	ILE	THR	ASN
	N423	ASN	ILE	GLY	ASN	GLY	LYS
	R424	THR	TYR	LEU	ASN	LEU	ILE
L605		ASN	LEU	LEU	ILE	GLY	TYR
	N429	ASN	ILE	GLY	PHE	LEU	ILE
H608		ASN	PRO	LEU	THR	GLU	GLY
	L446	ASN	VAL	GLY	ILE	ILE	VAL
F626	L447	LYS	SER	VAL	TYR	SER	GLU
	F448	PHE	LEU	ILE	ARG	PRO	ILE
L629	I449	ILE	VAL	LEU	SER	PHE	ASN
		ASN	ASP	TYR	ILE	ILE	LYS
R640	I453	ILE	LEU	LEU	ASN	VAL	TYR
		SER	ILE	ILE	PHE	PHE	ILE
L643	I459	TYR	ASN	VAL	PHE	GLY	LEU
P644	N460	ASN	ILE	ILE	MET	TYR	SER
R645	I461	PHE	ASN	PHE	ASN	LYS	LYS
		PHE	ILE	TYR	TRP	PHE	ASN
Y657	I469	ILE	ILE	LEU	LEU	PHE	ASN
F658		ILE	PHE	GLY	TYR	VAL	ASN
Y659	L472	LEU	TYR	ILE	ILE	TYR	ASP
L660		ILE	ILE	GLN	MET	ILE	TYR
P661	H476	ILE	LEU	ILE	ILE	MET	ILE
		LEU	LEU	TYR	TYR	ASP	TRP
	I480	ASN	LEU	VAL	TYR	LEU	ILE
		ILE	TYR	TYR	GLY	ASN	ILE
	I484	ILE	THR	ILE	VAL	ASN	CYS
		ASN	LEU	SER	ARG	TYR	ASN
		ILE	LEU	PHE	MET	TRP	CYS
L670	I500	SER	LEU	SER	THR	LEU	THR

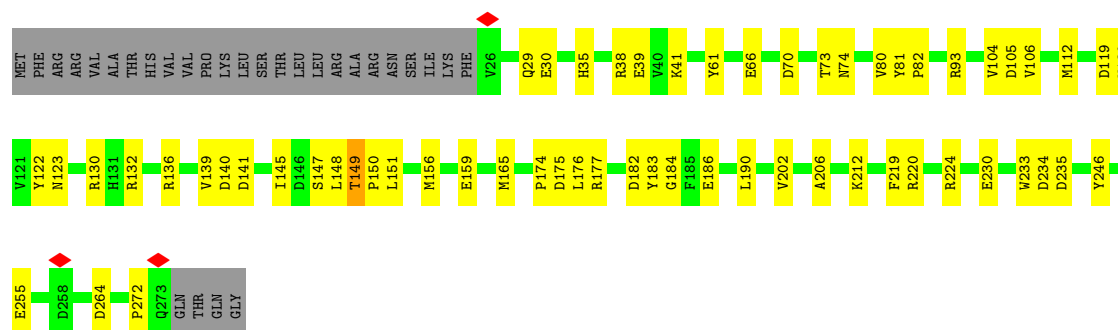
Chain N2: 80% 19%

Y160	K161	G167	G168	Y169	I171	P172		Y175	I178	L185	M186		I189	N199	P200	H209	L219	F220	L221	T247	I248	N252	L253	F257	I260	L263	L267	L281	W282	I299	ASN						
LEU	TYR	N4	N5	S15	I16	I17	I18	G19	V27	E32	N40	D42	N43	F44	I50	L58	I62	S68	T69	Q79	L80	I81	I87	I88	N89	T90	T100	V116	I120	L121	I122	E126	T139	W143	I147	S155	F156

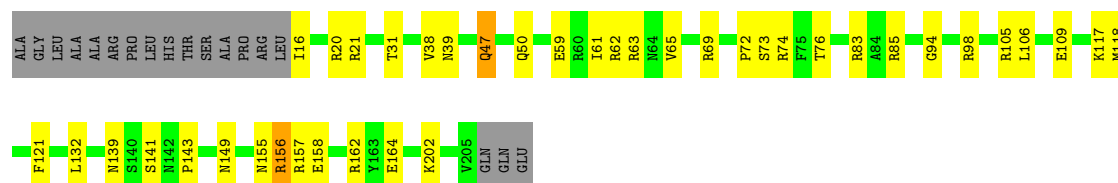
[illegible]



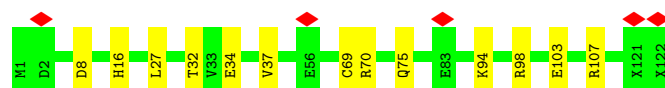
• Molecule 50: NDUFS3



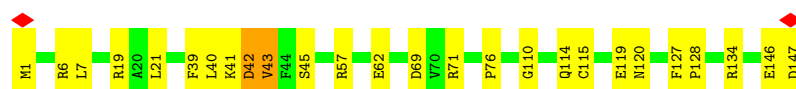
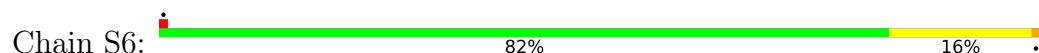
• Molecule 51: NDUFS4



• Molecule 52: NDUFS5

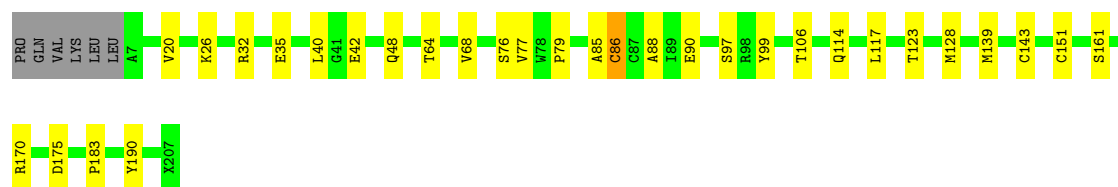


• Molecule 53: NDUFS6



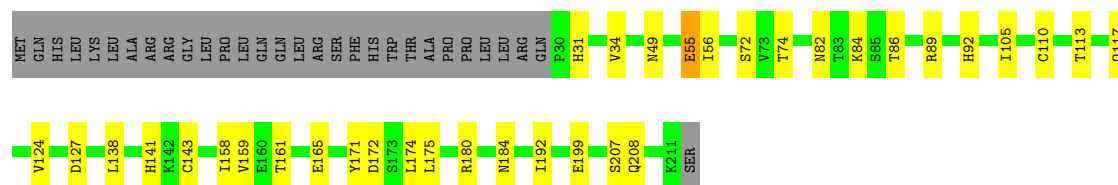
• Molecule 54: NDUFS7

Chain S7:  82% 14%



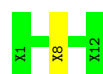
- Molecule 55: NDUFS8

Chain S8:  69% 16% 14%



- Molecule 56: UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK

Chain U1:  92% 8%



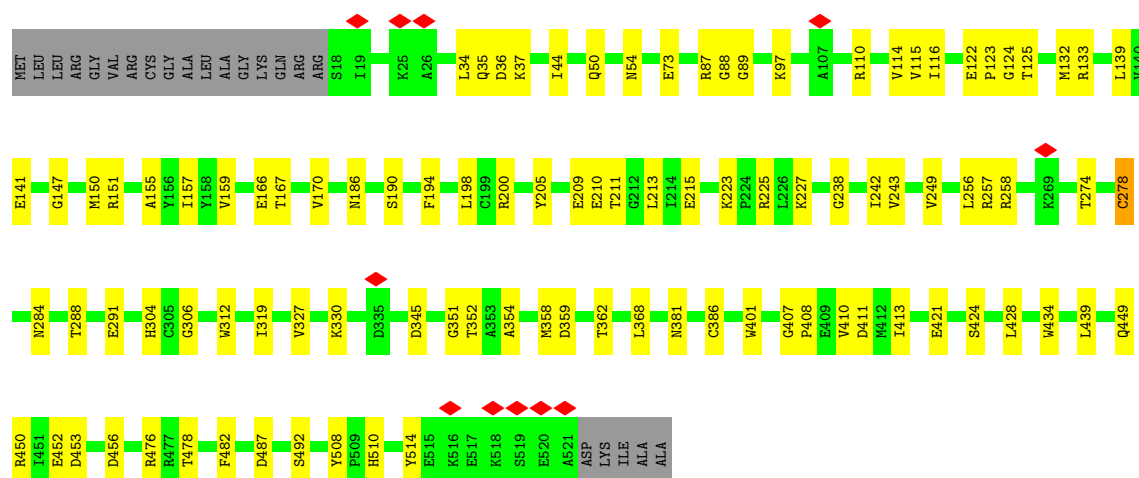
- Molecule 56: UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK-UNK

Chain U2:  8% 100%



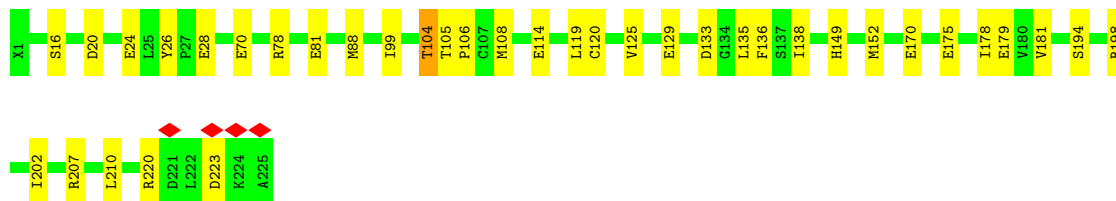
- Molecule 57: NDUFV1

Chain V1:  77% 19%



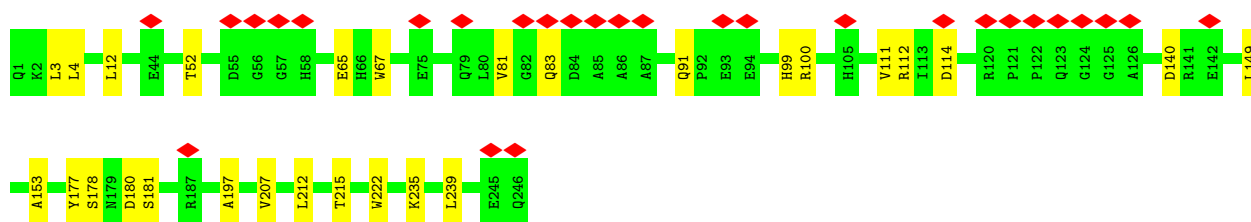
● Molecule 58: NDUFV2

Chain V2:  84% 16%



● Molecule 59: NDUEG7

Chain E7:  11% 89% 11%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	86599	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	33.507	Depositor
Minimum map value	-20.416	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.993	Depositor
Recommended contour level	4.5	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: NAI, FMN, FES, ZN, K, NDP, PC1, U10, CDL, SF4, 3PE, ZMP, 2MR

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	0.19	0/2858	0.33	0/3878
2	1B	0.17	0/4306	0.27	0/5854
3	2B	0.20	0/958	0.33	0/1306
4	4L	0.18	0/924	0.26	0/1261
5	A1	0.15	0/1108	0.27	0/1511
6	A2	0.15	0/1530	0.28	0/2089
7	A3	0.17	0/1079	0.28	0/1453
8	A5	0.16	0/1282	0.25	0/1737
9	A6	0.16	0/3395	0.26	0/4608
10	A7	0.17	0/1194	0.30	0/1619
11	A8	0.15	0/1879	0.25	0/2543
12	A9	0.17	0/3920	0.29	0/5335
13	AB	0.15	0/704	0.23	0/951
14	AC	0.15	0/736	0.22	0/1000
15	AL	0.17	0/2317	0.29	0/3136
16	AM	0.16	0/1533	0.23	0/2079
17	AN	0.15	0/2382	0.24	0/3249
18	B2	0.15	0/947	0.22	0/1291
19	B3	0.16	0/326	0.30	0/441
20	B4	0.17	0/1419	0.27	0/1922
21	B5	0.18	0/1111	0.27	0/1505
22	B6	0.17	0/803	0.26	0/1087
23	B7	0.14	0/877	0.24	0/1172
24	B8	0.17	0/1273	0.26	0/1733
25	B9	0.17	0/1274	0.28	0/1728
26	BL	0.17	0/1266	0.27	0/1710
27	BM	0.18	0/876	0.27	0/1192
28	C4	0.17	0/1592	0.29	0/2158
29	E1	0.15	0/3596	0.26	0/4879
30	E2	0.14	0/3658	0.25	0/4983
31	E3	0.14	0/3320	0.26	0/4520
32	E4	0.16	0/2850	0.25	0/3884

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
33	E5	0.12	0/2004	0.34	0/2721
34	E6	0.15	0/2954	0.26	0/4004
35	E8	0.15	0/1747	0.28	0/2367
36	EA	0.19	0/858	0.27	0/1163
37	EB	0.13	0/650	0.22	0/863
38	EC	0.16	0/676	0.28	0/925
39	ED	0.14	0/1176	0.24	0/1590
40	FX	0.20	0/2035	0.33	0/2763
41	G1	0.18	0/3374	0.29	0/4589
42	G2	0.17	0/1832	0.29	0/2476
43	G3	0.16	0/1957	0.30	0/2646
44	N1	0.18	0/2672	0.32	0/3639
45	N2	0.20	0/2582	0.32	0/3530
46	N3	0.20	0/1068	0.36	0/1456
46	N6	0.16	0/1275	0.31	0/1730
47	N4	0.20	0/4105	0.33	0/5594
48	N5	0.19	0/4963	0.33	0/6758
49	S2	0.20	0/3244	0.32	0/4403
50	S3	0.19	0/2112	0.32	0/2874
51	S4	0.16	0/1573	0.29	0/2107
52	S5	0.13	0/960	0.23	0/1291
53	S6	0.17	0/1232	0.30	0/1659
54	S7	0.19	0/1558	0.29	0/2120
55	S8	0.19	0/1485	0.29	0/2010
57	V1	0.17	0/3990	0.30	0/5394
58	V2	0.16	0/1787	0.29	0/2428
59	E7	0.14	0/1931	0.27	0/2618
All	All	0.17	0/113093	0.29	0/153532

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	34	0
2	1B	4198	4159	4175	54	0
3	2B	1070	989	1009	23	0
4	4L	890	878	880	19	0
5	A1	1071	1026	1030	7	0
6	A2	1493	1474	1478	16	0
7	A3	1050	1039	1041	9	0
8	A5	1261	1248	1251	10	0
9	A6	3328	3280	3293	40	0
10	A7	1154	1118	1123	15	0
11	A8	1822	1726	1736	34	0
12	A9	3829	3850	3857	48	0
13	AB	694	673	677	10	0
14	AC	721	697	702	9	0
15	AL	2237	2172	2180	18	0
16	AM	1487	1448	1452	24	0
17	AN	2306	2267	2275	18	0
18	B2	913	857	858	11	0
19	B3	449	309	313	9	0
20	B4	1377	1358	1364	16	0
21	B5	1112	1069	1075	14	0
22	B6	773	747	751	6	0
23	B7	857	835	841	13	0
24	B8	1224	1127	1136	23	0
25	B9	1236	1207	1212	19	0
26	BL	1227	1179	1185	10	0
27	BM	910	827	830	13	0
28	C4	1545	1517	1519	21	0
29	E1	3512	3496	3510	50	0
30	E2	3563	3540	3554	37	0
31	E3	3255	3263	3279	37	0
32	E4	2770	2732	2742	32	0
33	E5	1977	2069	2075	41	0
34	E6	2871	2758	2767	27	0
35	E8	1691	1663	1668	26	0
36	EA	961	832	835	10	0
37	EB	774	631	636	7	0
38	EC	660	663	666	10	0
39	ED	1142	1131	1134	18	0
40	FX	1967	1849	1858	32	0
41	G1	3281	3139	3156	51	0
42	G2	1804	1846	1850	30	0
43	G3	1961	1944	1950	31	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
44	N1	2605	2726	2729	50	0
45	N2	2512	2589	2592	50	0
46	N3	1037	1057	1057	26	0
46	N6	1257	1385	1385	35	0
47	N4	4001	4214	4224	77	0
48	N5	4837	5032	5046	96	0
49	S2	3173	3101	3114	60	0
50	S3	2050	1928	1936	45	0
51	S4	1536	1502	1505	32	0
52	S5	991	895	898	10	0
53	S6	1200	1192	1198	24	0
54	S7	1545	1500	1503	25	0
55	S8	1451	1392	1397	30	0
56	U1	60	16	18	1	0
56	U2	60	16	17	0	0
57	V1	3897	3827	3837	70	0
58	V2	1759	1701	1710	27	0
59	E7	1888	1892	1903	19	0
60	1A	4	0	0	0	0
60	V2	4	0	0	0	0
61	1A	16	0	0	1	0
61	S7	8	0	0	0	0
61	S8	16	0	0	2	0
61	V1	8	0	0	1	0
62	1A	1	0	0	0	0
63	A1	80	111	111	2	0
63	A9	66	80	80	1	0
63	AL	50	77	77	0	0
63	AM	97	148	148	3	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	1	0
63	ED	54	88	88	2	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	1	0
63	N5	131	190	190	3	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	7	0
64	B3	65	74	74	3	0
64	B5	58	60	60	0	0
64	C4	163	223	223	2	0
64	E6	64	72	72	0	0
64	E7	68	80	80	0	0
64	EA	114	116	116	2	0
64	N4	98	149	149	2	0
64	N5	163	223	223	1	0
65	A9	48	26	26	0	0
66	AB	36	0	47	5	0
66	AC	36	0	47	4	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	0	0
68	N4	43	55	55	4	0
69	E7	1	0	0	0	0
69	S6	1	0	0	0	0
70	V1	31	19	19	2	0
71	V1	44	0	27	6	0
All	All	113966	112888	113418	1321	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 6.

The worst 5 of 1321 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)
47:N4:43:ILE:HD13	47:N4:472:LEU:HD21	1.39	1.04
33:E5:287:LEU:O	33:E5:289:LEU:N	1.98	0.96
47:N4:293:THR:HG21	47:N4:365:ILE:HD11	1.47	0.96
31:E3:222:TYR:HH	38:EC:60:HIS:HE2	0.98	0.91
9:A6:45:HIS:NE2	9:A6:117:GLU:OE1	2.03	0.91

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%)	337 (96%)	13 (4%)	0	100	100
2	1B	523/527 (99%)	507 (97%)	16 (3%)	0	100	100
3	2B	112/142 (79%)	106 (95%)	6 (5%)	0	100	100
4	4L	106/171 (62%)	104 (98%)	2 (2%)	0	100	100
5	A1	135/141 (96%)	125 (93%)	10 (7%)	0	100	100
6	A2	190/193 (98%)	186 (98%)	4 (2%)	0	100	100
7	A3	122/125 (98%)	118 (97%)	4 (3%)	0	100	100
8	A5	152/184 (83%)	148 (97%)	4 (3%)	0	100	100
9	A6	421/437 (96%)	404 (96%)	17 (4%)	0	100	100
10	A7	134/136 (98%)	124 (92%)	10 (8%)	0	100	100
11	A8	221/223 (99%)	215 (97%)	6 (3%)	0	100	100
12	A9	482/489 (99%)	466 (97%)	16 (3%)	0	100	100
13	AB	86/134 (64%)	86 (100%)	0	0	100	100
14	AC	90/134 (67%)	90 (100%)	0	0	100	100
15	AL	263/281 (94%)	248 (94%)	15 (6%)	0	100	100
16	AM	182/198 (92%)	177 (97%)	5 (3%)	0	100	100
17	AN	285/287 (99%)	281 (99%)	4 (1%)	0	100	100
18	B2	103/145 (71%)	102 (99%)	1 (1%)	0	100	100
19	B3	32/62 (52%)	31 (97%)	0	1 (3%)	3	12
20	B4	169/171 (99%)	159 (94%)	10 (6%)	0	100	100
21	B5	132/140 (94%)	129 (98%)	3 (2%)	0	100	100
22	B6	89/91 (98%)	87 (98%)	2 (2%)	0	100	100
23	B7	95/97 (98%)	94 (99%)	1 (1%)	0	100	100
24	B8	145/176 (82%)	143 (99%)	2 (1%)	0	100	100
25	B9	149/158 (94%)	140 (94%)	9 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
26	BL	142/144 (99%)	138 (97%)	4 (3%)	0	100	100
27	BM	99/112 (88%)	96 (97%)	3 (3%)	0	100	100
28	C4	181/185 (98%)	177 (98%)	4 (2%)	0	100	100
29	E1	448/483 (93%)	432 (96%)	16 (4%)	0	100	100
30	E2	464/467 (99%)	450 (97%)	13 (3%)	1 (0%)	43	70
31	E3	430/434 (99%)	423 (98%)	7 (2%)	0	100	100
32	E4	349/368 (95%)	340 (97%)	9 (3%)	0	100	100
33	E5	266/290 (92%)	245 (92%)	21 (8%)	0	100	100
34	E6	340/371 (92%)	334 (98%)	6 (2%)	0	100	100
35	E8	203/205 (99%)	194 (96%)	9 (4%)	0	100	100
36	EA	96/126 (76%)	91 (95%)	5 (5%)	0	100	100
37	EB	73/101 (72%)	73 (100%)	0	0	100	100
38	EC	83/101 (82%)	76 (92%)	7 (8%)	0	100	100
39	ED	136/151 (90%)	132 (97%)	4 (3%)	0	100	100
40	FX	235/325 (72%)	224 (95%)	11 (5%)	0	100	100
41	G1	416/436 (95%)	405 (97%)	11 (3%)	0	100	100
42	G2	234/267 (88%)	222 (95%)	12 (5%)	0	100	100
43	G3	253/261 (97%)	239 (94%)	14 (6%)	0	100	100
44	N1	308/670 (46%)	290 (94%)	18 (6%)	0	100	100
45	N2	294/300 (98%)	278 (95%)	16 (5%)	0	100	100
46	N3	119/293 (41%)	116 (98%)	3 (2%)	0	100	100
46	N6	152/293 (52%)	149 (98%)	3 (2%)	0	100	100
47	N4	476/478 (100%)	464 (98%)	12 (2%)	0	100	100
48	N5	582/584 (100%)	562 (97%)	20 (3%)	0	100	100
49	S2	391/395 (99%)	373 (95%)	18 (5%)	0	100	100
50	S3	246/277 (89%)	238 (97%)	8 (3%)	0	100	100
51	S4	188/208 (90%)	178 (95%)	10 (5%)	0	100	100
52	S5	110/122 (90%)	106 (96%)	4 (4%)	0	100	100
53	S6	145/147 (99%)	142 (98%)	3 (2%)	0	100	100
54	S7	195/207 (94%)	188 (96%)	7 (4%)	0	100	100
55	S8	180/212 (85%)	176 (98%)	4 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
57	V1	502/526 (95%)	481 (96%)	21 (4%)	0	100	100
58	V2	220/225 (98%)	215 (98%)	5 (2%)	0	100	100
59	E7	244/246 (99%)	236 (97%)	8 (3%)	0	100	100
All	All	13568/15237 (89%)	13090 (96%)	476 (4%)	2 (0%)	100	100

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
19	B3	34	ARG
30	E2	370	THR

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%)	303 (98%)	7 (2%)	44	74
2	1B	453/454 (100%)	452 (100%)	1 (0%)	87	96
3	2B	109/111 (98%)	109 (100%)	0	100	100
4	4L	96/151 (64%)	95 (99%)	1 (1%)	68	87
5	A1	115/118 (98%)	113 (98%)	2 (2%)	53	80
6	A2	159/160 (99%)	158 (99%)	1 (1%)	78	92
7	A3	104/104 (100%)	104 (100%)	0	100	100
8	A5	134/152 (88%)	132 (98%)	2 (2%)	57	82
9	A6	278/358 (78%)	276 (99%)	2 (1%)	76	90
12	A9	5/424 (1%)	5 (100%)	0	100	100
15	AL	127/242 (52%)	125 (98%)	2 (2%)	55	81
16	AM	131/168 (78%)	129 (98%)	2 (2%)	57	82
17	AN	193/241 (80%)	192 (100%)	1 (0%)	81	93
18	B2	97/131 (74%)	97 (100%)	0	100	100
19	B3	30/31 (97%)	30 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
20	B4	144/144 (100%)	143 (99%)	1 (1%)	76	90
21	B5	108/108 (100%)	108 (100%)	0	100	100
22	B6	82/82 (100%)	82 (100%)	0	100	100
23	B7	93/93 (100%)	92 (99%)	1 (1%)	65	86
24	B8	127/148 (86%)	127 (100%)	0	100	100
25	B9	132/139 (95%)	130 (98%)	2 (2%)	57	82
26	BL	132/132 (100%)	130 (98%)	2 (2%)	57	82
27	BM	93/93 (100%)	93 (100%)	0	100	100
28	C4	166/167 (99%)	162 (98%)	4 (2%)	43	73
29	E1	381/404 (94%)	375 (98%)	6 (2%)	55	81
30	E2	379/380 (100%)	371 (98%)	8 (2%)	47	75
31	E3	339/341 (99%)	337 (99%)	2 (1%)	78	92
32	E4	302/317 (95%)	300 (99%)	2 (1%)	76	90
33	E5	200/205 (98%)	198 (99%)	2 (1%)	68	87
34	E6	293/314 (93%)	290 (99%)	3 (1%)	68	87
35	E8	179/179 (100%)	176 (98%)	3 (2%)	53	80
36	EA	84/86 (98%)	84 (100%)	0	100	100
37	EB	70/70 (100%)	69 (99%)	1 (1%)	59	83
38	EC	73/86 (85%)	72 (99%)	1 (1%)	59	83
39	ED	121/133 (91%)	120 (99%)	1 (1%)	73	89
40	FX	212/276 (77%)	208 (98%)	4 (2%)	50	77
41	G1	348/365 (95%)	343 (99%)	5 (1%)	59	83
42	G2	192/214 (90%)	188 (98%)	4 (2%)	47	75
43	G3	202/202 (100%)	202 (100%)	0	100	100
44	N1	295/639 (46%)	291 (99%)	4 (1%)	59	83
45	N2	285/289 (99%)	281 (99%)	4 (1%)	59	83
46	N3	116/281 (41%)	112 (97%)	4 (3%)	32	64
46	N6	147/281 (52%)	144 (98%)	3 (2%)	48	76
47	N4	455/455 (100%)	447 (98%)	8 (2%)	51	78
48	N5	546/546 (100%)	539 (99%)	7 (1%)	61	84
49	S2	335/336 (100%)	327 (98%)	8 (2%)	43	73

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
50	S3	224/250 (90%)	220 (98%)	4 (2%)	51	78
51	S4	159/172 (92%)	155 (98%)	4 (2%)	42	72
52	S5	102/102 (100%)	102 (100%)	0	100	100
53	S6	130/130 (100%)	128 (98%)	2 (2%)	57	82
54	S7	132/171 (77%)	130 (98%)	2 (2%)	57	82
55	S8	109/187 (58%)	107 (98%)	2 (2%)	51	78
57	V1	361/427 (84%)	358 (99%)	3 (1%)	73	89
58	V2	190/190 (100%)	185 (97%)	5 (3%)	40	71
59	E7	192/192 (100%)	190 (99%)	2 (1%)	68	87
All	All	10571/12511 (84%)	10436 (99%)	135 (1%)	59	84

5 of 135 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
51	S4	16	ILE
53	S6	42	ASP
58	V2	120	CYS
32	E4	304	ARG
32	E4	206	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 230 such sidechains are listed below:

Mol	Chain	Res	Type
34	E6	147	GLN
58	V2	42	GLN
42	G2	50	HIS
57	V1	282	HIS
49	S2	45	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	$\# Z > 2$
49	2MR	S2	154	49	10,12,13	2.50	2 (20%)	5,13,15	0.90	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
49	2MR	S2	154	49	-	2/10/13/15	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
49	S2	154	2MR	CZ-NH2	5.42	1.44	1.33
49	S2	154	2MR	CZ-NE	5.12	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
49	S2	154	2MR	CG-CD-NE-CZ
49	S2	154	2MR	N-CA-CB-CG

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 65 ligands modelled in this entry, 3 are monoatomic - leaving 62 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
61	SF4	1A	403	1	0,12,12	-	-	-		
64	CDL	AM	216	-	71,71,99	0.35	0	77,83,111	0.37	0
63	PC1	E8	301	-	53,53,53	0.29	0	59,61,61	0.32	0
63	PC1	N5	606	-	35,35,53	0.35	0	41,43,61	0.33	0
64	CDL	B5	201	-	57,57,99	0.38	0	63,69,111	0.34	0
64	CDL	EA	201	-	58,58,99	0.39	0	64,70,111	0.36	0
64	CDL	C4	202	-	93,93,99	0.31	0	99,105,111	0.30	0
71	NAI	V1	581	-	47,48,48	0.29	0	64,73,73	0.37	0
64	CDL	AL	302	-	67,67,99	0.36	0	73,79,111	0.32	0
63	PC1	A1	202	-	48,48,53	0.31	0	54,56,61	0.33	0
61	SF4	1A	402	1	0,12,12	-	-	-		
63	PC1	AM	218	-	48,48,53	0.31	0	54,56,61	0.30	0
63	PC1	E8	302	-	53,53,53	0.30	0	59,61,61	0.29	0
63	PC1	E8	303	-	32,32,53	0.36	0	38,40,61	0.39	0
63	PC1	N5	601	-	53,53,53	0.29	0	59,61,61	0.30	0
66	ZMP	AB	150	13	30,35,36	0.74	1 (3%)	34,42,45	0.92	1 (2%)
63	PC1	A9	561	-	32,32,53	0.36	0	38,40,61	0.36	0
63	PC1	B5	203	-	53,53,53	0.30	0	59,61,61	0.33	0
64	CDL	EA	202	-	54,54,99	0.39	0	60,66,111	0.35	0
67	3PE	AN	302	-	50,50,50	0.30	0	53,55,55	0.30	0
65	NDP	A9	559	-	51,52,52	0.37	0	71,80,80	0.40	0
67	3PE	N4	504	-	40,40,50	0.34	0	43,45,55	0.32	0
64	CDL	N4	501	-	97,97,99	0.30	0	103,109,111	0.28	0
68	U10	N4	505	-	43,43,63	2.48	20 (46%)	54,55,79	1.79	15 (27%)
64	CDL	E7	301	-	67,67,99	0.36	0	73,79,111	0.31	0
63	PC1	E8	304	-	29,29,53	0.38	0	35,37,61	0.34	0
61	SF4	V1	580	57	0,12,12	-	-	-		
63	PC1	B5	202	-	53,53,53	0.29	0	59,61,61	0.33	0
60	FES	V2	301	58	0,4,4	-	-	-		
64	CDL	AM	215	-	71,71,99	0.35	0	77,83,111	0.30	0
63	PC1	AM	220	-	47,47,53	0.31	0	53,55,61	0.27	0
61	SF4	S8	298	55	0,12,12	-	-	-		

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
63	PC1	AN	301	-	47,47,53	0.31	0	53,55,61	0.30	0
63	PC1	N1	701	-	48,48,53	0.30	0	54,56,61	0.30	0
61	SF4	S7	301	54	0,12,12	-	-	-		
63	PC1	N3	301	-	41,41,53	0.33	0	47,49,61	0.40	0
64	CDL	N5	603	-	69,69,99	0.35	0	75,81,111	0.32	0
67	3PE	G1	516	-	39,39,50	0.34	0	42,44,55	0.32	0
70	FMN	V1	579	-	33,33,33	0.32	0	48,50,50	0.32	0
63	PC1	A1	203	-	30,30,53	0.37	0	36,38,61	0.37	0
63	PC1	A9	560	-	32,32,53	0.37	0	38,40,61	0.36	0
63	PC1	C4	203	-	37,37,53	0.34	0	43,45,61	0.33	0
64	CDL	A3	201	-	57,57,99	0.39	0	63,69,111	0.34	0
64	CDL	B3	102	-	64,64,99	0.37	0	70,76,111	0.35	0
60	FES	1A	401	1	0,4,4	-	-	-		
63	PC1	N5	605	-	40,40,53	0.32	0	46,48,61	0.32	0
64	CDL	AL	304	-	69,69,99	0.36	0	75,81,111	0.31	0
63	PC1	N4	502	-	38,38,53	0.34	0	44,46,61	0.33	0
67	3PE	N5	607	-	50,50,50	0.31	0	53,55,55	0.32	0
63	PC1	AL	301	-	49,49,53	0.31	0	55,57,61	0.30	0
64	CDL	E6	431	-	63,63,99	0.37	0	69,75,111	0.33	0
64	CDL	C4	204	-	68,68,99	0.35	0	74,80,111	0.34	0
61	SF4	S8	297	55	0,12,12	-	-	-		
64	CDL	AL	303	-	63,63,99	0.37	0	69,75,111	0.32	0
63	PC1	E4	401	-	50,50,53	0.29	0	56,58,61	0.30	0
63	PC1	ED	201	-	53,53,53	0.29	0	59,61,61	0.32	0
63	PC1	N4	503	-	32,32,53	0.37	0	38,40,61	0.36	0
66	ZMP	AC	201	14	30,35,36	0.70	1 (3%)	34,42,45	0.86	1 (2%)
63	PC1	N1	702	-	39,39,53	0.33	0	45,47,61	0.31	0
64	CDL	AM	217	-	71,71,99	0.35	0	77,83,111	0.33	0
63	PC1	N2	301	-	36,36,53	0.35	0	42,44,61	0.36	0
64	CDL	N5	608	-	92,92,99	0.31	0	98,104,111	0.30	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
61	SF4	1A	403	1	-	-	0/6/5/5
64	CDL	AM	216	-	-	16/82/82/110	-
63	PC1	E8	301	-	-	7/57/57/57	-
63	PC1	N5	606	-	-	6/39/39/57	-
64	CDL	B5	201	-	-	7/68/68/110	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
64	CDL	EA	201	-	-	11/69/69/110	-
64	CDL	C4	202	-	-	20/104/104/110	-
71	NAI	V1	581	-	-	7/29/72/72	0/5/5/5
64	CDL	AL	302	-	-	13/78/78/110	-
63	PC1	A1	202	-	-	17/52/52/57	-
61	SF4	1A	402	1	-	-	0/6/5/5
63	PC1	AM	218	-	-	17/52/52/57	-
63	PC1	E8	302	-	-	20/57/57/57	-
63	PC1	E8	303	-	-	4/36/36/57	-
63	PC1	N5	601	-	-	10/57/57/57	-
66	ZMP	AB	150	13	-	16/40/42/43	-
63	PC1	A9	561	-	-	5/36/36/57	-
63	PC1	B5	203	-	-	20/57/57/57	-
64	CDL	EA	202	-	-	15/65/65/110	-
67	3PE	AN	302	-	-	8/54/54/54	-
65	NDP	A9	559	-	-	3/34/77/77	0/5/5/5
67	3PE	N4	504	-	-	9/44/44/54	-
64	CDL	N4	501	-	-	18/108/108/110	-
68	U10	N4	505	-	-	8/39/63/87	0/1/1/1
64	CDL	E7	301	-	-	19/78/78/110	-
63	PC1	E8	304	-	-	10/33/33/57	-
61	SF4	V1	580	57	-	-	0/6/5/5
63	PC1	B5	202	-	-	19/57/57/57	-
60	FES	V2	301	58	-	-	0/1/1/1
64	CDL	AM	215	-	-	21/82/82/110	-
63	PC1	AM	220	-	-	9/51/51/57	-
70	FMN	V1	579	-	-	2/18/18/18	0/3/3/3
63	PC1	AN	301	-	-	11/51/51/57	-
63	PC1	N1	701	-	-	16/52/52/57	-
63	PC1	N3	301	-	-	10/45/45/57	-
67	3PE	G1	516	-	-	9/43/43/54	-
64	CDL	N5	603	-	-	10/80/80/110	-
61	SF4	S7	301	54	-	-	0/6/5/5
61	SF4	S8	298	55	-	-	0/6/5/5
63	PC1	A1	203	-	-	6/34/34/57	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
63	PC1	A9	560	-	-	10/36/36/57	-
63	PC1	C4	203	-	-	9/41/41/57	-
64	CDL	A3	201	-	-	6/68/68/110	-
64	CDL	B3	102	-	-	15/75/75/110	-
60	FES	1A	401	1	-	-	0/1/1/1
63	PC1	N5	605	-	-	13/44/44/57	-
64	CDL	AL	304	-	-	27/80/80/110	-
63	PC1	N4	502	-	-	11/42/42/57	-
67	3PE	N5	607	-	-	10/54/54/54	-
63	PC1	AL	301	-	-	11/53/53/57	-
64	CDL	E6	431	-	-	24/74/74/110	-
64	CDL	C4	204	-	-	9/79/79/110	-
64	CDL	AL	303	-	-	20/74/74/110	-
61	SF4	S8	297	55	-	-	0/6/5/5
63	PC1	E4	401	-	-	13/54/54/57	-
63	PC1	ED	201	-	-	10/57/57/57	-
63	PC1	N4	503	-	-	6/36/36/57	-
66	ZMP	AC	201	14	-	24/40/42/43	-
63	PC1	N1	702	-	-	10/43/43/57	-
64	CDL	AM	217	-	-	20/82/82/110	-
63	PC1	N2	301	-	-	14/40/40/57	-
64	CDL	N5	608	-	-	17/103/103/110	-

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	10.41	1.53	1.35
68	N4	505	U10	C4-C3	4.81	1.53	1.36
68	N4	505	U10	C7-C8	2.84	1.55	1.50
68	N4	505	U10	C31-C29	2.75	1.57	1.51
66	AB	150	ZMP	C9-C10	-2.73	1.48	1.50

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
68	N4	505	U10	C30-C29-C31	4.29	122.67	115.23
68	N4	505	U10	C7-C6-C1	-4.08	117.90	124.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
68	N4	505	U10	C7-C8-C9	-4.00	119.94	126.83
68	N4	505	U10	C7-C6-C5	3.56	122.66	118.52
68	N4	505	U10	C15-C14-C16	3.19	120.77	115.23

There are no chirality outliers.

5 of 678 torsion outliers are listed below:

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

There are no ring outliers.

33 monomers are involved in 62 short contacts:

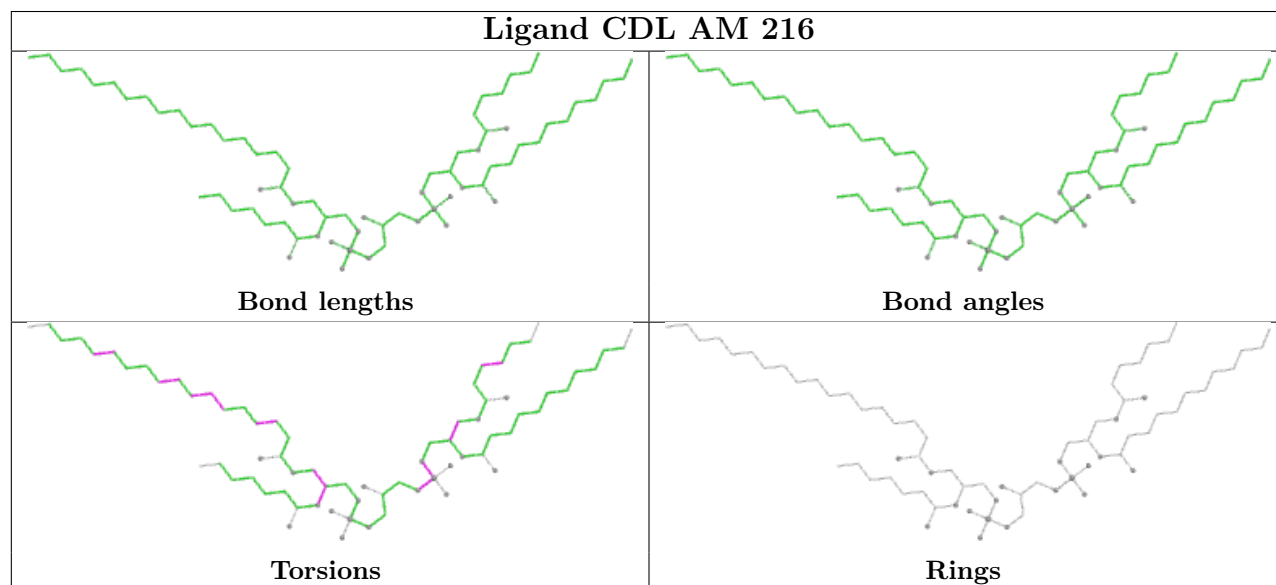
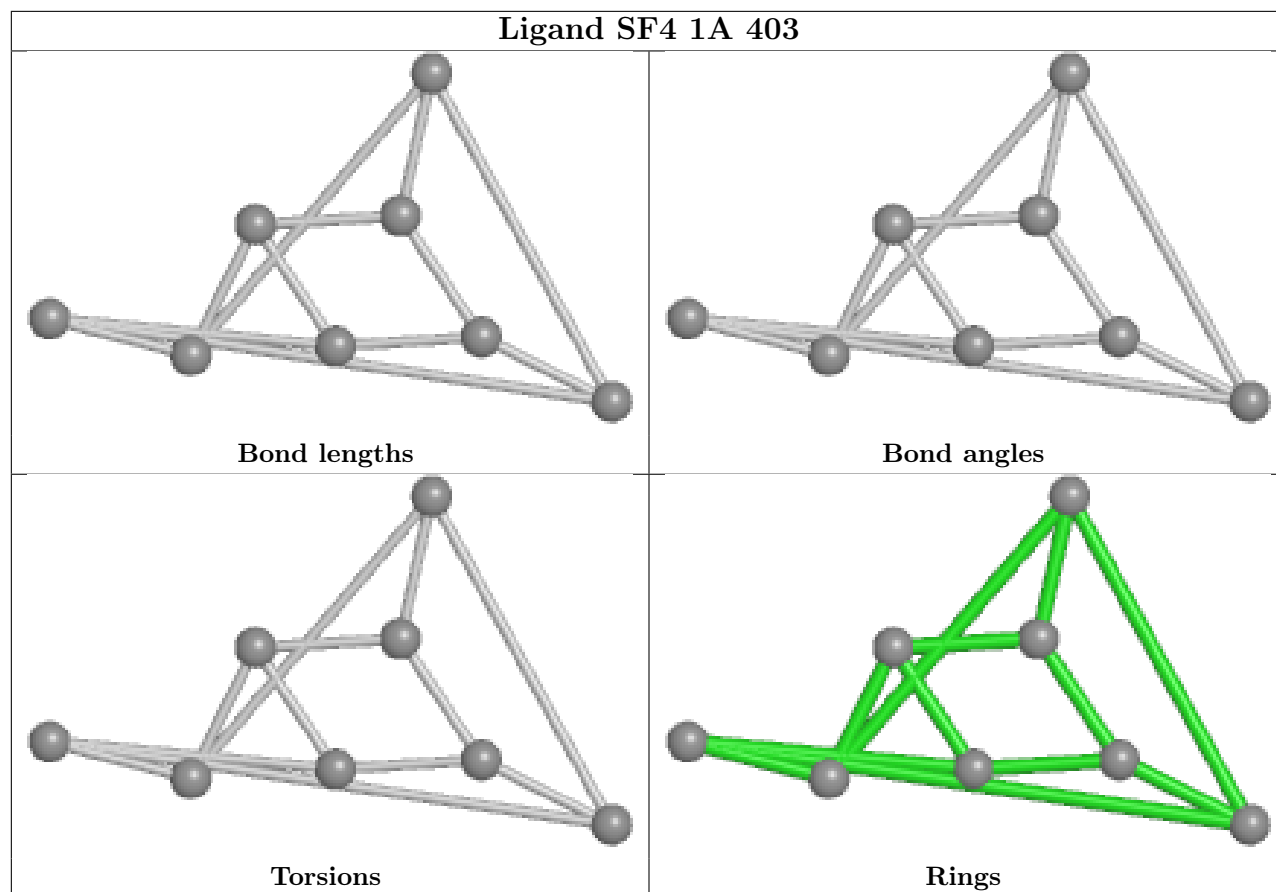
Mol	Chain	Res	Type	Clashes	Symm-Clashes
61	1A	403	SF4	1	0
64	AM	216	CDL	2	0
63	E8	301	PC1	1	0
64	EA	201	CDL	1	0
64	C4	202	CDL	1	0
71	V1	581	NAI	6	0
63	A1	202	PC1	1	0
63	AM	218	PC1	1	0
63	N5	601	PC1	3	0
66	AB	150	ZMP	5	0
63	B5	203	PC1	2	0
64	EA	202	CDL	1	0
64	N4	501	CDL	2	0
68	N4	505	U10	4	0
61	V1	580	SF4	1	0
64	AM	215	CDL	4	0
63	AM	220	PC1	2	0
64	N5	603	CDL	1	0
67	G1	516	3PE	2	0
70	V1	579	FMN	2	0
63	A1	203	PC1	1	0
63	A9	560	PC1	1	0
63	C4	203	PC1	1	0

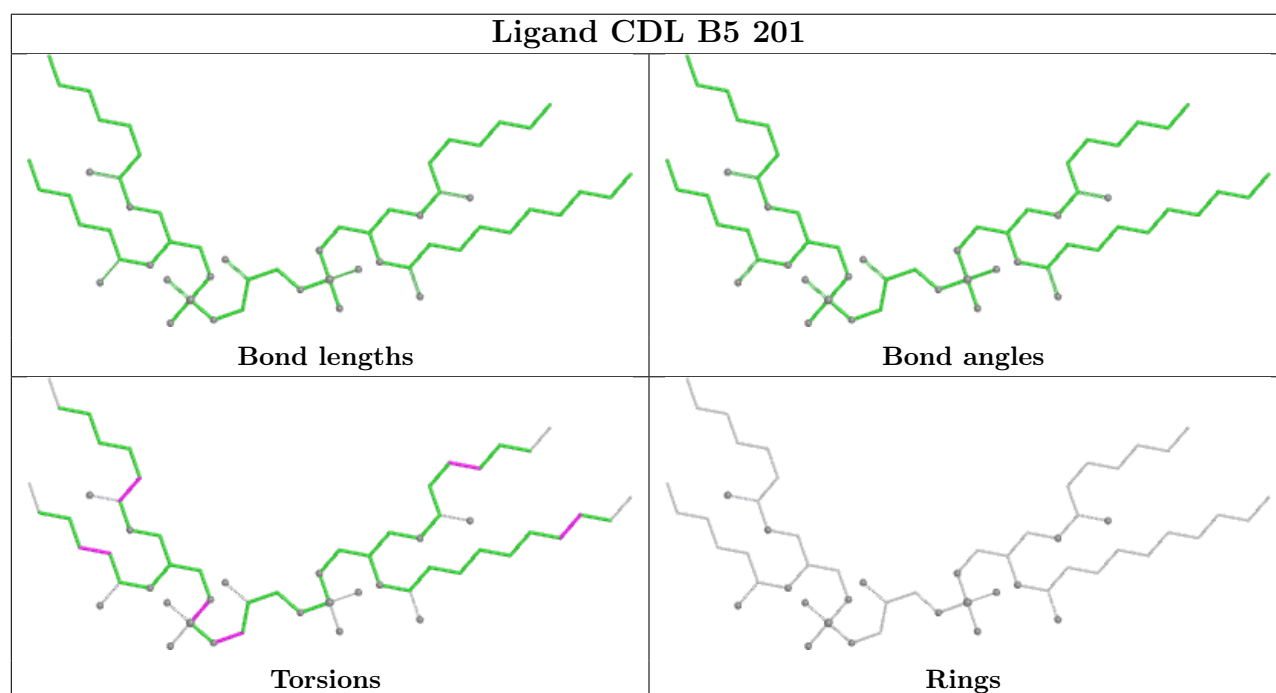
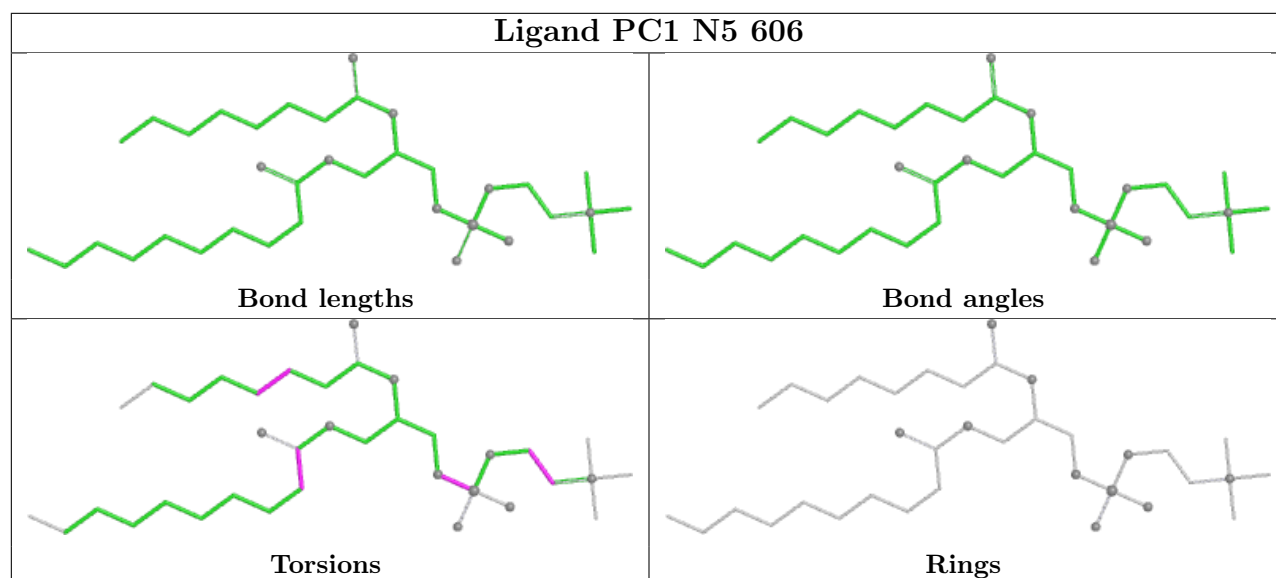
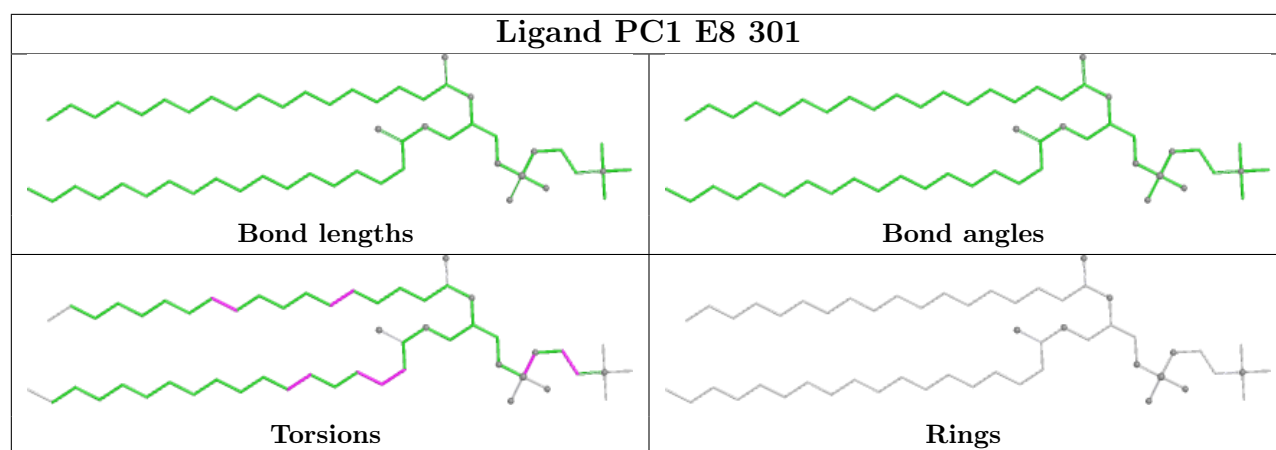
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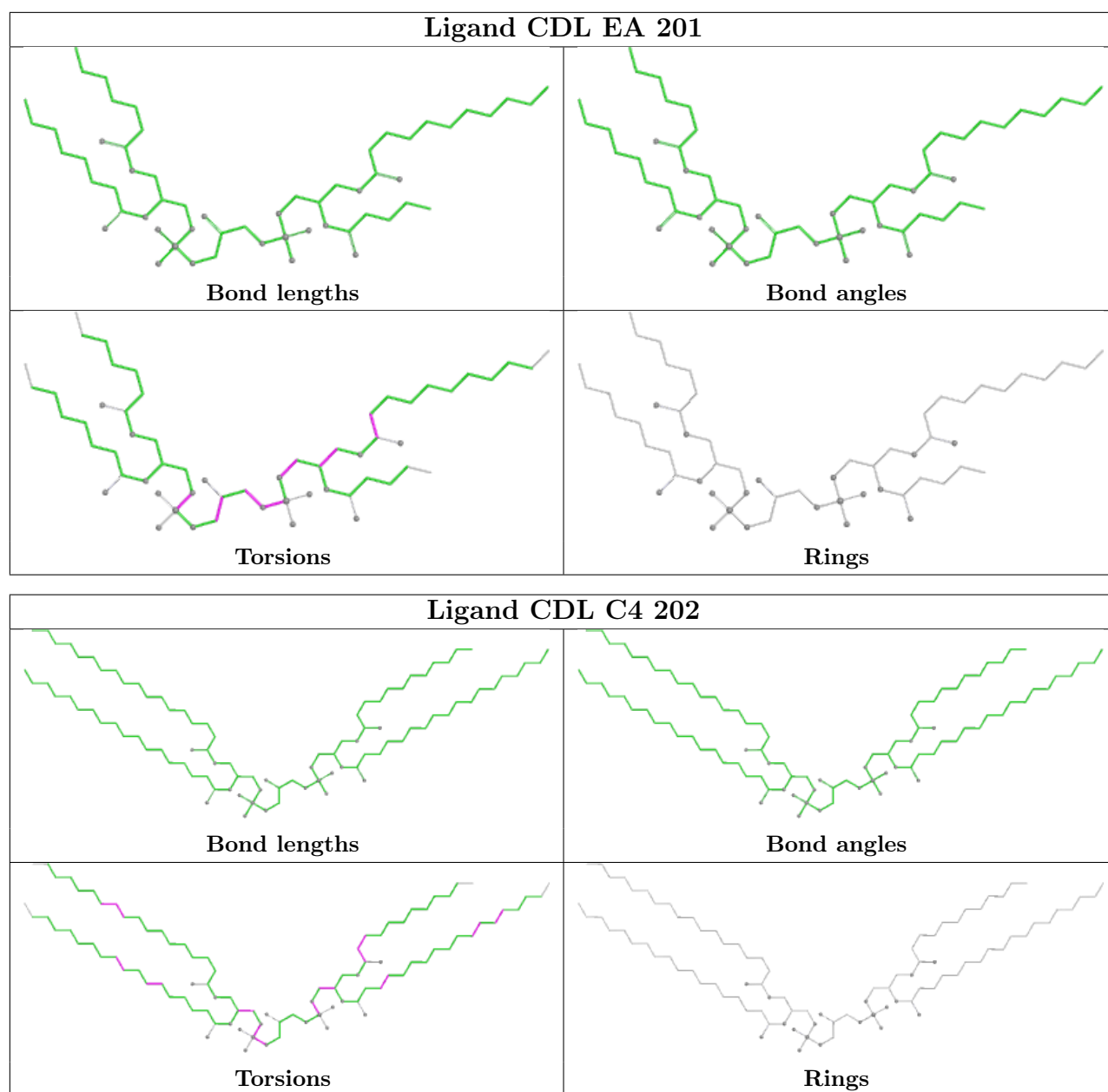
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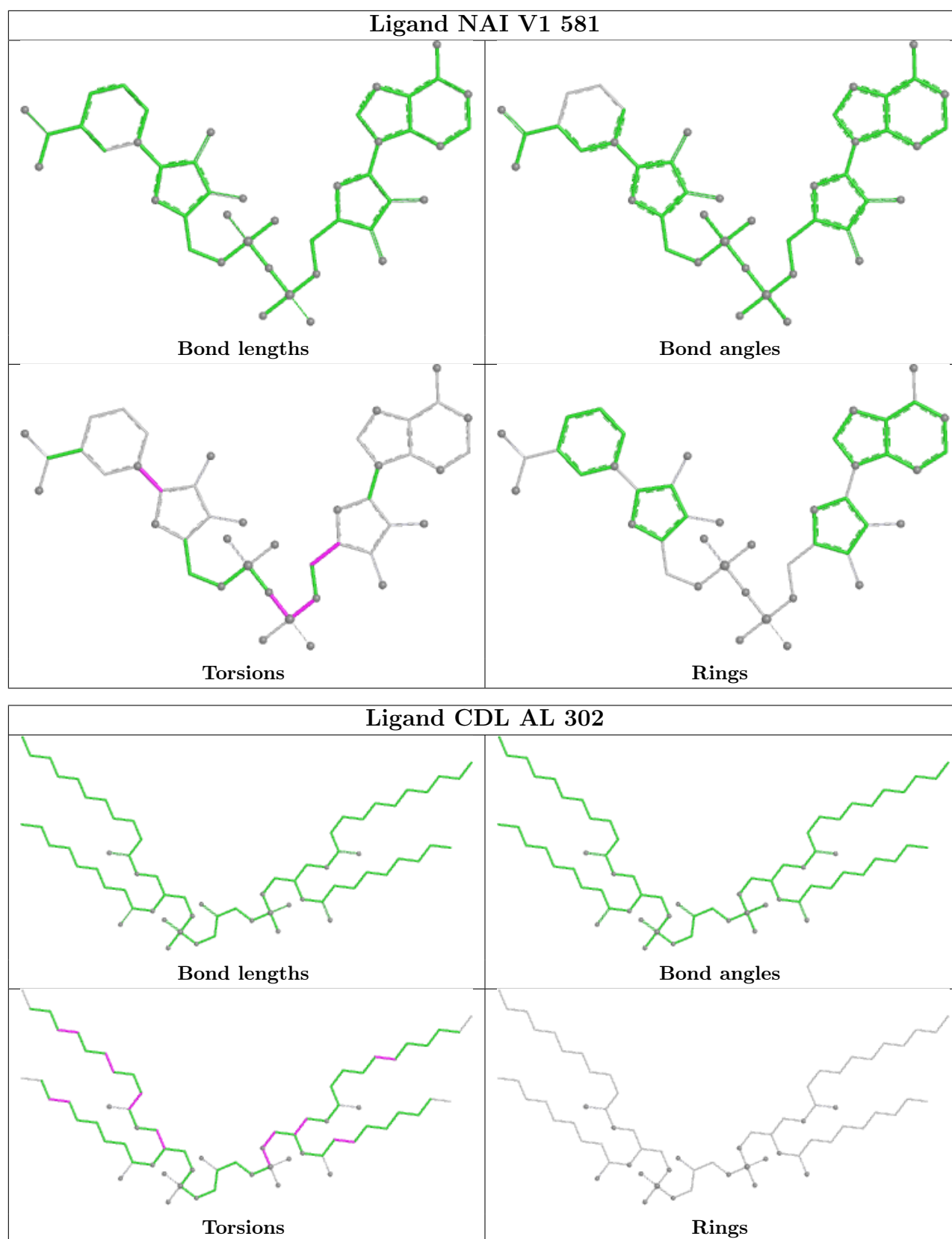
Mol	Chain	Res	Type	Clashes	Symm-Clashes
64	A3	201	CDL	1	0
64	B3	102	CDL	3	0
64	AL	304	CDL	2	0
63	N4	502	PC1	1	0
64	C4	204	CDL	1	0
61	S8	297	SF4	2	0
63	ED	201	PC1	2	0
66	AC	201	ZMP	4	0
64	AM	217	CDL	1	0
63	N2	301	PC1	1	0

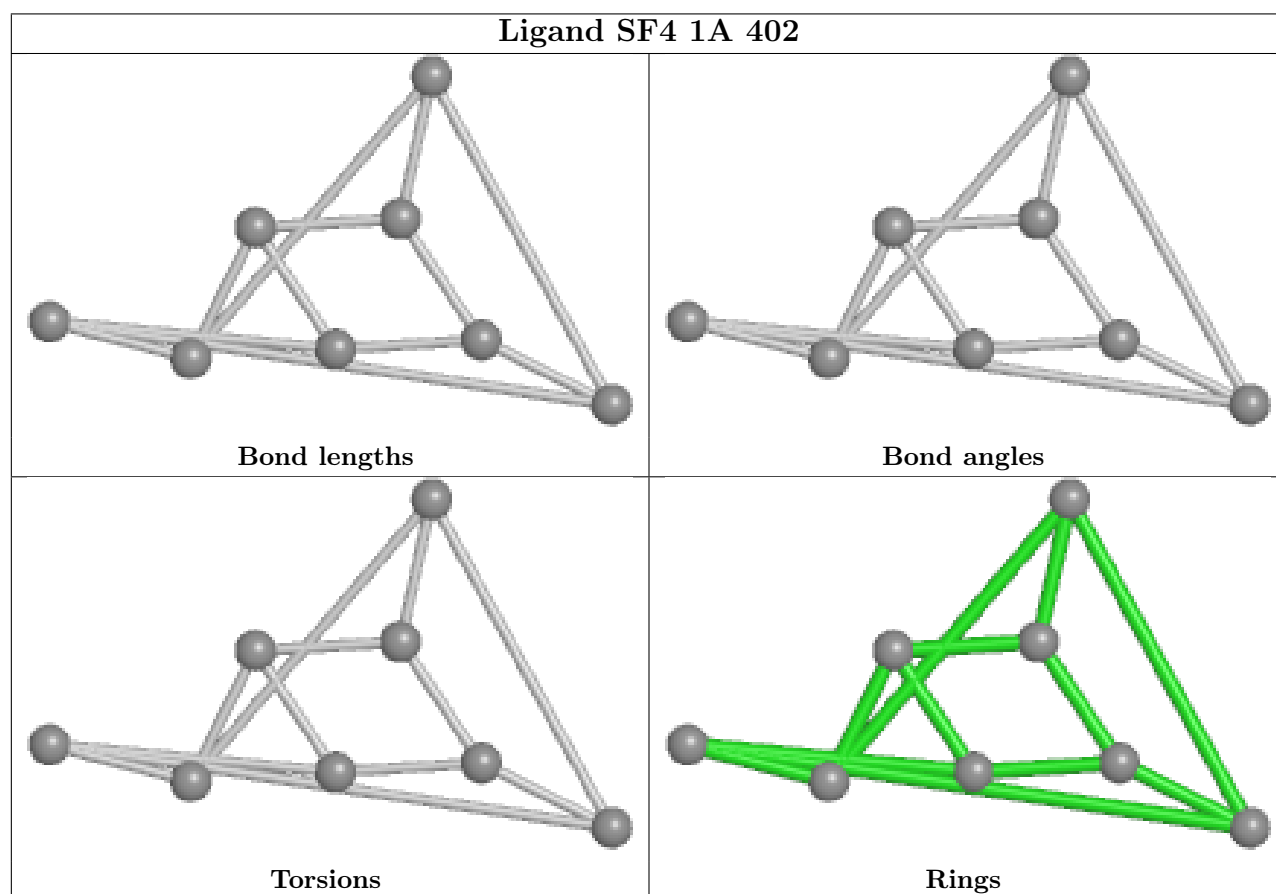
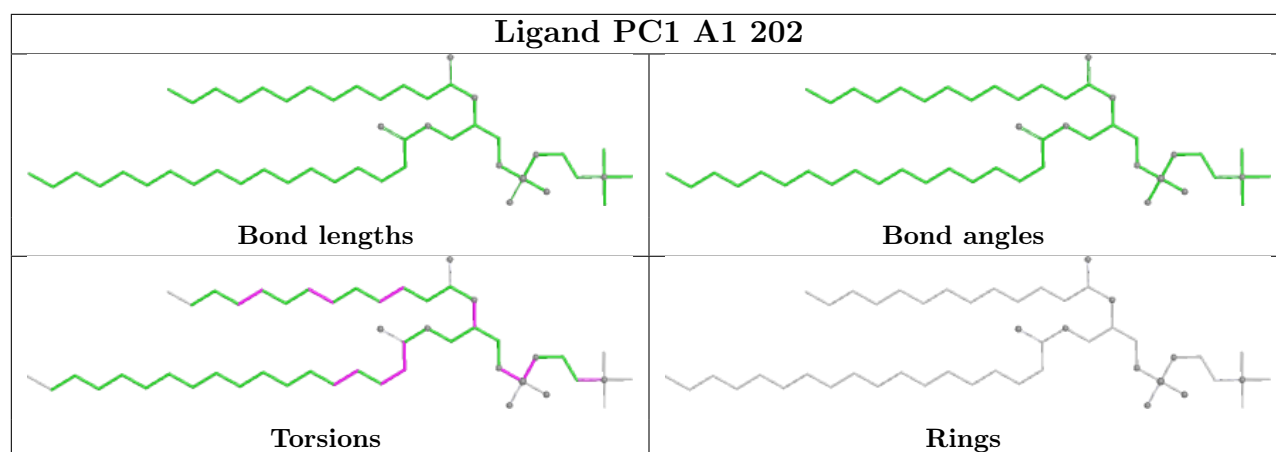
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

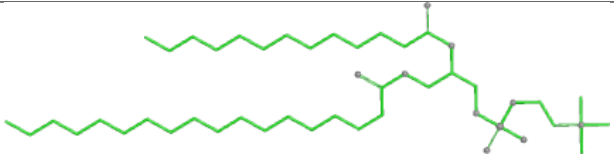
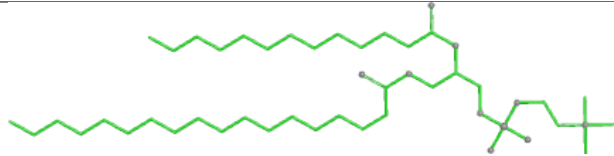
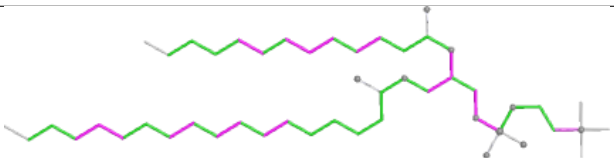
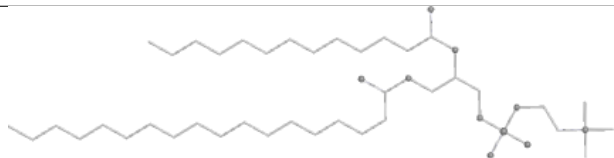


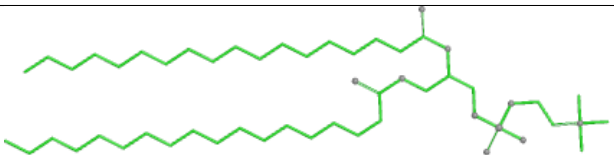
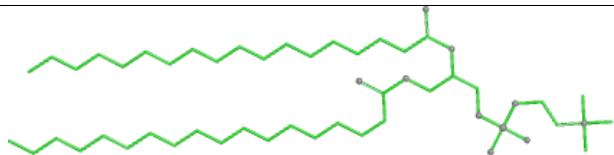
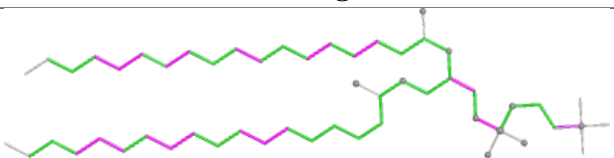
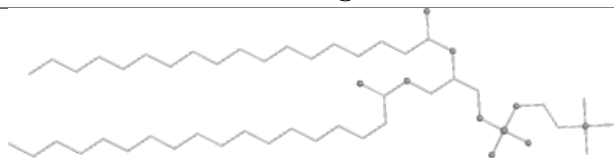


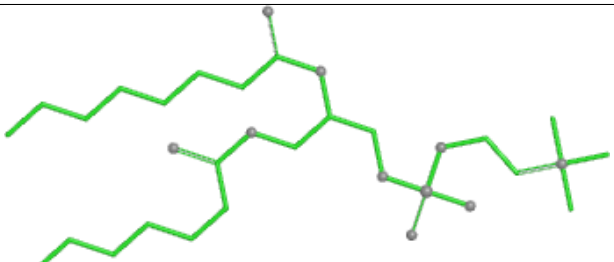
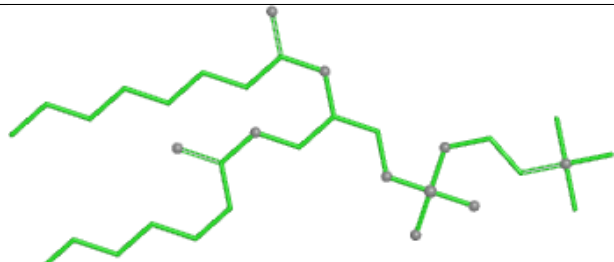
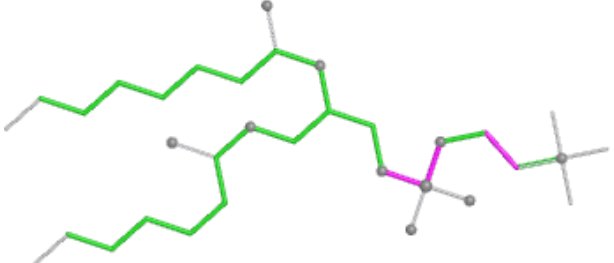
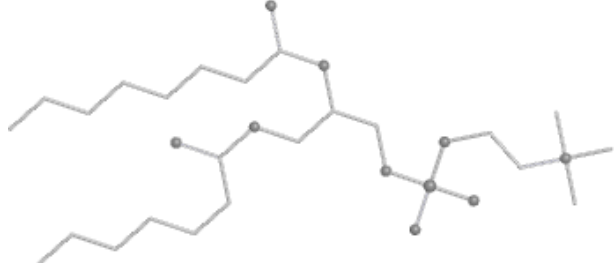


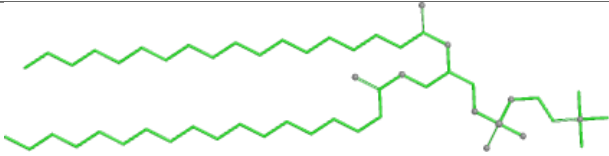
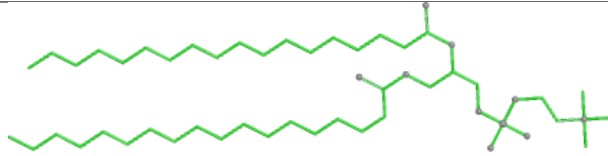
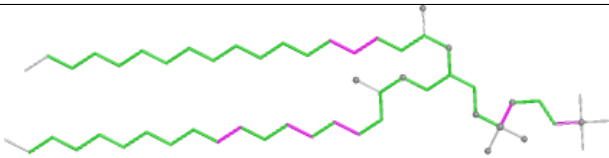
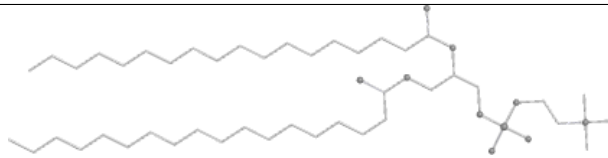
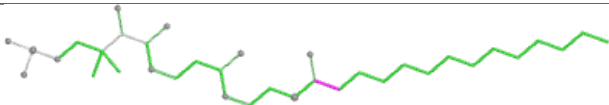
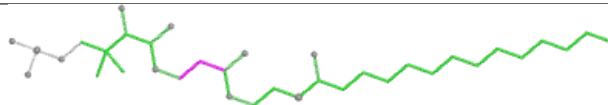
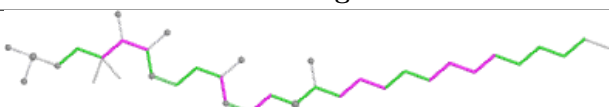
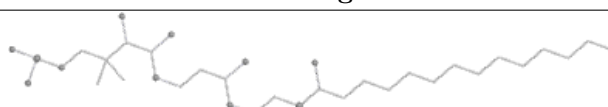
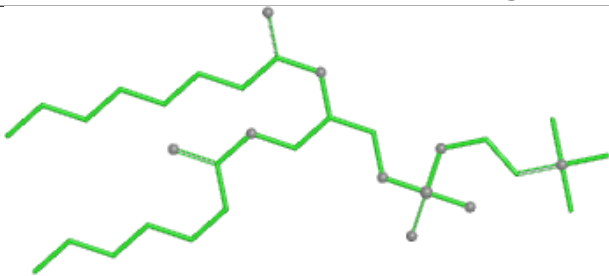
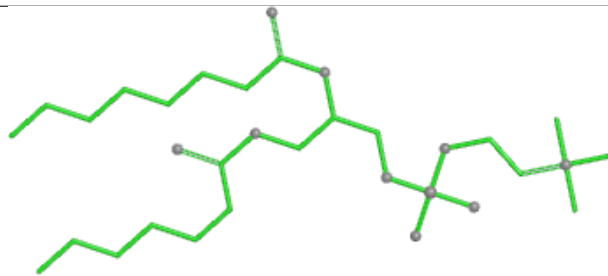
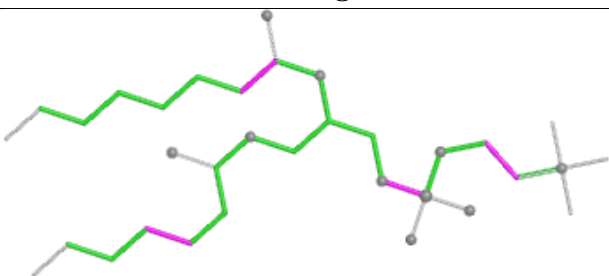
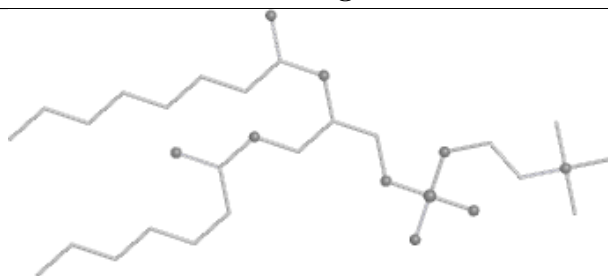


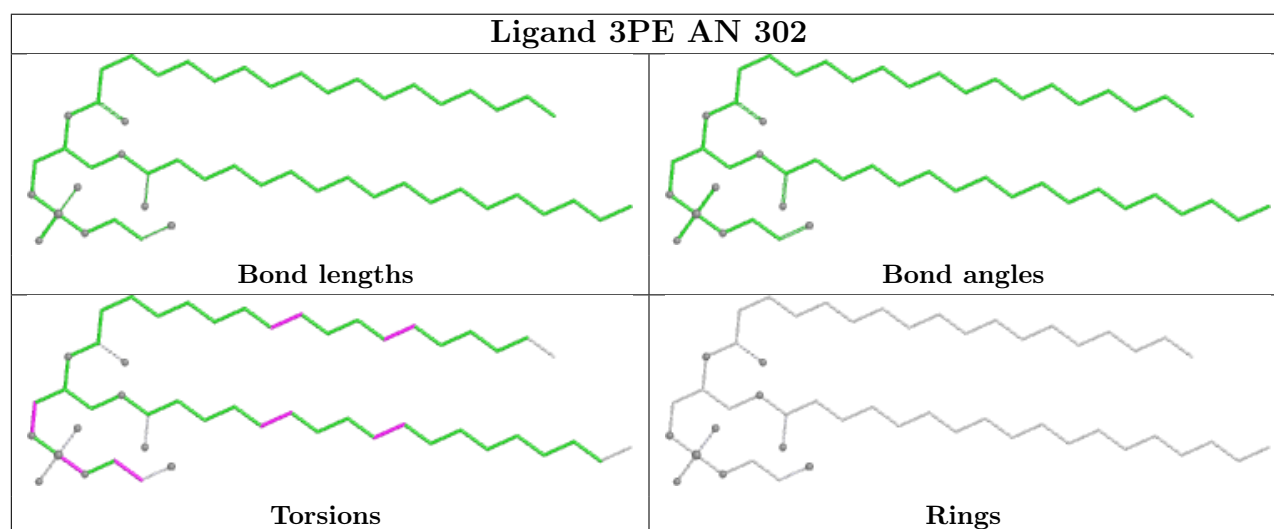
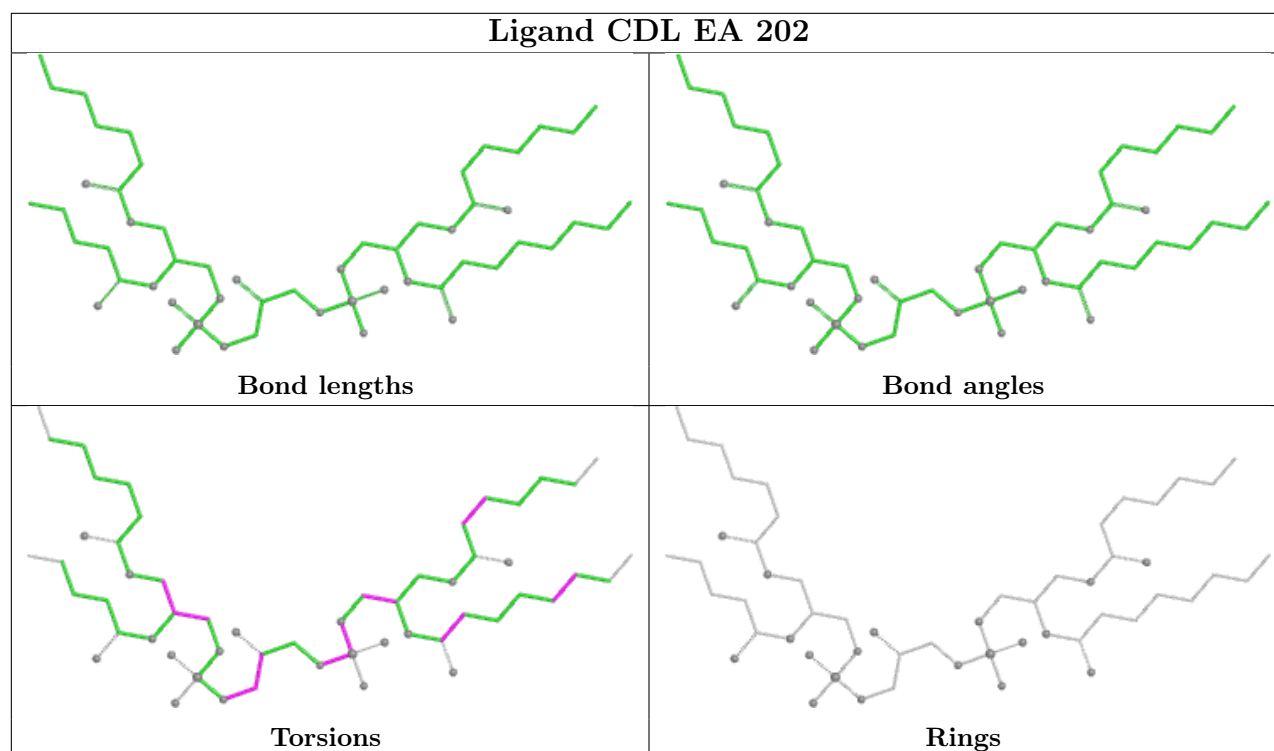
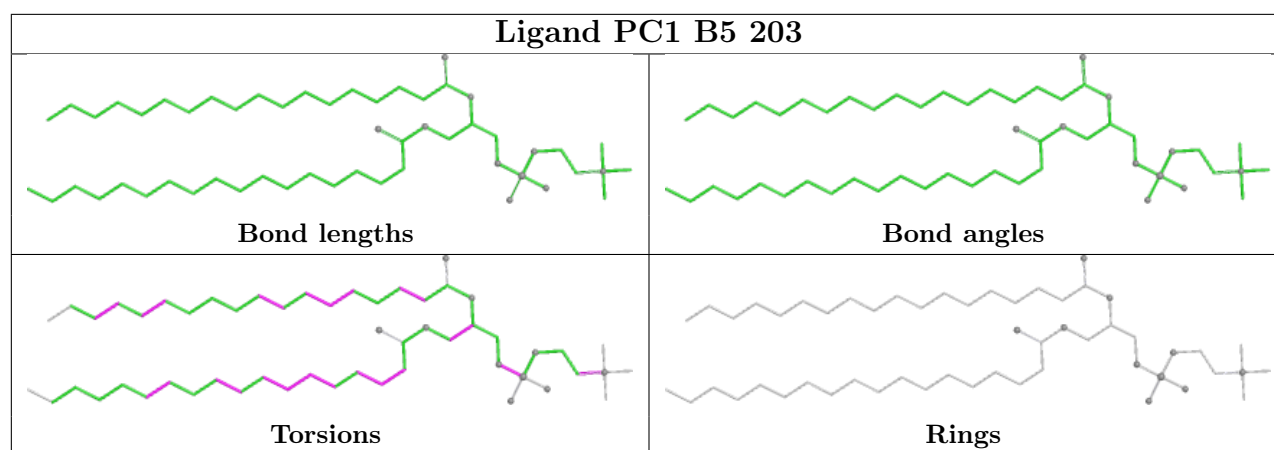


Ligand PC1 AM 218	
	
Bond lengths	Bond angles
	
Torsions	Rings

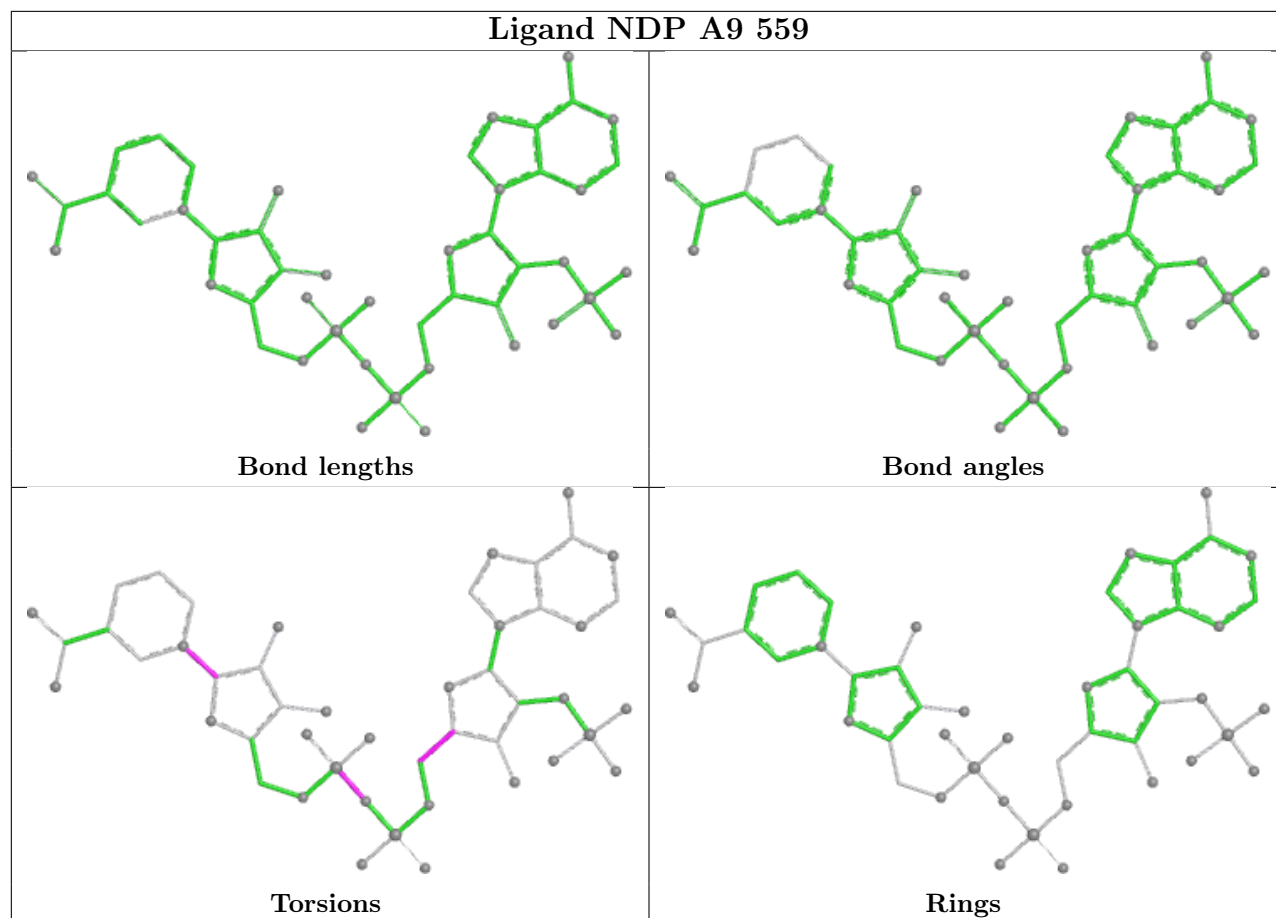
Ligand PC1 E8 302	
	
Bond lengths	Bond angles
	
Torsions	Rings

Ligand PC1 E8 303	
	
Bond lengths	Bond angles
	
Torsions	Rings

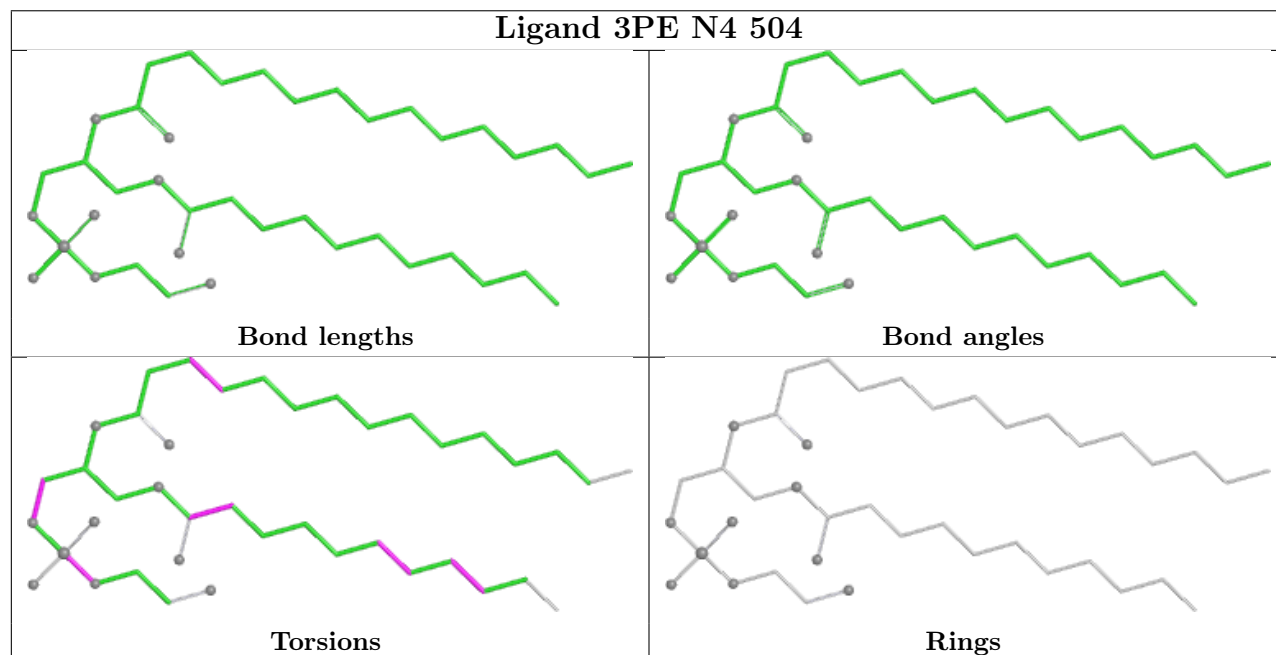
Ligand PC1 N5 601	
 Bond lengths	 Bond angles
 Torsions	 Rings
Ligand ZMP AB 150	
 Bond lengths	 Bond angles
 Torsions	 Rings
Ligand PC1 A9 561	
 Bond lengths	 Bond angles
 Torsions	 Rings

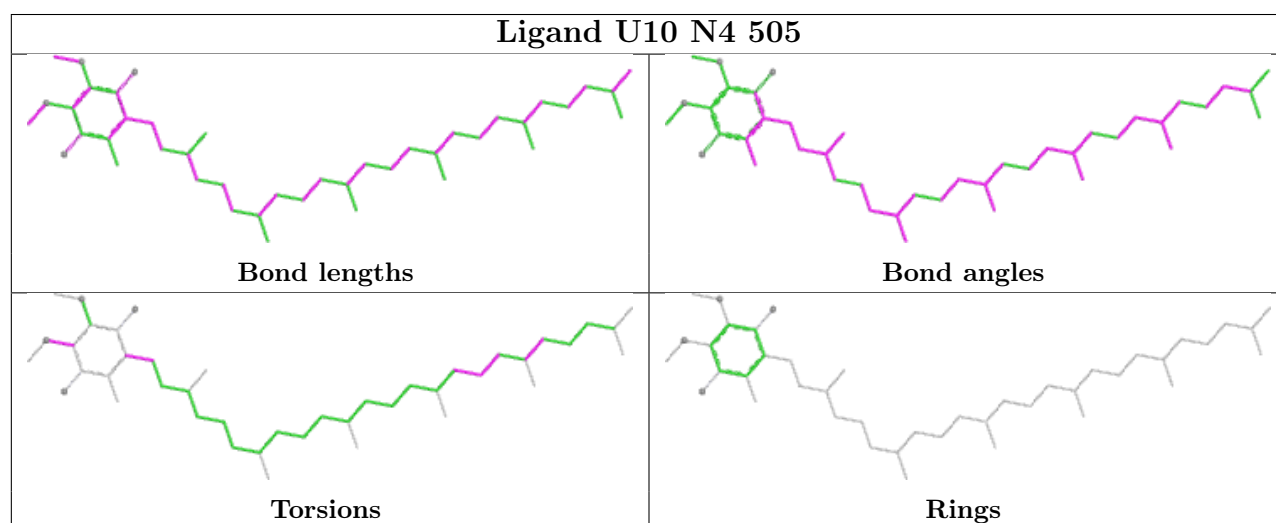
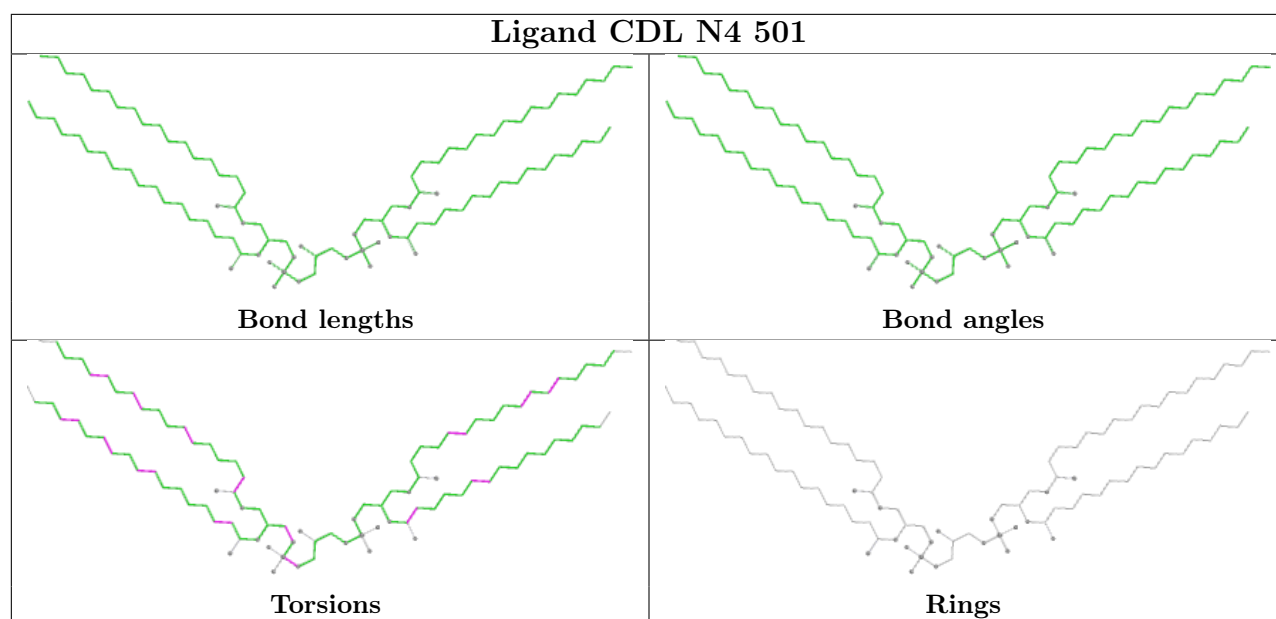


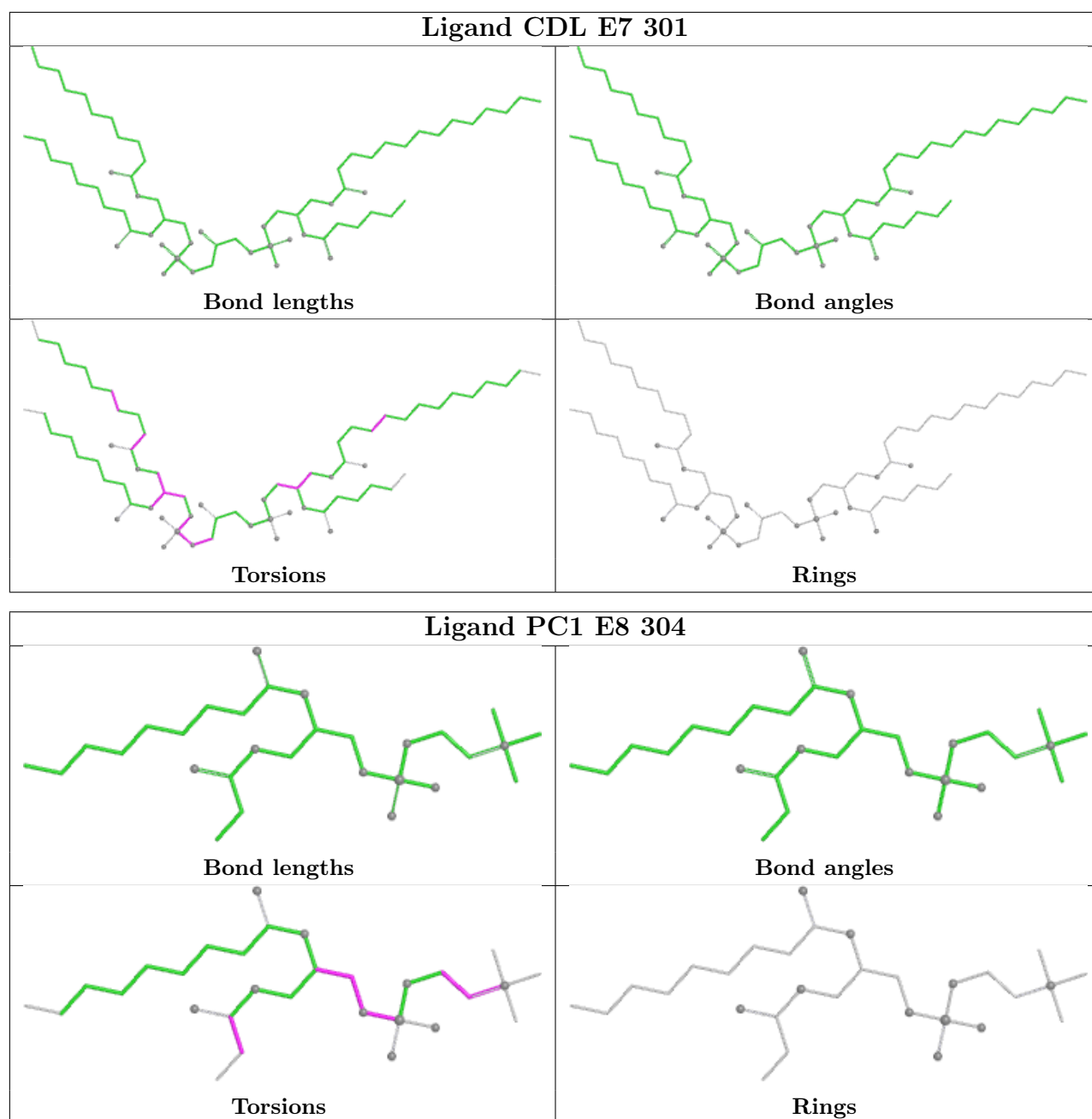
Ligand NDP A9 559

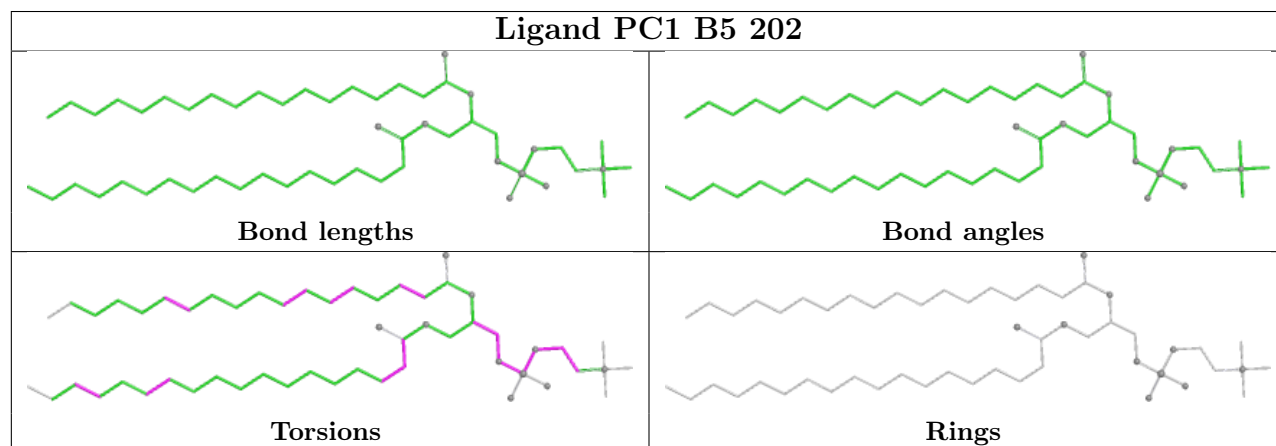
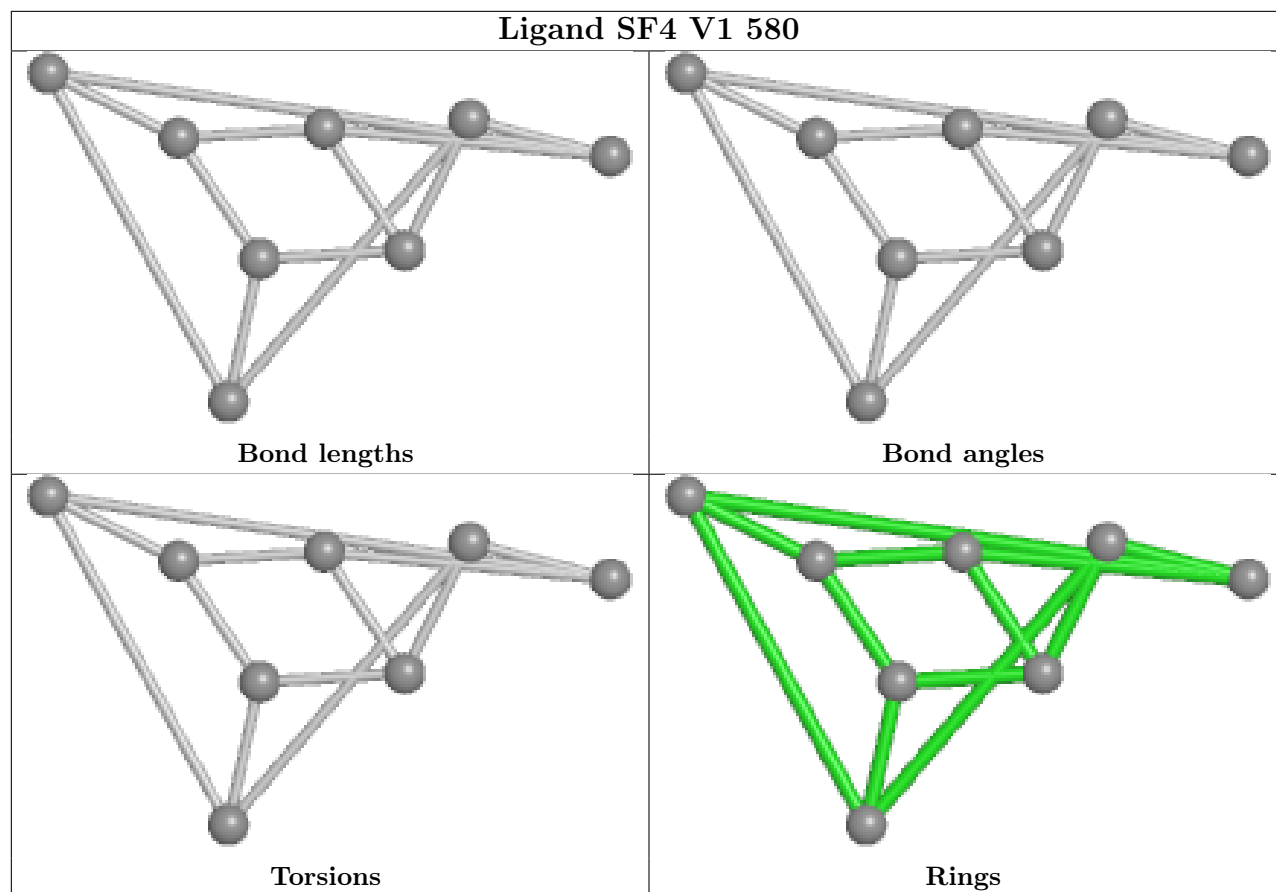


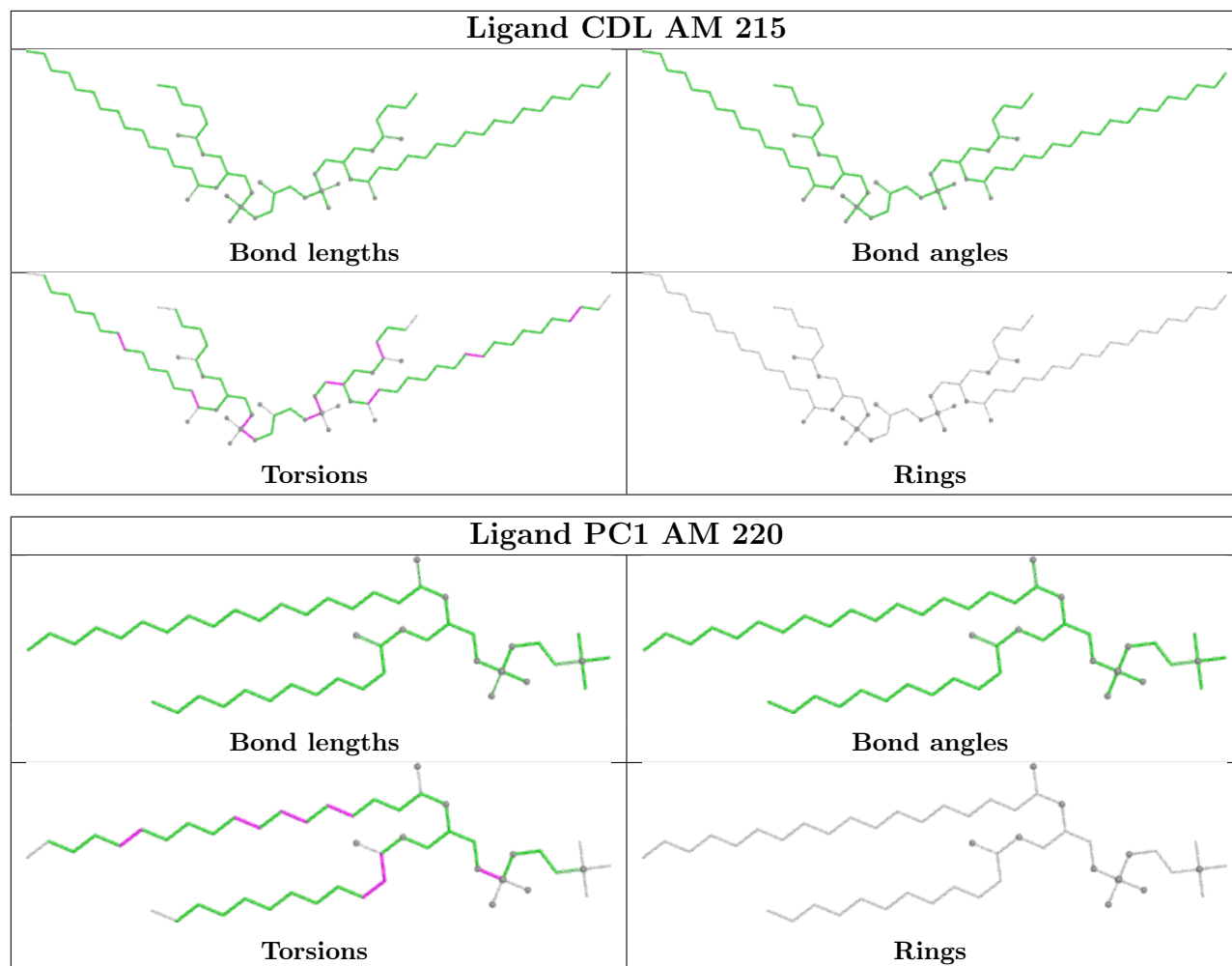
Ligand 3PE N4 504

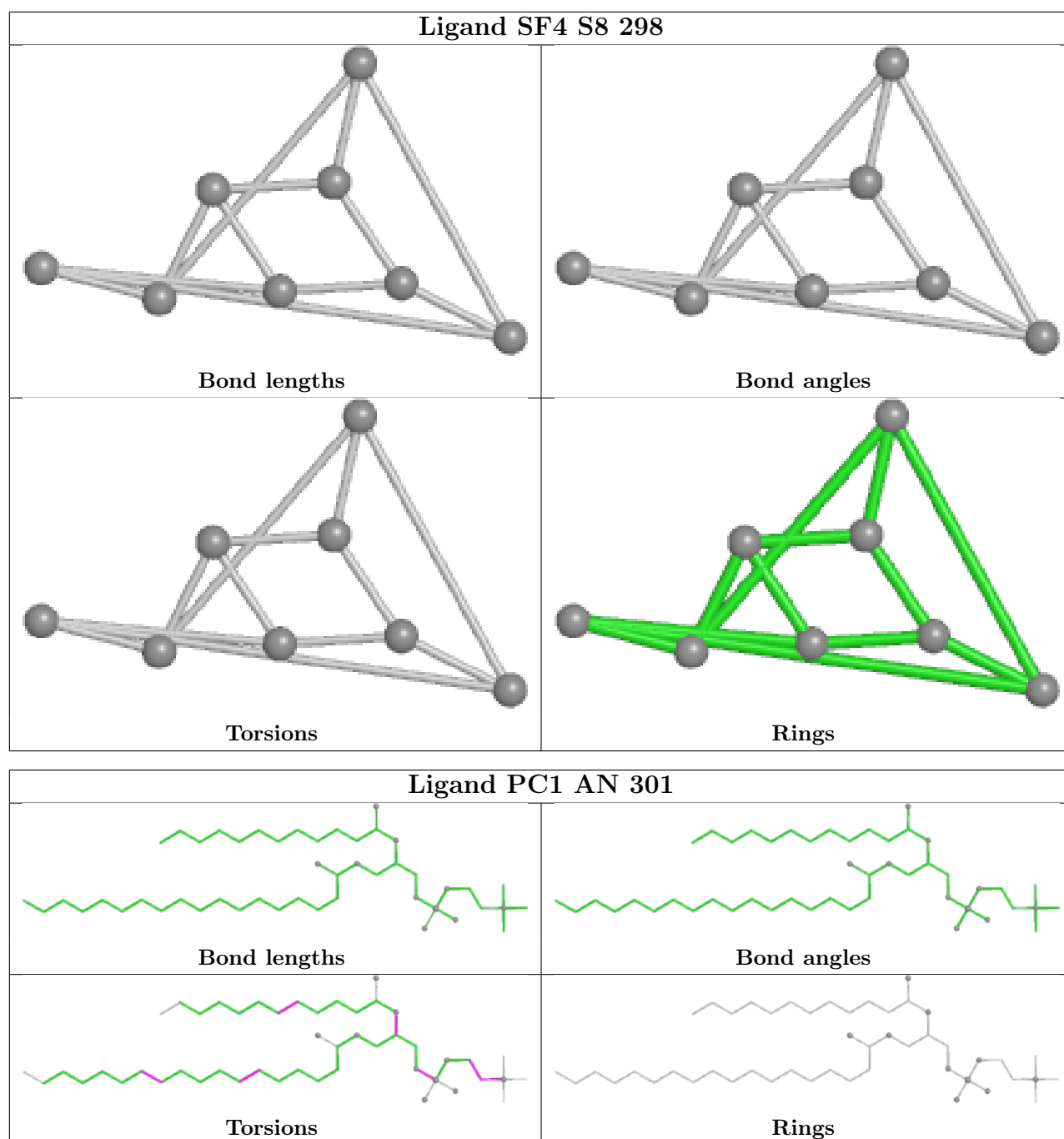


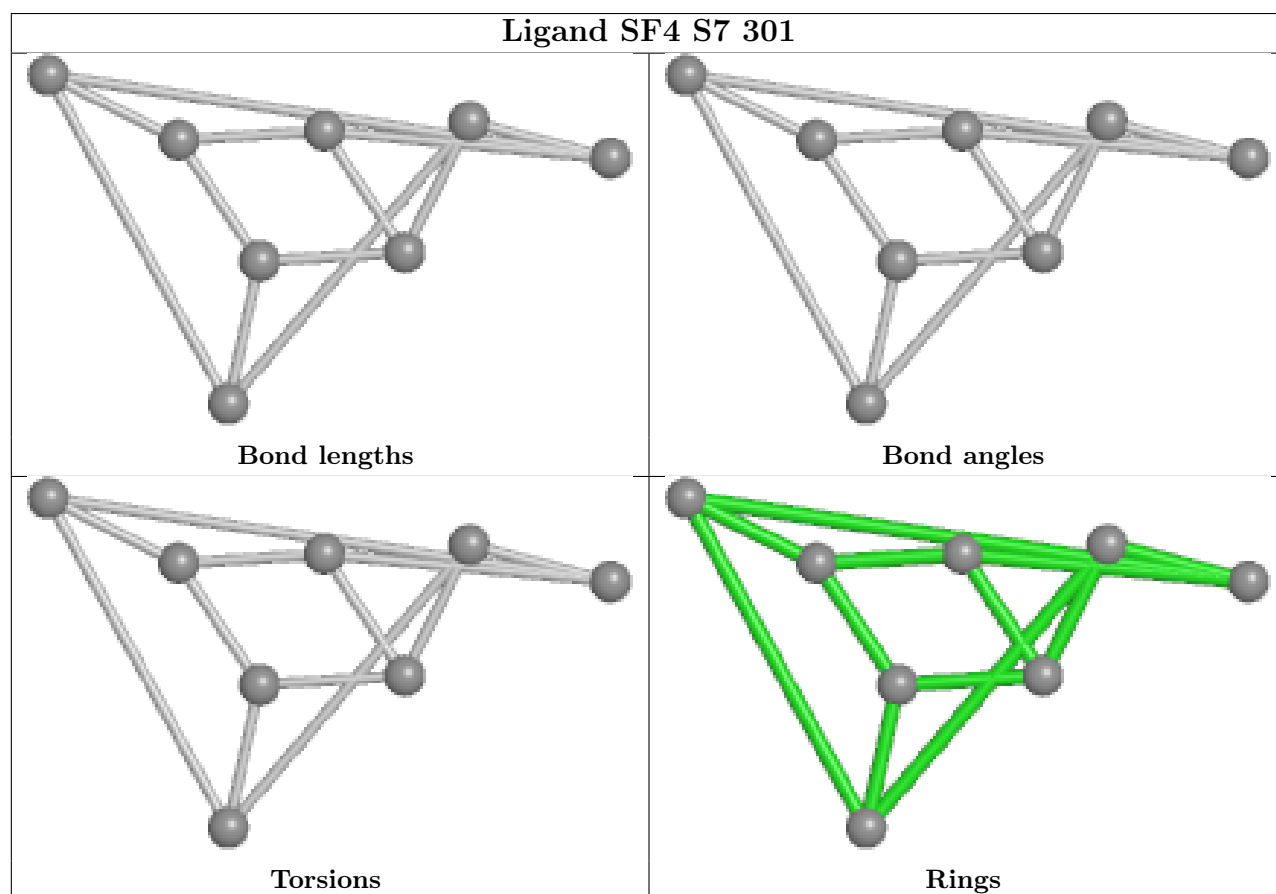
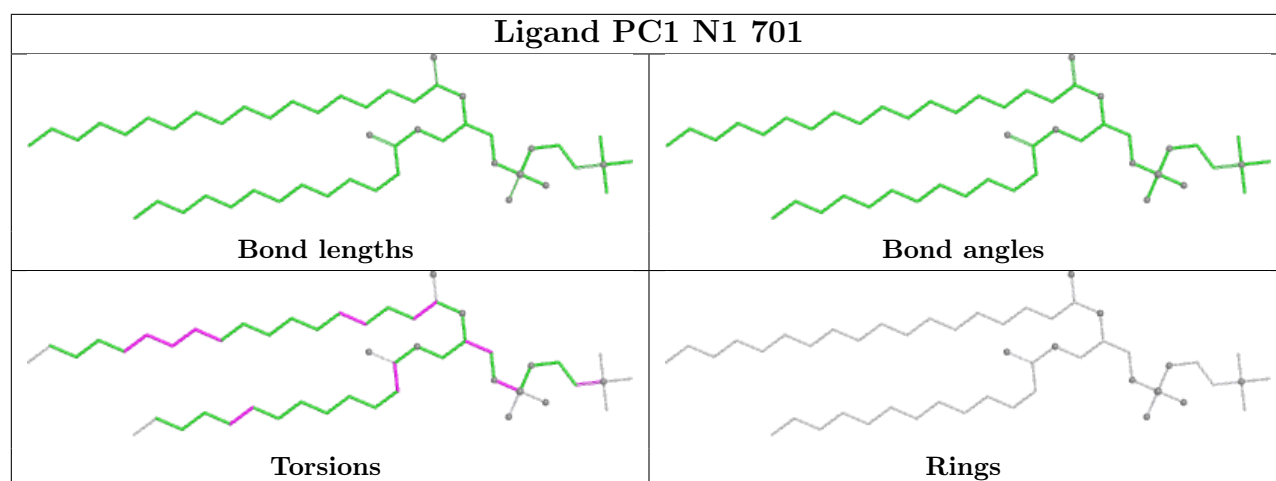


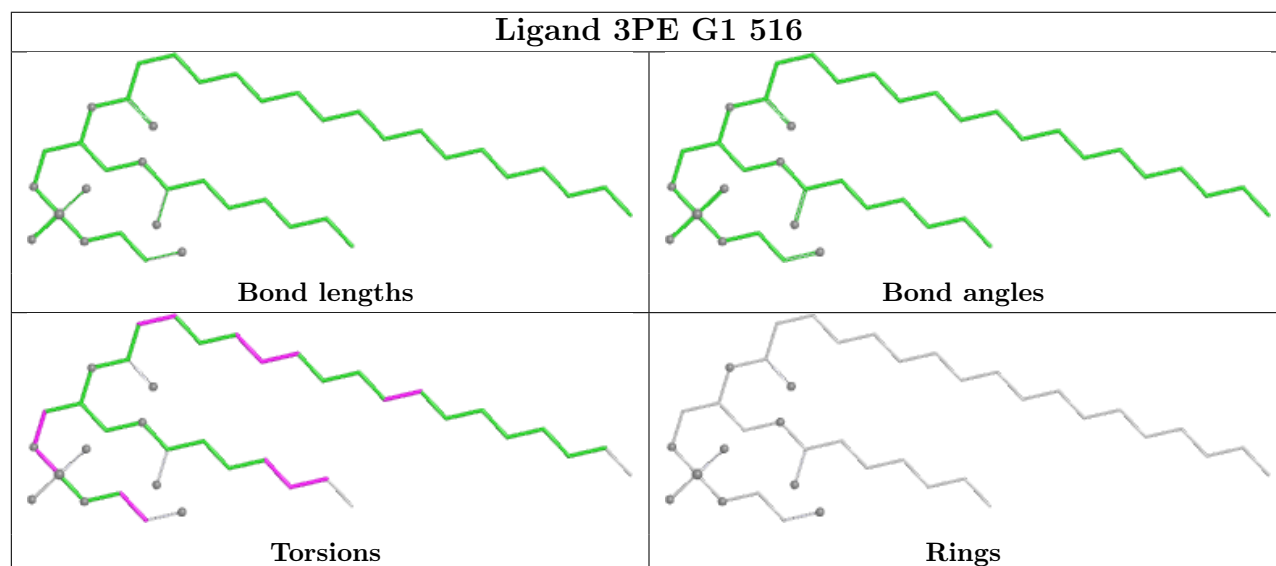
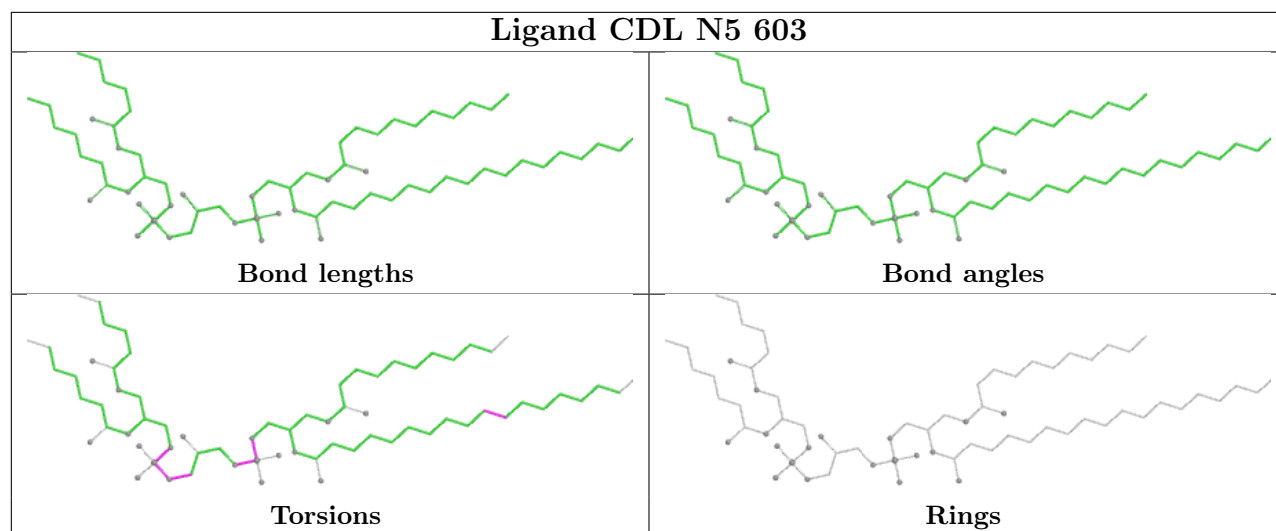
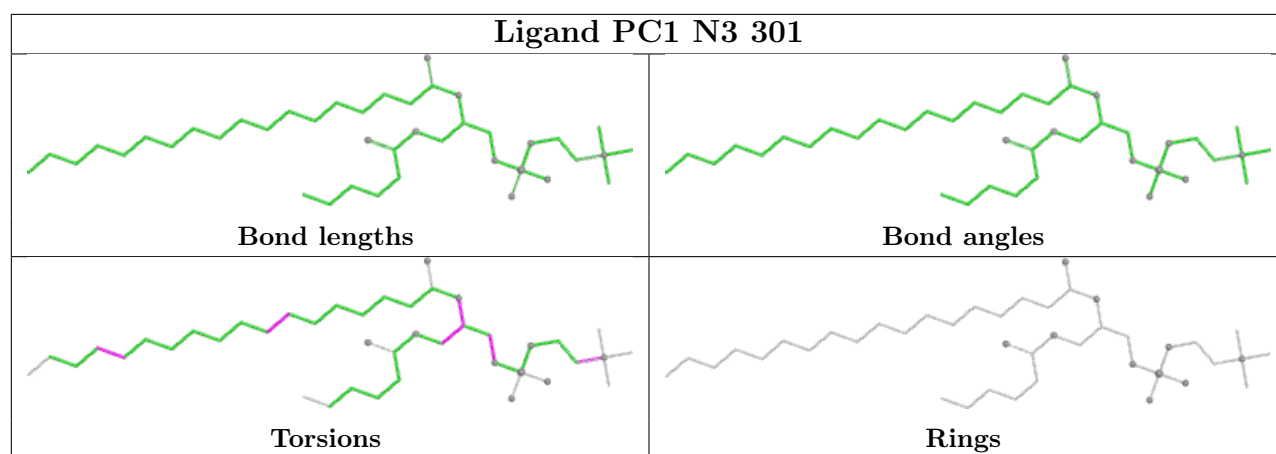




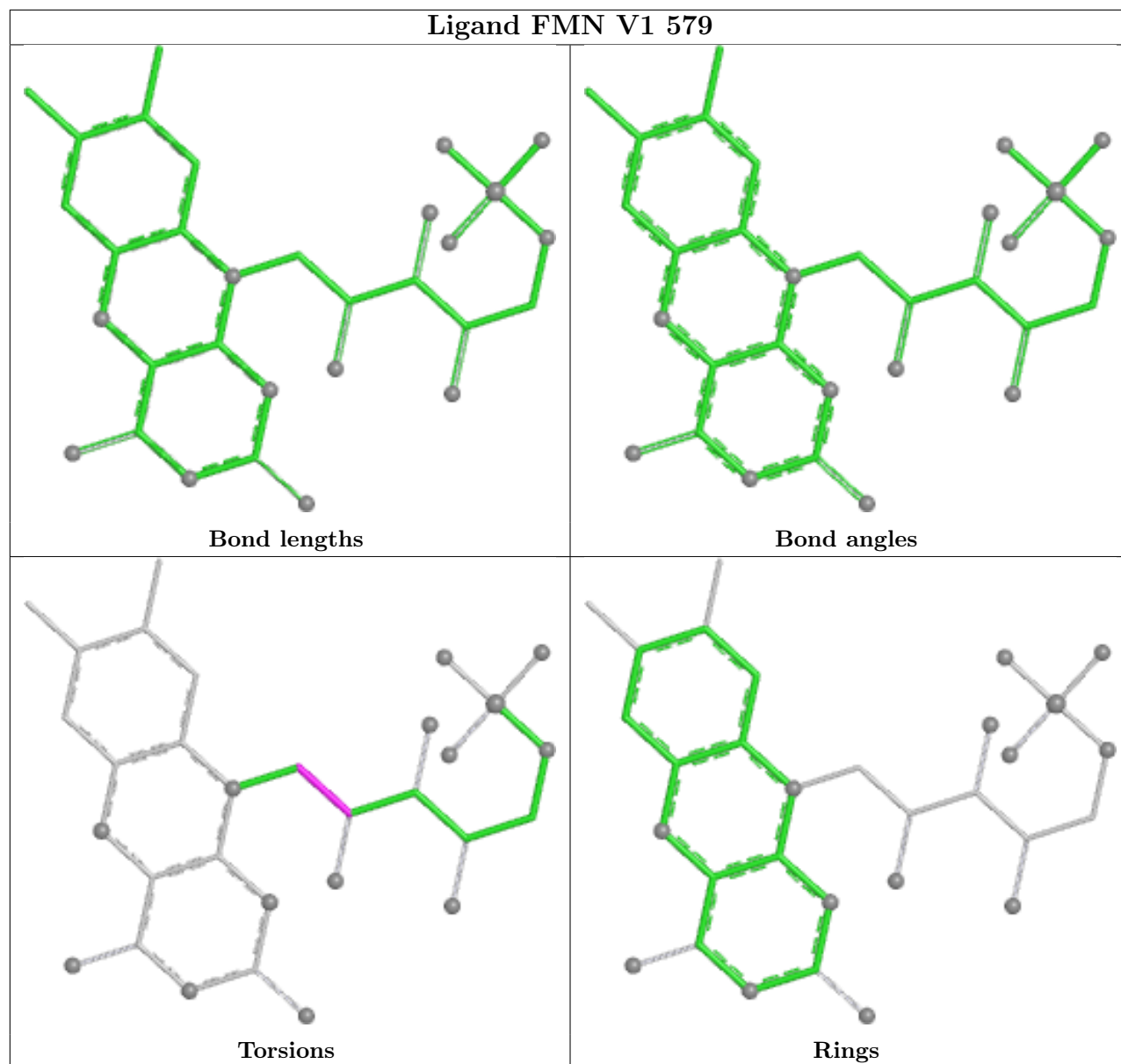


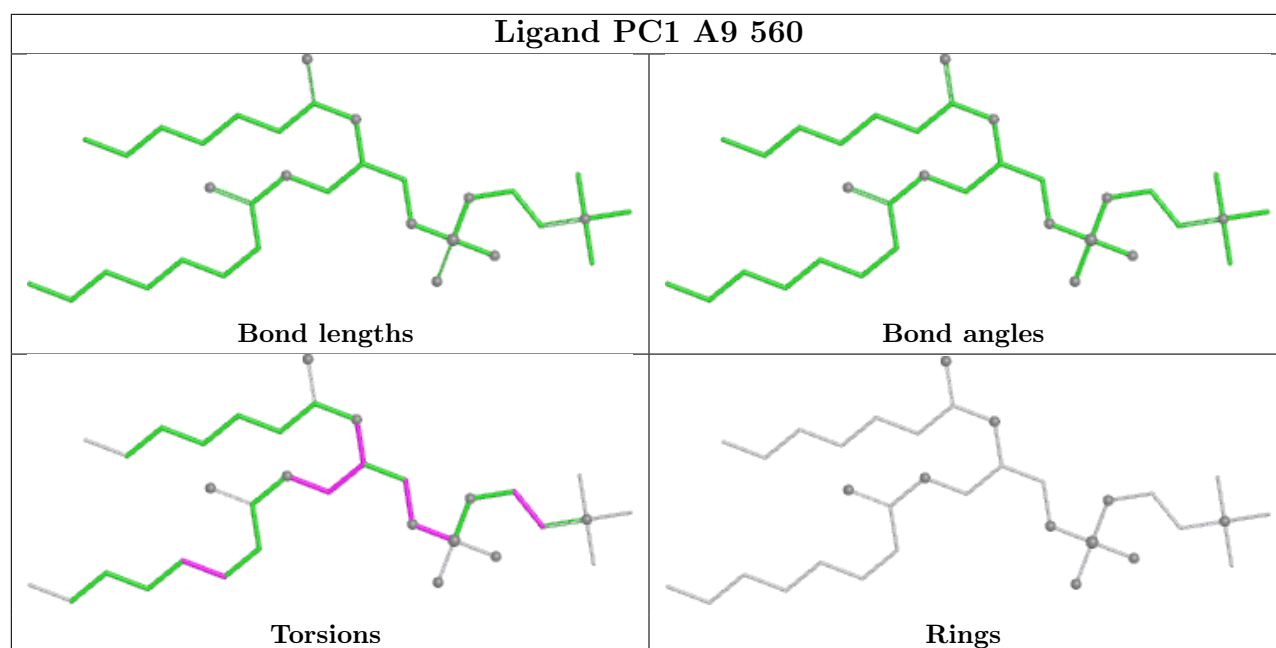
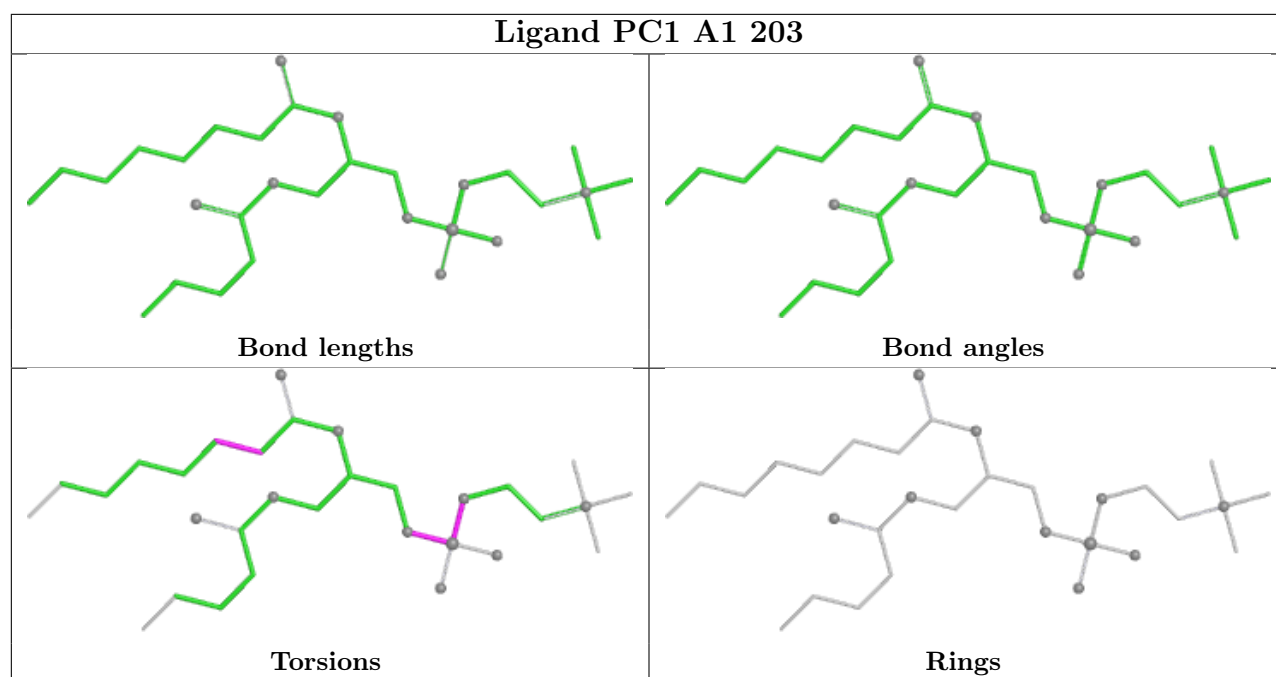


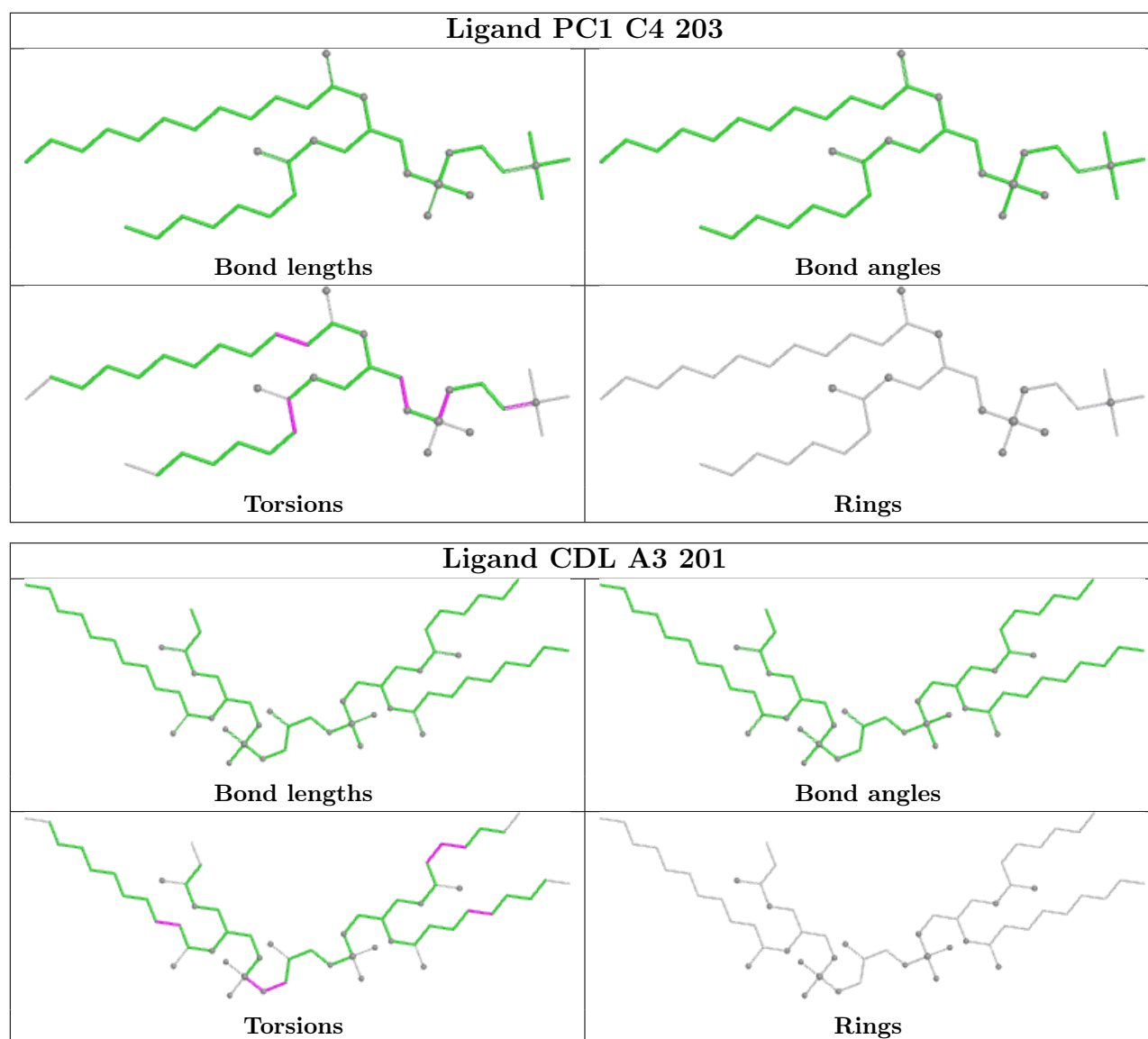


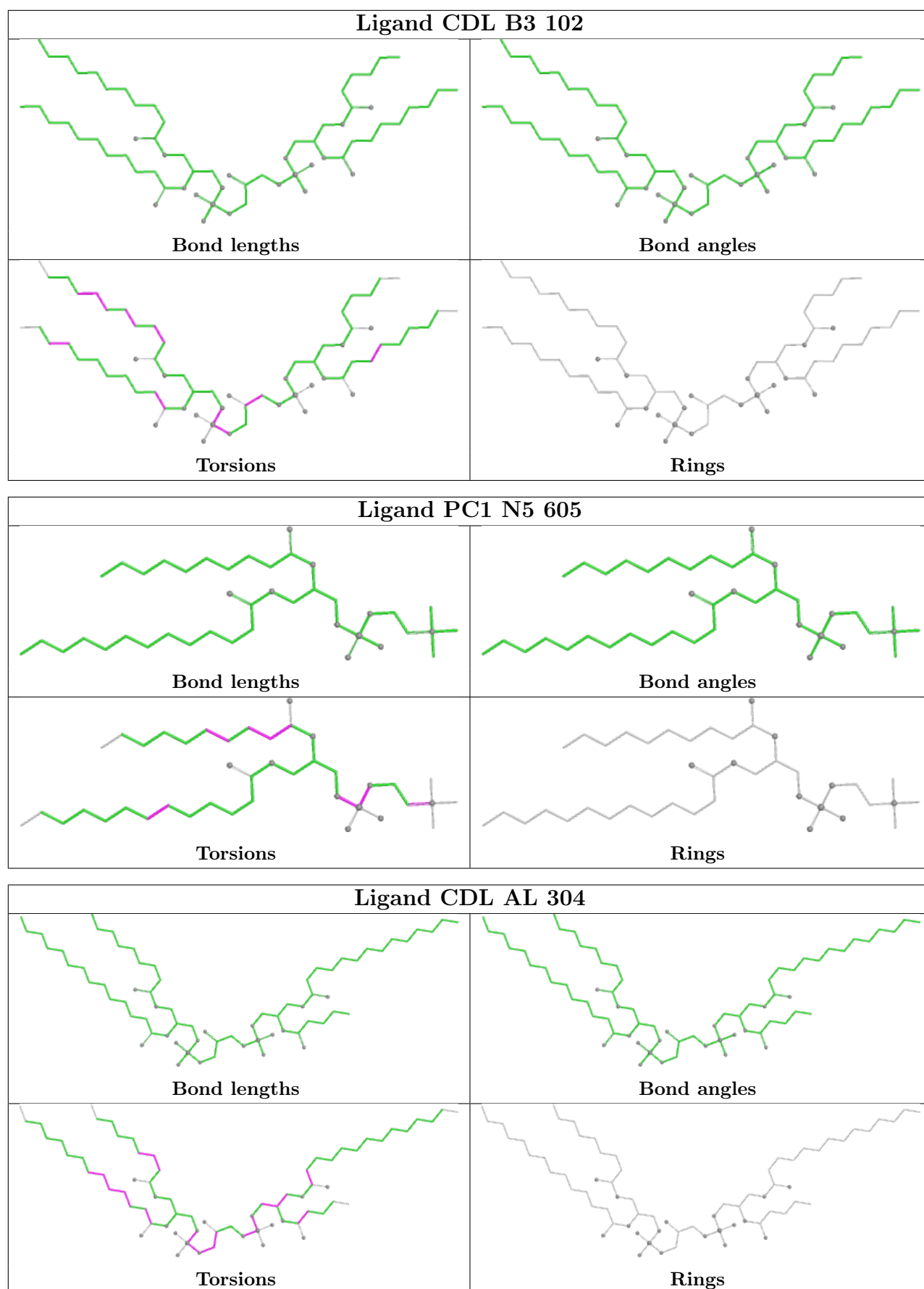


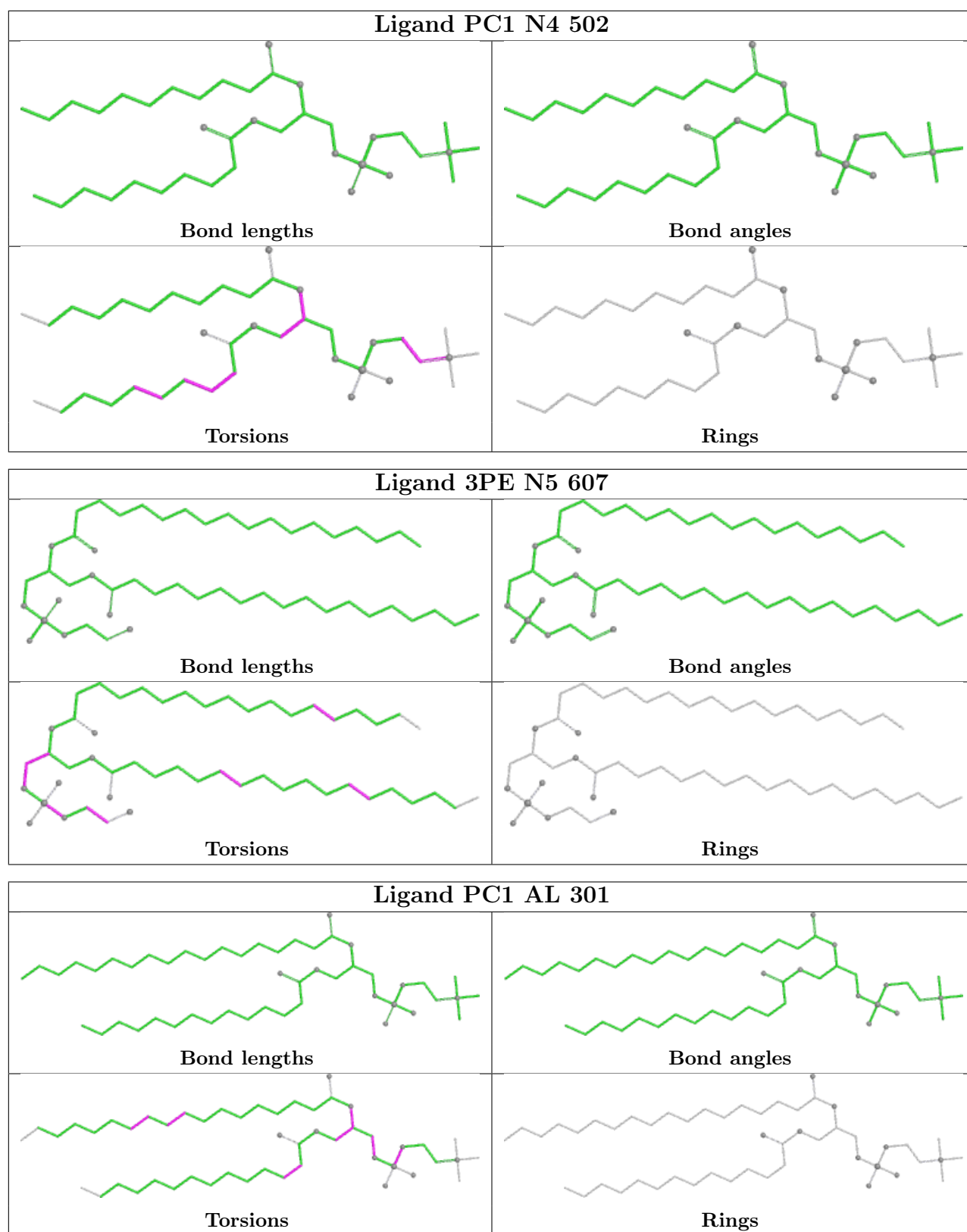
Ligand FMN V1 579

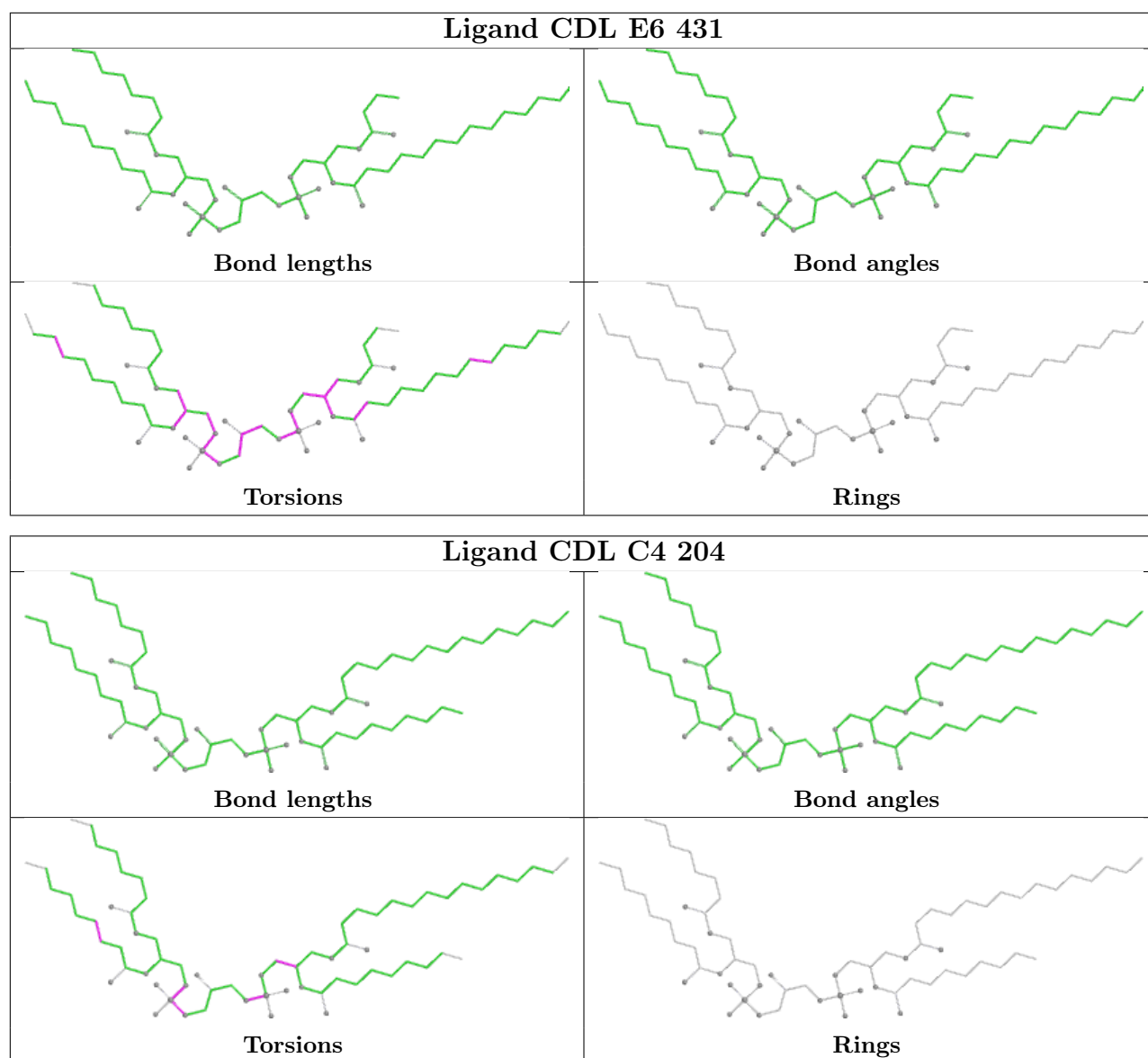


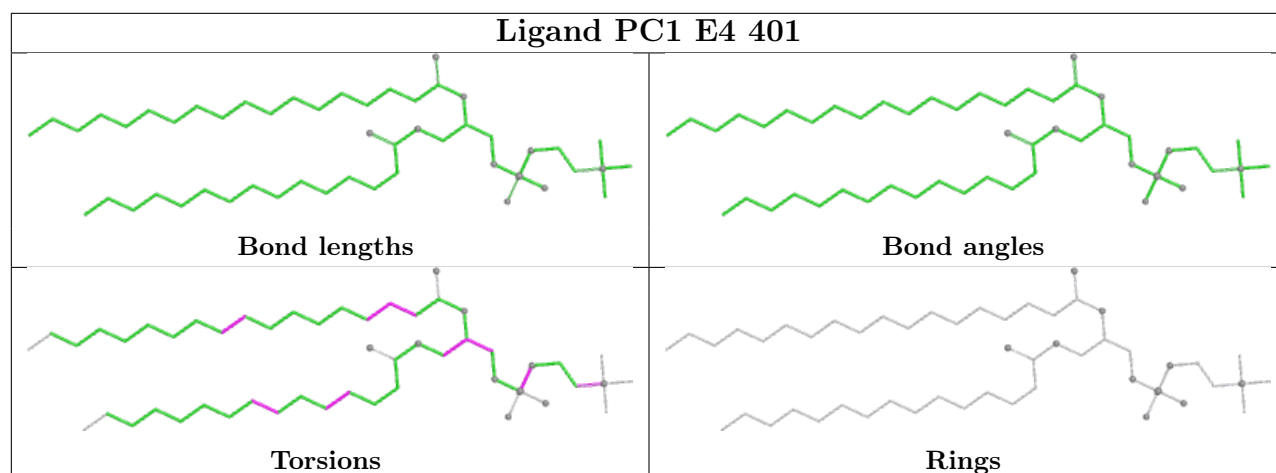
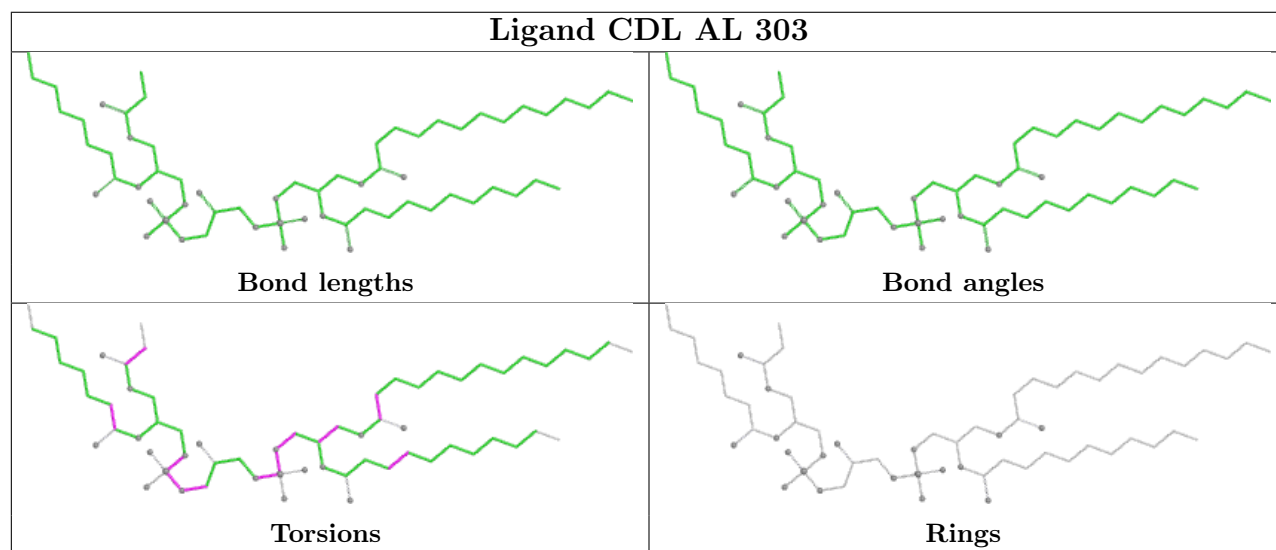
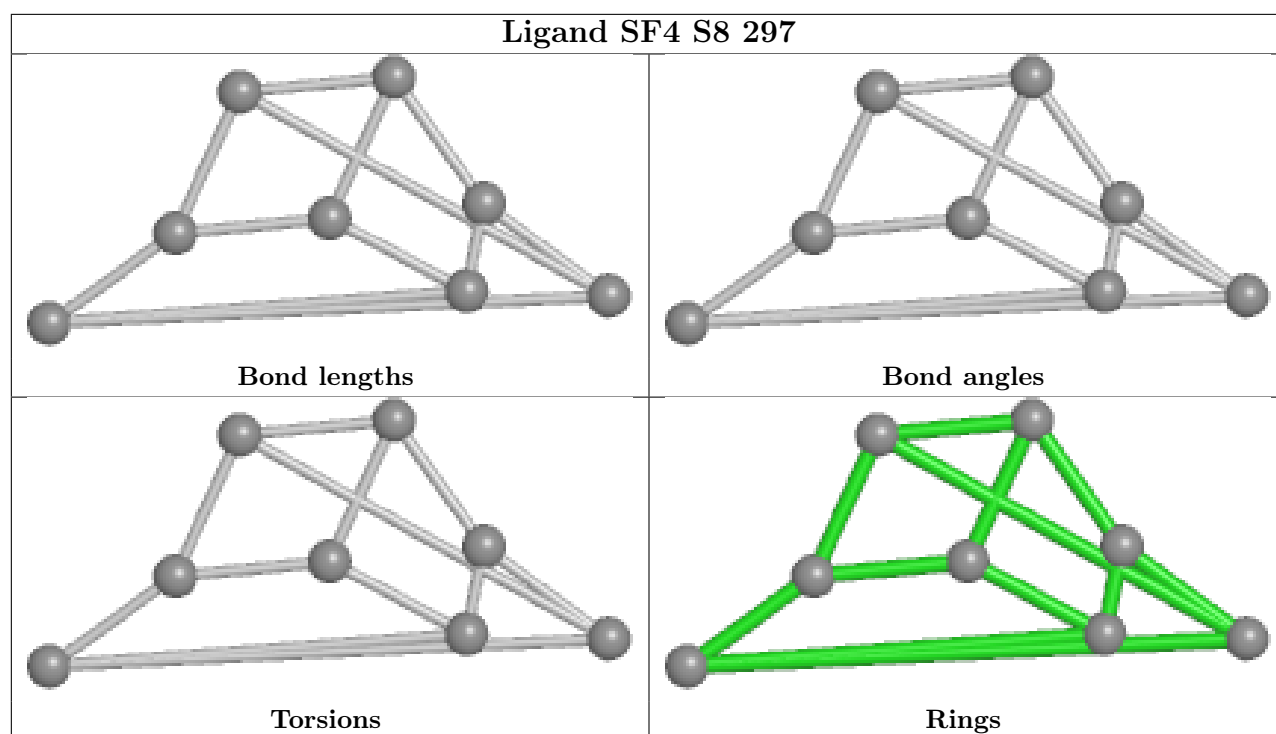


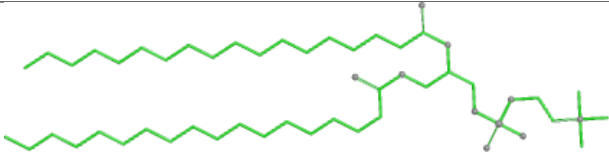
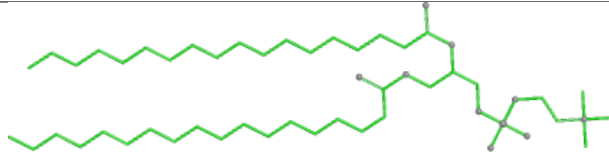
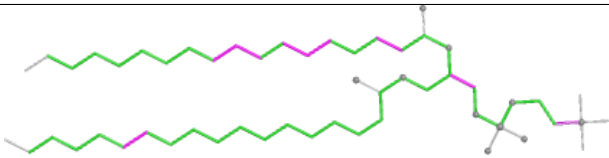
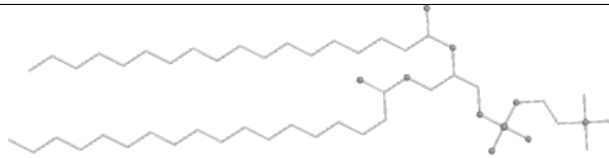
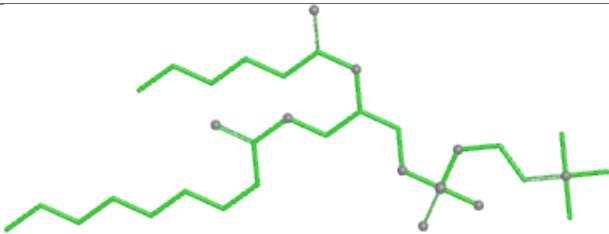
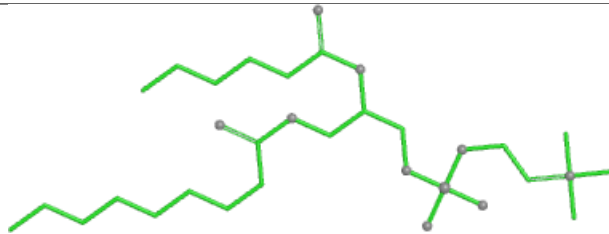
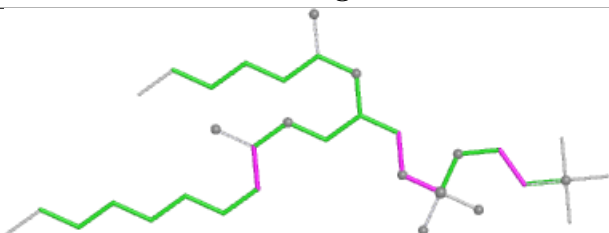
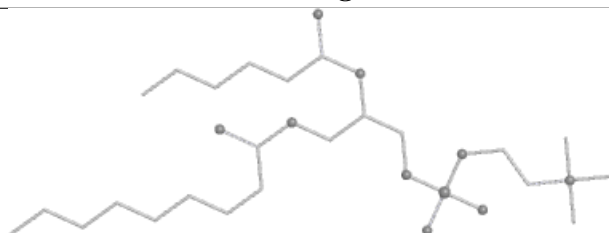
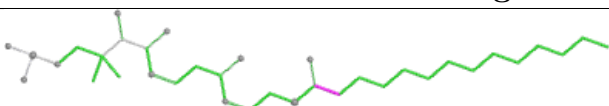
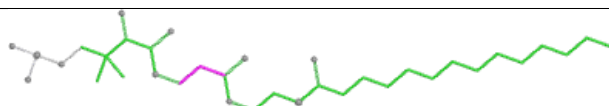
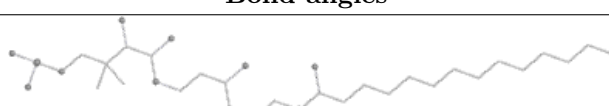


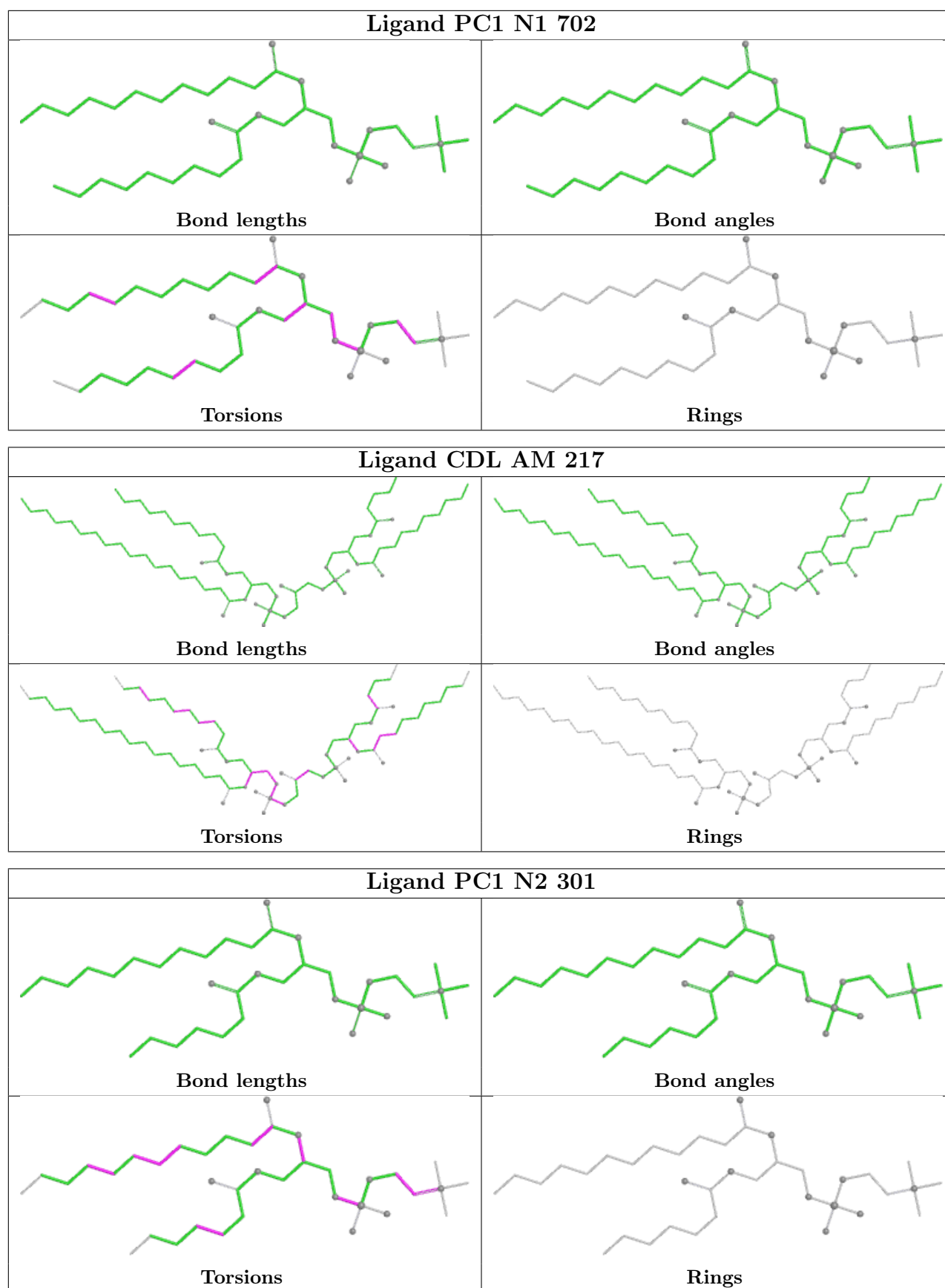


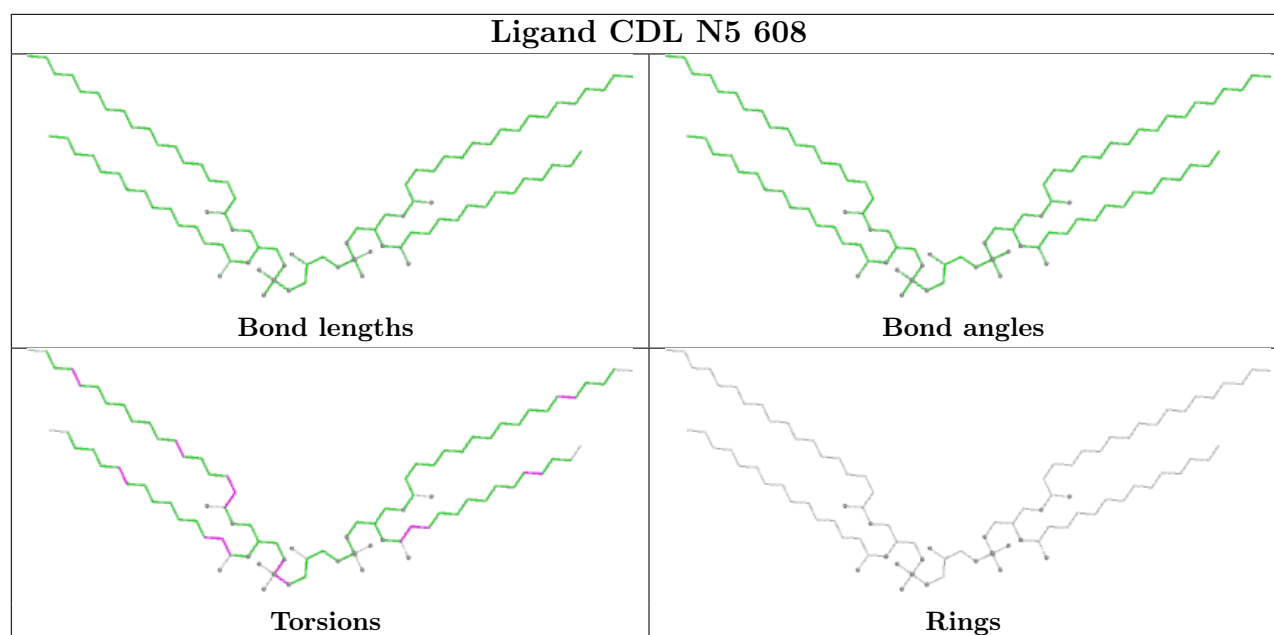






Ligand PC1 ED 201	
 Bond lengths	 Bond angles
 Torsions	 Rings
Ligand PC1 N4 503	
 Bond lengths	 Bond angles
 Torsions	 Rings
Ligand ZMP AC 201	
 Bond lengths	 Bond angles
 Torsions	 Rings





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.

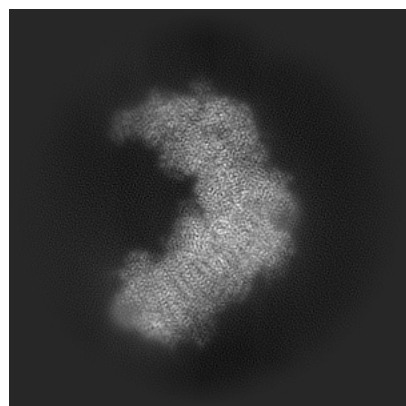
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-36108. These allow visual inspection of the internal detail of the map and identification of artifacts.

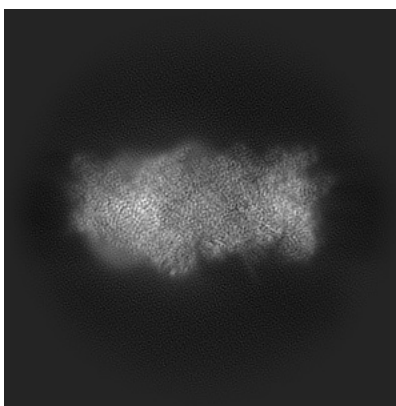
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

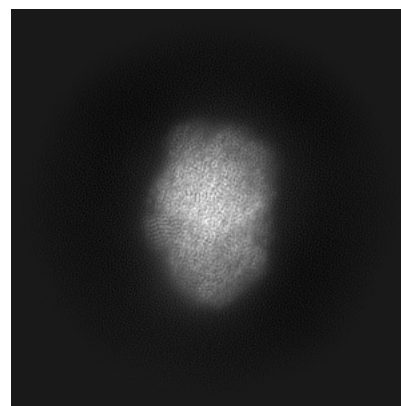
6.1.1 Primary map



X

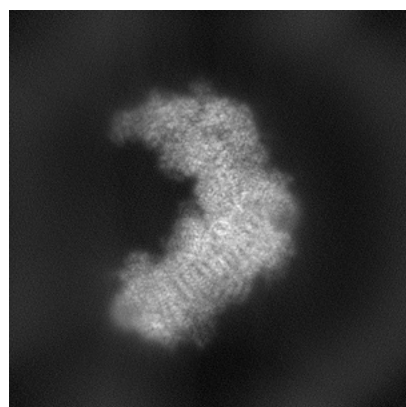


Y

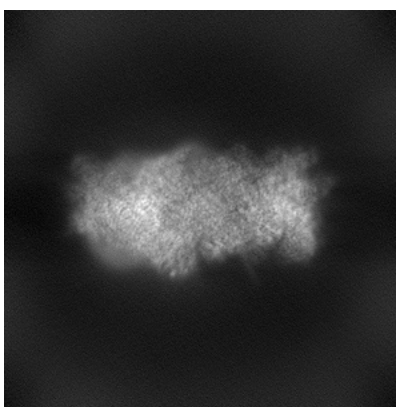


Z

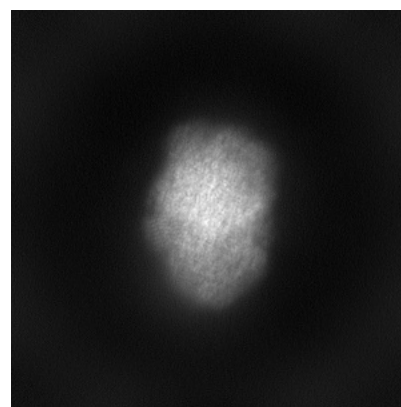
6.1.2 Raw map



X



Y

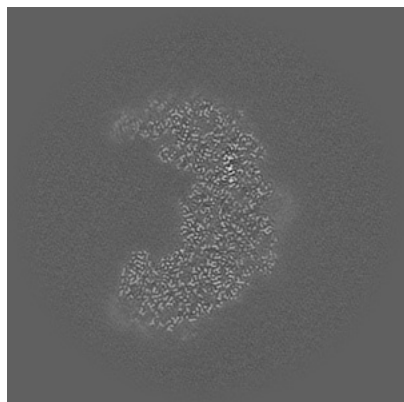


Z

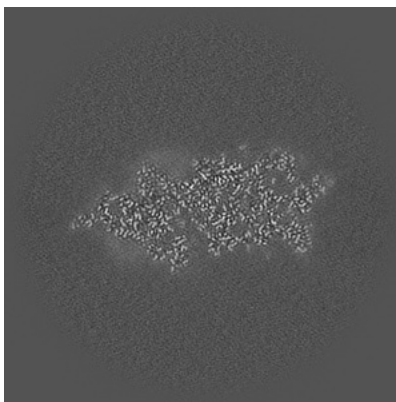
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

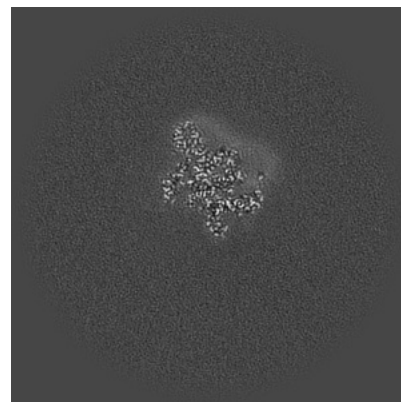
6.2.1 Primary map



X Index: 240

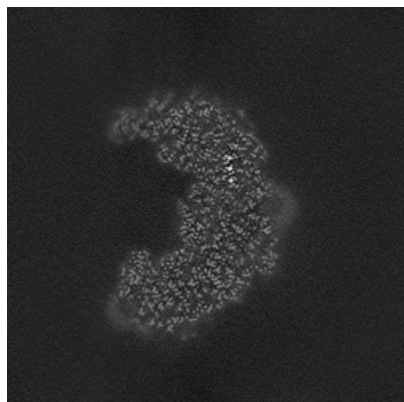


Y Index: 240

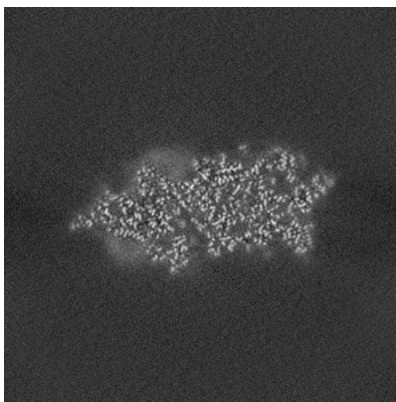


Z Index: 240

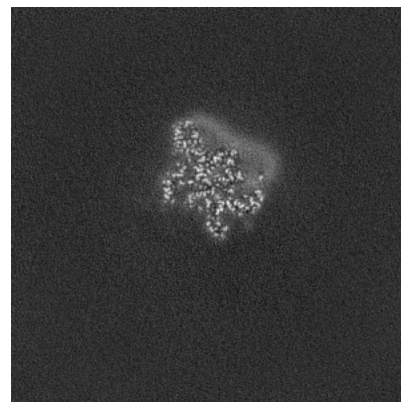
6.2.2 Raw map



X Index: 240



Y Index: 240

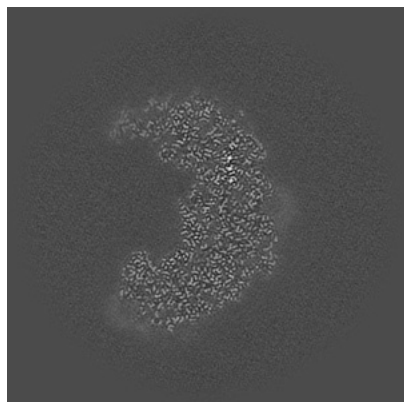


Z Index: 240

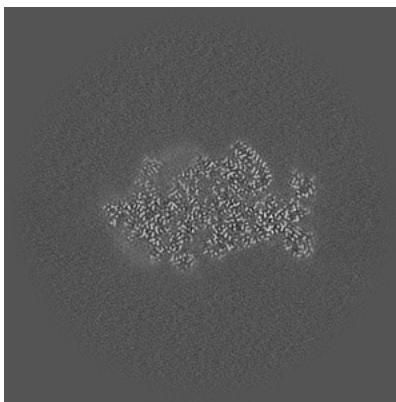
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

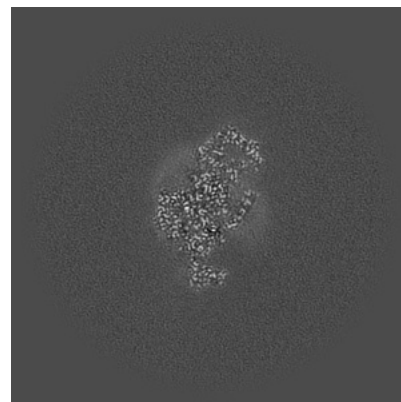
6.3.1 Primary map



X Index: 241

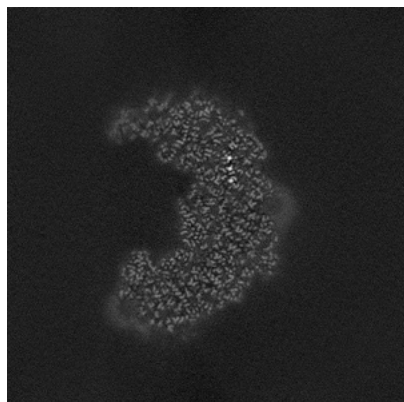


Y Index: 254

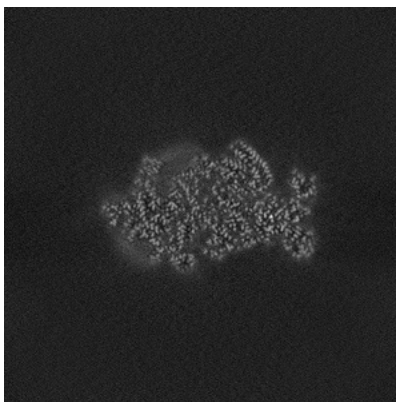


Z Index: 181

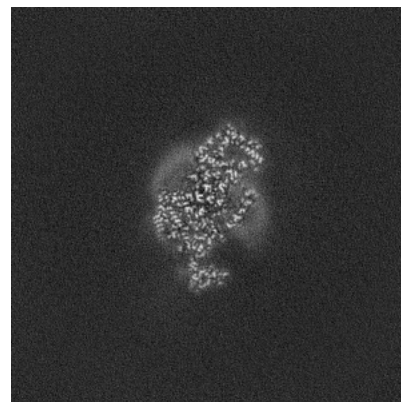
6.3.2 Raw map



X Index: 241



Y Index: 254

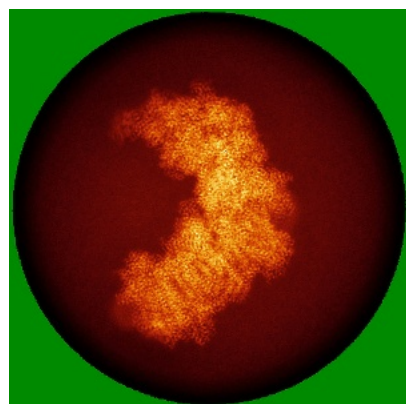


Z Index: 182

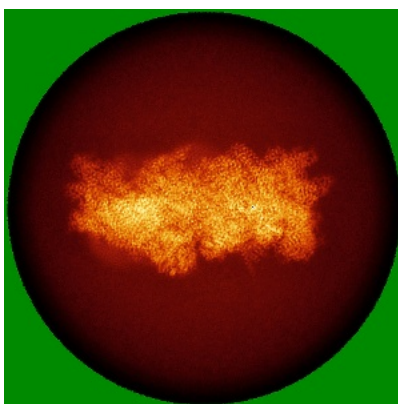
The images above show the largest variance slices of the map in three orthogonal directions.

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

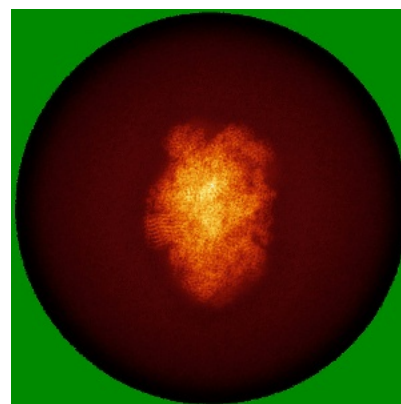
6.4.1 Primary map



X



Y

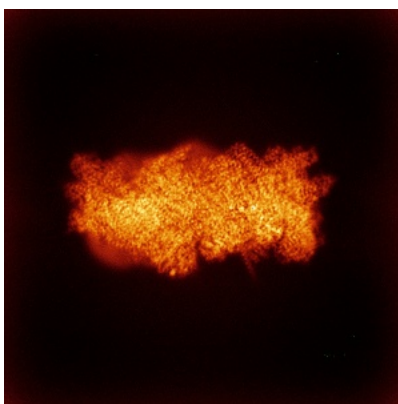


Z

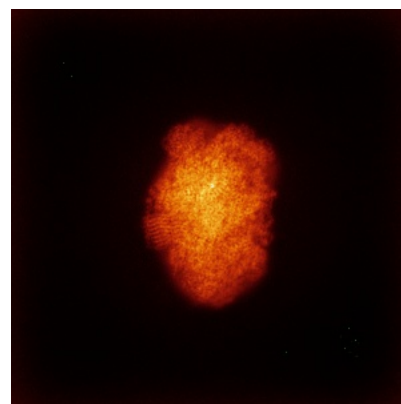
6.4.2 Raw map



X



Y

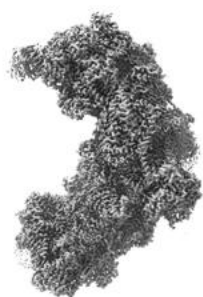


Z

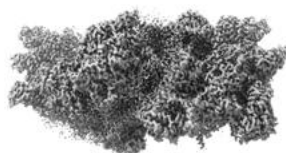
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 4.5. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

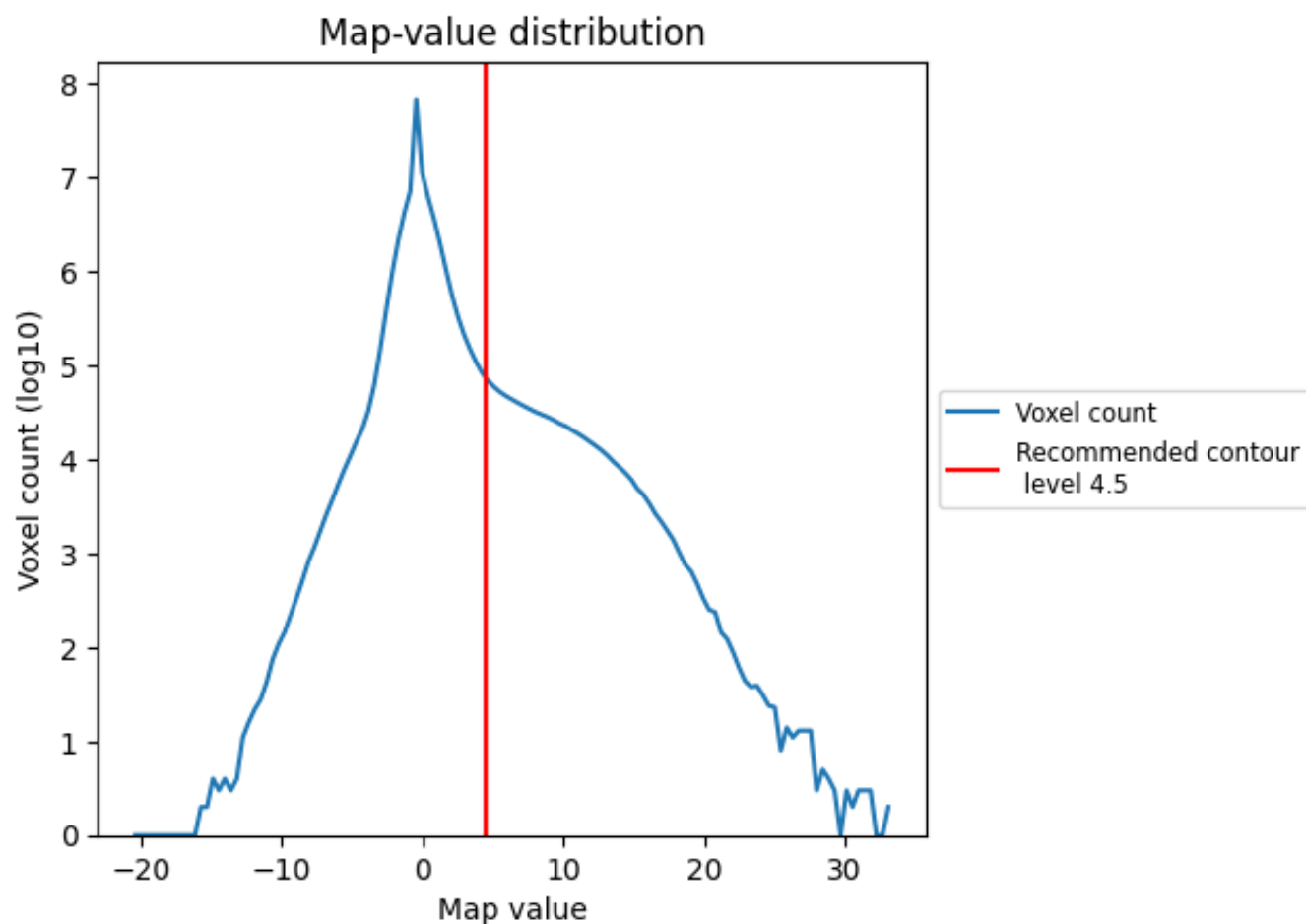
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

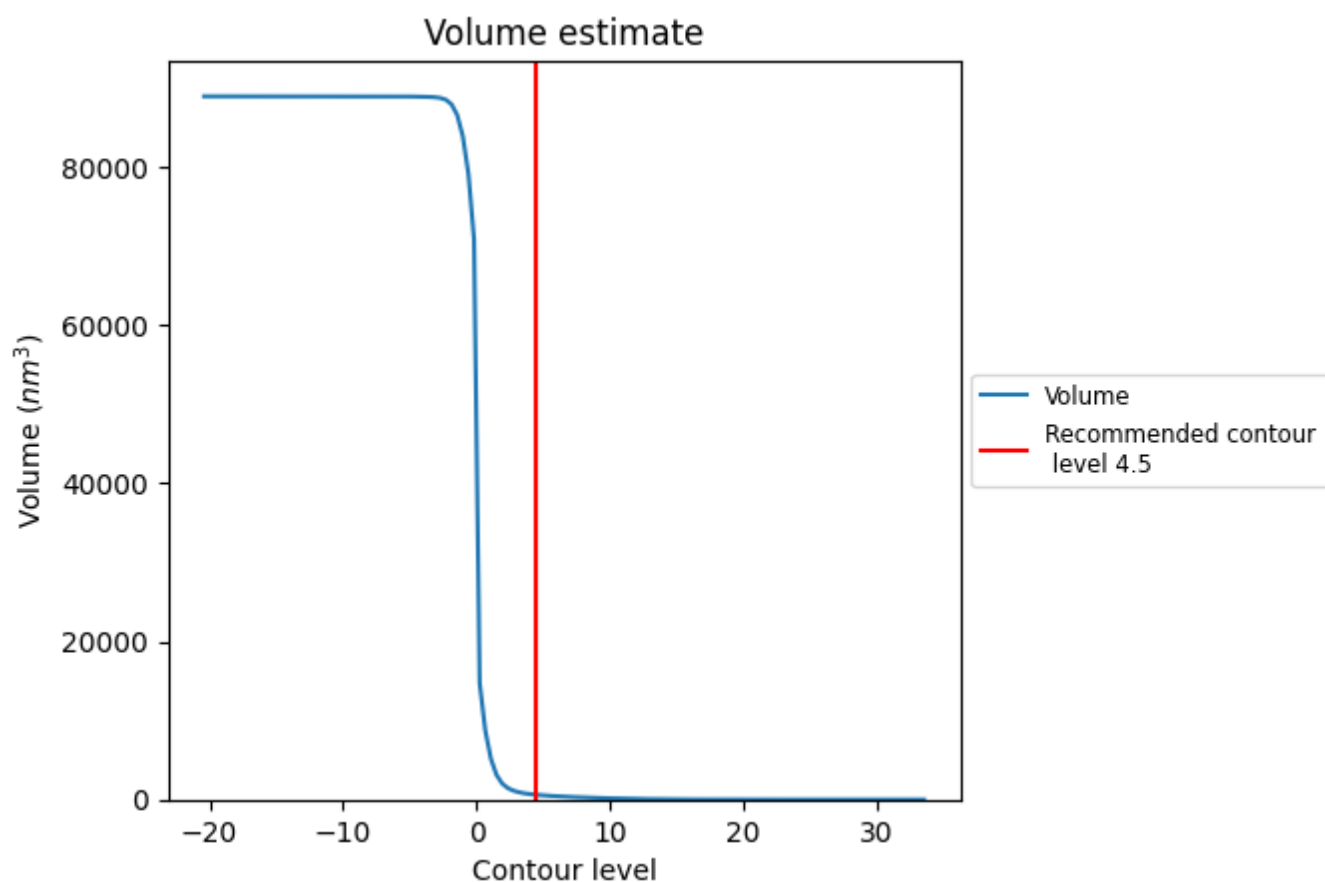
This section contains the results of statistical analysis of the map.

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.

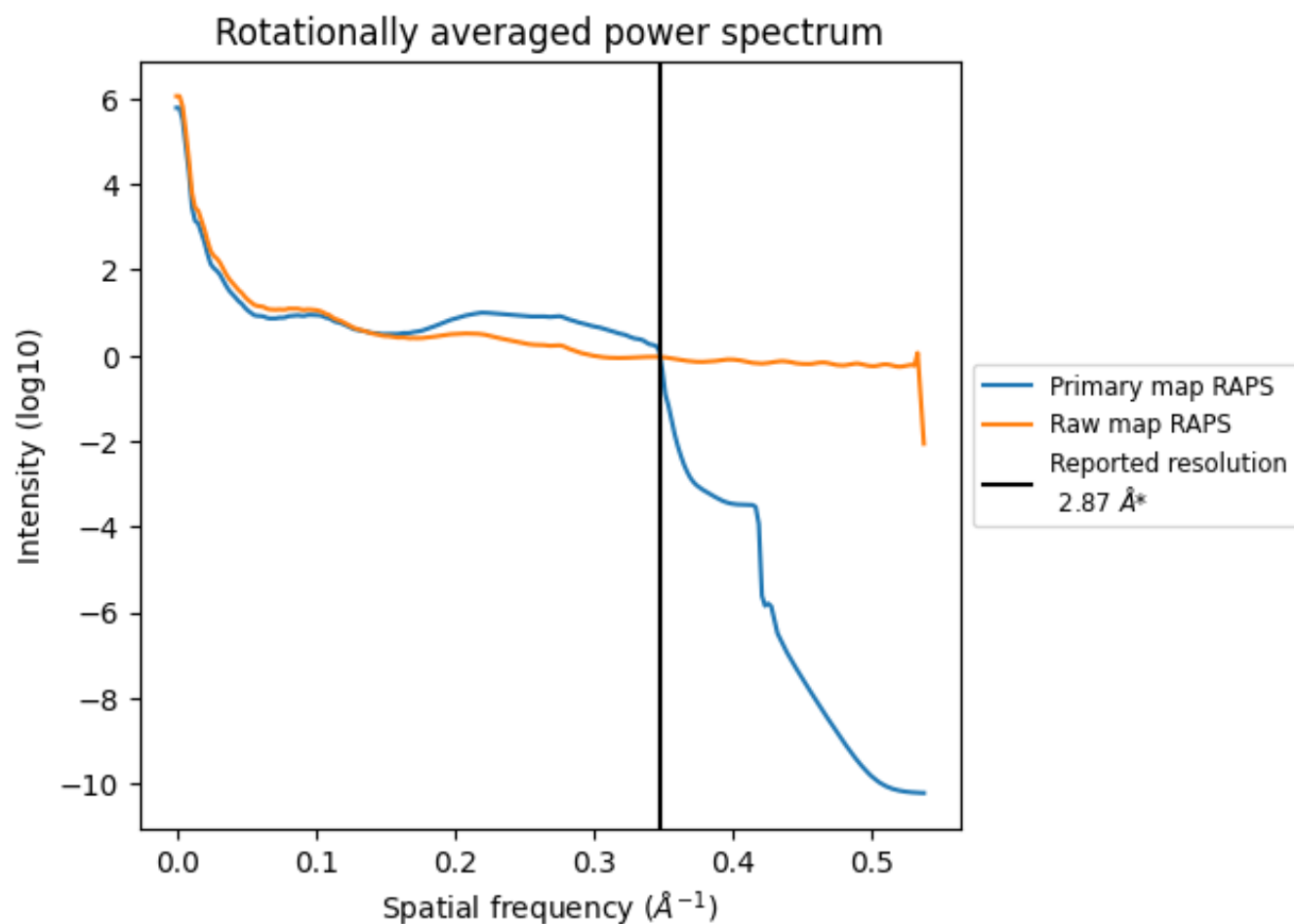
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 596 nm³; this corresponds to an approximate mass of 538 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 ⓘ

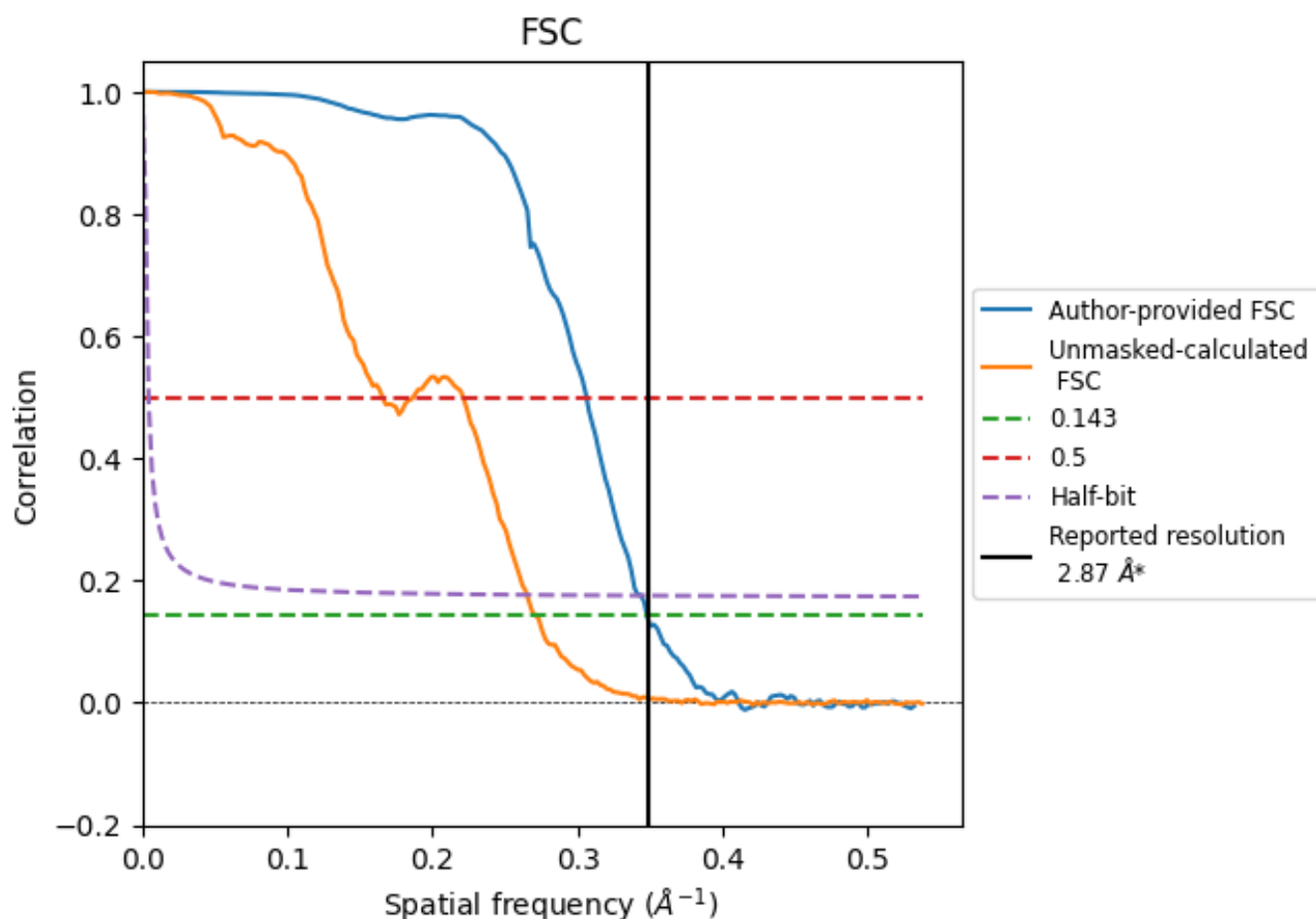


*Reported resolution corresponds to spatial frequency of $0.348 \text{ \AA}^{-1}$

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 0.348 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

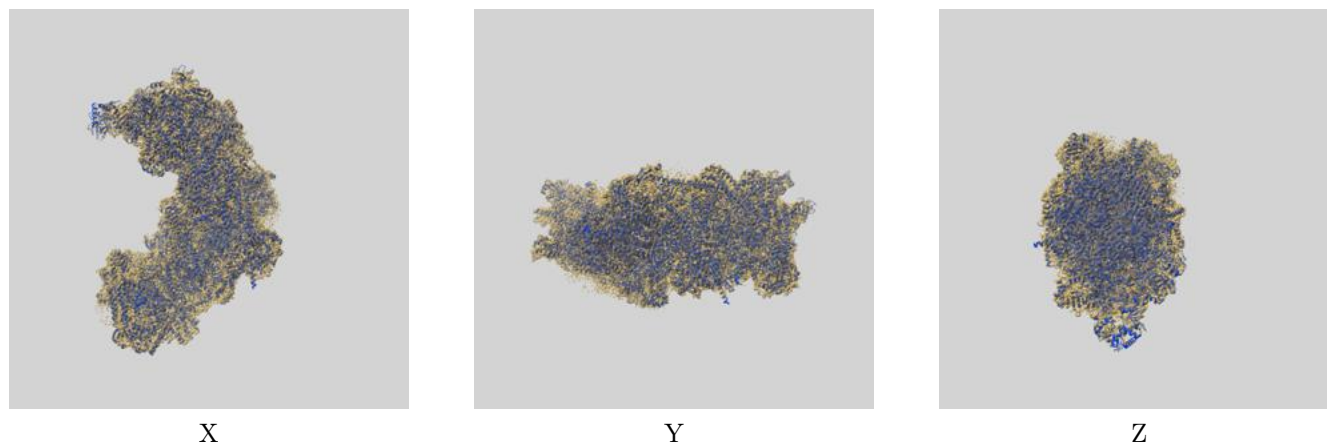
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.87	-	-
Author-provided FSC curve	2.87	3.27	2.91
Unmasked-calculated*	3.69	6.04	3.77

*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 3.69 differs from the reported value 2.87 by more than 10 %

9 Map-model fit [i](#)

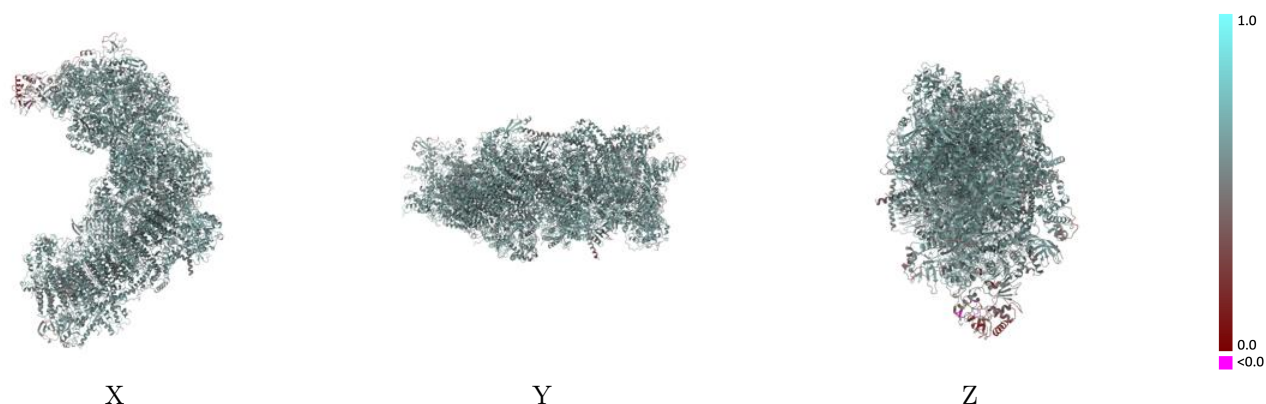
This section contains information regarding the fit between EMDB map EMD-36108 and PDB model 8J9I. Per-residue inclusion information can be found in section [3](#) on page [23](#).

9.1 Map-model overlay [i](#)



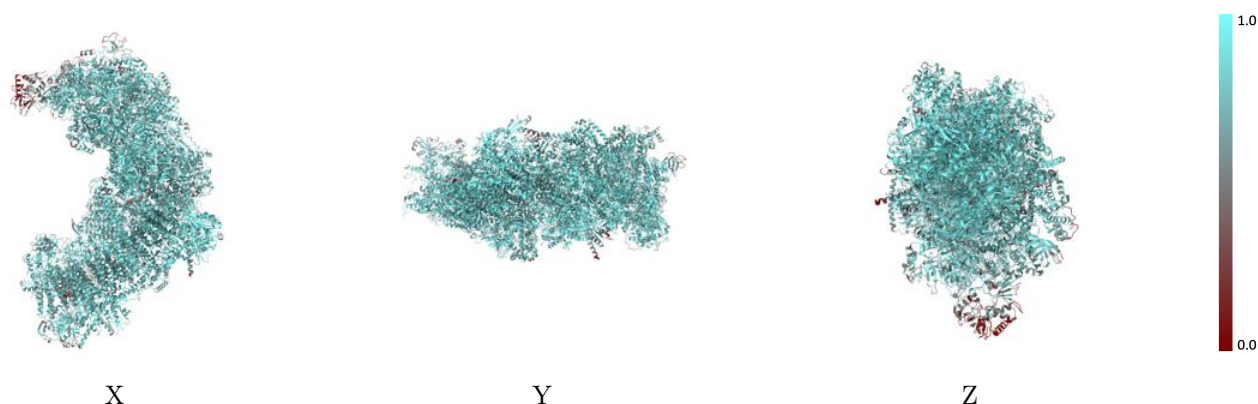
The images above show the 3D surface view of the map at the recommended contour level 4.5 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



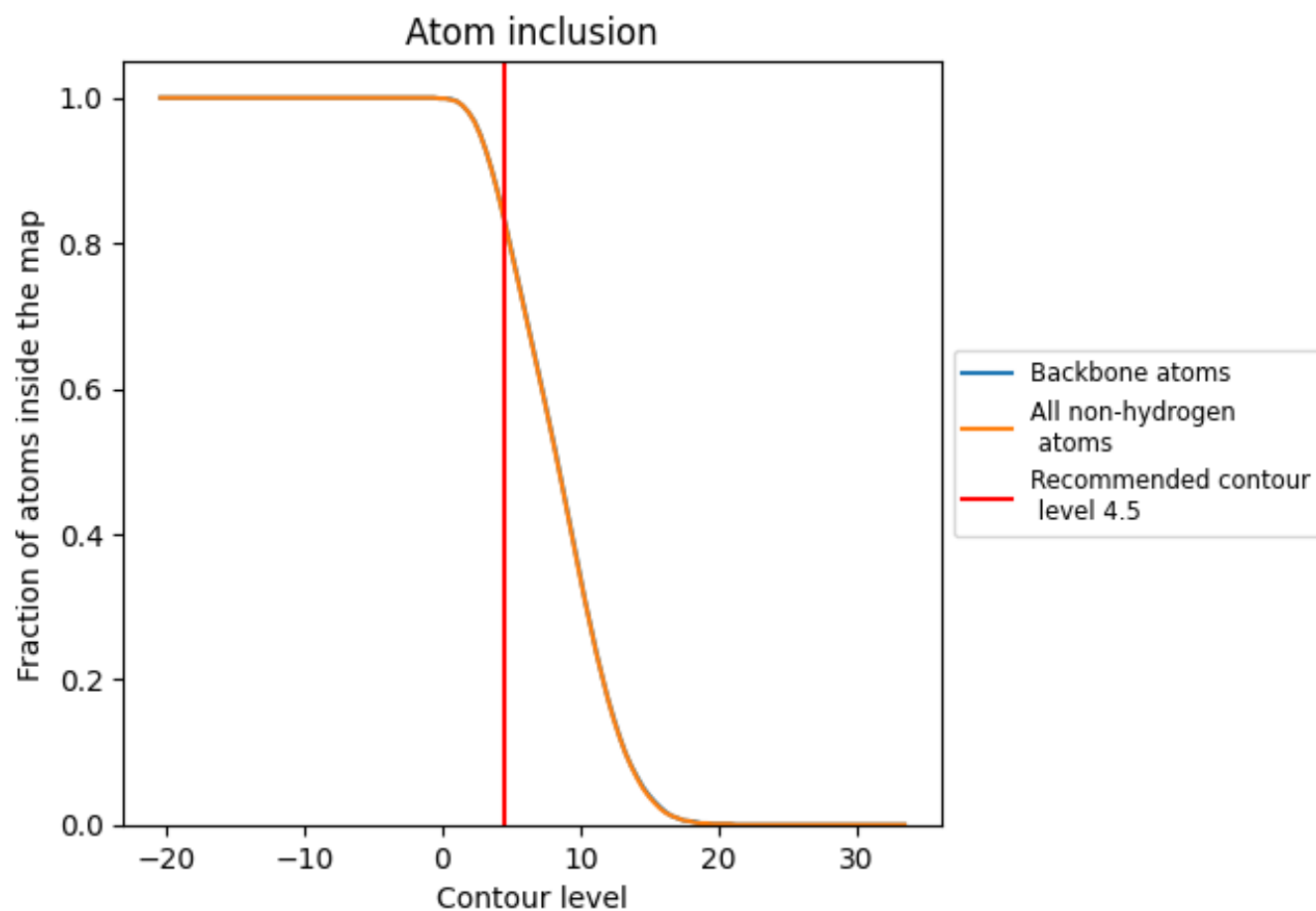
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (4.5).

9.4 Atom inclusion [i](#)



At the recommended contour level, 83% of all backbone atoms, 83% 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 (4.5) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8320	 0.5840
1A	 0.8850	 0.6050
1B	 0.8780	 0.6010
2B	 0.8970	 0.6010
4L	 0.8940	 0.6050
A1	 0.8480	 0.5840
A2	 0.8250	 0.5930
A3	 0.8540	 0.5910
A5	 0.8090	 0.5780
A6	 0.7810	 0.5720
A7	 0.8870	 0.6100
A8	 0.8570	 0.5850
A9	 0.8550	 0.5980
AB	 0.7880	 0.5790
AC	 0.8040	 0.5790
AL	 0.8630	 0.5990
AM	 0.8100	 0.5780
AN	 0.8380	 0.5800
B2	 0.8460	 0.5920
B3	 0.8530	 0.5760
B4	 0.8550	 0.5800
B5	 0.8360	 0.5770
B6	 0.8610	 0.5910
B7	 0.8300	 0.5830
B8	 0.8910	 0.6040
B9	 0.8590	 0.5990
BL	 0.8520	 0.5900
BM	 0.8460	 0.5910
C4	 0.8310	 0.5730
E1	 0.7900	 0.5690
E2	 0.7420	 0.5540
E3	 0.7540	 0.5640
E4	 0.8450	 0.5930
E5	 0.3800	 0.3750
E6	 0.7900	 0.5690



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Chain	Atom inclusion	Q-score
E7	 0.6840	 0.5290
E8	 0.7620	 0.5640
EA	 0.8770	 0.5980
EB	 0.8160	 0.5790
EC	 0.7800	 0.5610
ED	 0.7820	 0.5670
FX	 0.8710	 0.5990
G1	 0.8610	 0.6000
G2	 0.8310	 0.5840
G3	 0.8310	 0.5890
N1	 0.8780	 0.5960
N2	 0.8980	 0.6030
N3	 0.8840	 0.6000
N4	 0.8800	 0.5940
N5	 0.8580	 0.5880
N6	 0.8590	 0.5890
S2	 0.9040	 0.6120
S3	 0.8900	 0.6100
S4	 0.8630	 0.6040
S5	 0.8350	 0.5820
S6	 0.8910	 0.6100
S7	 0.9390	 0.6140
S8	 0.9200	 0.6160
U1	 0.9000	 0.5850
U2	 0.8000	 0.5230
V1	 0.8420	 0.5900
V2	 0.8580	 0.5890