



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 01:17 PM UTC

PDB ID : 5ZCP / pdb_00005zcp
Title : azide-bound cytochrome c oxidase structure determined using the crystals exposed to 20 mM azide solution for 2 days
Authors : Shimada, A.; Hatano, K.; Tadehara, H.; Tsukihara, T.
Deposited on : 2018-02-19
Resolution : 1.65 Å(reported)

This is a wwPDB X-ray Structure 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/XrayValidationReportHelp>

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:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

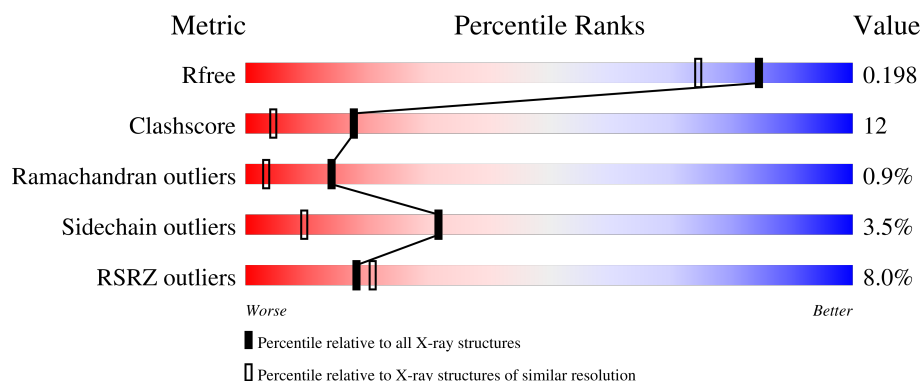
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2563 (1.66-1.66)
Clashscore	190562	2662 (1.66-1.66)
Ramachandran outliers	187476	2621 (1.66-1.66)
Sidechain outliers	187428	2621 (1.66-1.66)
RSRZ outliers	180081	2564 (1.66-1.66)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower 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 electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	514	0% 74% 22% .
1	N	514	2% 67% 29% .
2	B	227	12% 64% 28% 7% .
2	O	227	11% 70% 26% .
3	C	261	3% 74% 24% ..

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Mol	Chain	Length	Quality of chain
3	P	261	
4	D	147	
4	Q	147	
5	E	109	
5	R	109	
6	F	98	
6	S	98	
7	G	85	
7	T	85	
8	H	85	
8	U	85	
9	I	73	
9	V	73	
10	J	59	
10	W	59	
11	K	56	
11	X	56	
12	L	47	
12	Y	47	
13	M	46	
13	Z	46	

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
18	AZI	A	607[A]	-	X	X	-
18	AZI	A	607[B]	-	X	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
18	AZI	N	608[A]	-	X	-	-
18	AZI	N	608[B]	-	X	-	-
21	EDO	A	616	-	X	-	-
21	EDO	D	202	-	-	X	-
27	CDL	N	601	-	-	X	-
27	CDL	P	305	-	-	X	-

2 Entry composition

There are 30 unique types of molecules in this entry. The entry contains 33609 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, 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 Cytochrome c oxidase subunit 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	514	Total	C	N	O	S	0	22	0
			4193	2793	649	709	42			
1	N	514	Total	C	N	O	S	0	20	0
			4179	2786	647	704	42			

- Molecule 2 is a protein called Cytochrome c oxidase subunit 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	227	Total	C	N	O	S	0	9	0
			1899	1234	292	353	20			
2	O	227	Total	C	N	O	S	0	5	0
			1870	1215	288	347	20			

- Molecule 3 is a protein called Cytochrome c oxidase subunit 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	259	Total	C	N	O	S	0	9	0
			2185	1457	349	363	16			
3	P	259	Total	C	N	O	S	0	9	0
			2185	1457	349	363	16			

- Molecule 4 is a protein called Cytochrome c oxidase subunit 4 isoform 1, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	144	Total	C	N	O	S	0	4	0
			1233	803	204	222	4			
4	Q	144	Total	C	N	O	S	0	3	0
			1224	797	202	221	4			

- Molecule 5 is a protein called Cytochrome c oxidase subunit 5A, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	105	Total	C	N	O	S	0	0	0
			852	544	144	162	2			
5	R	105	Total	C	N	O	S	0	1	0
			863	550	148	163	2			

- Molecule 6 is a protein called Cytochrome c oxidase subunit 5B, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	98	Total	C	N	O	S	0	4	0
			778	481	139	152	6			
6	S	98	Total	C	N	O	S	0	2	0
			763	473	136	148	6			

- Molecule 7 is a protein called Cytochrome c oxidase subunit 6A2, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
7	G	84	Total 686	C 440	N 130	O 114	P 1	S 1	0	1	0
7	T	84	Total 686	C 440	N 130	O 114	P 1	S 1	0	1	0

- Molecule 8 is a protein called Cytochrome c oxidase subunit 6B1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	79	Total	C	N	O	S	0	0	0
			662	417	121	119	5			
8	U	79	Total	C	N	O	S	0	0	0
			662	417	121	119	5			

- Molecule 9 is a protein called Cytochrome c oxidase subunit 6C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	I	73	Total	C	N	O	S	0	0	0
			601	390	107	100	4			
9	V	73	Total	C	N	O	S	0	0	0
			601	390	107	100	4			

- Molecule 10 is a protein called Cytochrome c oxidase subunit 7A1, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	J	58	Total	C	N	O	S	0	0	0
			460	297	78	82	3			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	W	58	Total	C	N	O	S	0	1	0
			469	302	79	85	3			

- Molecule 11 is a protein called Cytochrome c oxidase subunit 7B, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	K	49	Total	C	N	O	S	0	0	0
			384	250	65	67	2			
11	X	49	Total	C	N	O	S	0	1	0
			391	255	66	68	2			

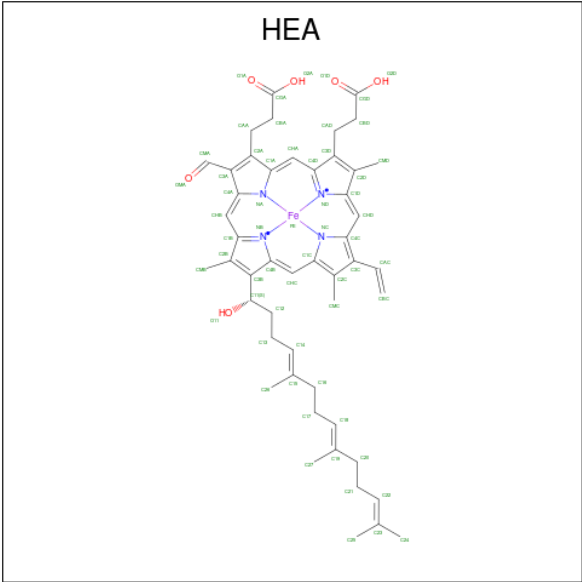
- Molecule 12 is a protein called Cytochrome c oxidase subunit 7C, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	L	46	Total	C	N	O	S	0	0	0
			380	254	64	60	2			
12	Y	46	Total	C	N	O	S	0	1	0
			388	259	65	61	3			

- Molecule 13 is a protein called Cytochrome c oxidase subunit 8B, mitochondrial.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
13	M	43	Total	C	N	O	0	0	0
			335	223	53	59			
13	Z	43	Total	C	N	O	0	0	0
			335	223	53	59			

- Molecule 14 is HEME-A (CCD ID: HEA) (formula: $C_{49}H_{56}FeN_4O_6$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
14	A	1	Total 60	C 49	Fe 1	N 4	O 6	0	0
14	A	1	Total 120	C 98	Fe 2	N 8	O 12	0	1
14	N	1	Total 60	C 49	Fe 1	N 4	O 6	0	0
14	N	1	Total 120	C 98	Fe 2	N 8	O 12	0	1

- Molecule 15 is COPPER (II) ION (CCD ID: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
15	A	1	Total	Cu	0	0
			1	1		
15	N	1	Total	Cu	0	0
			1	1		

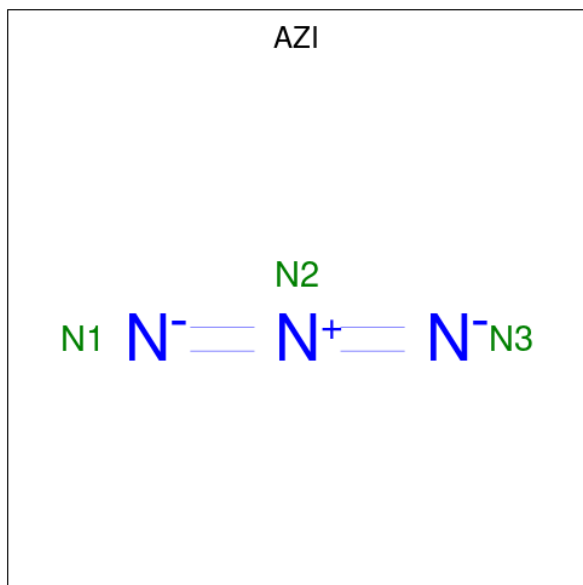
- Molecule 16 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
16	A	1	Total	Mg	0	0
			1	1		
16	N	1	Total	Mg	0	0
			1	1		

- Molecule 17 is SODIUM ION (CCD ID: NA) (formula: Na).

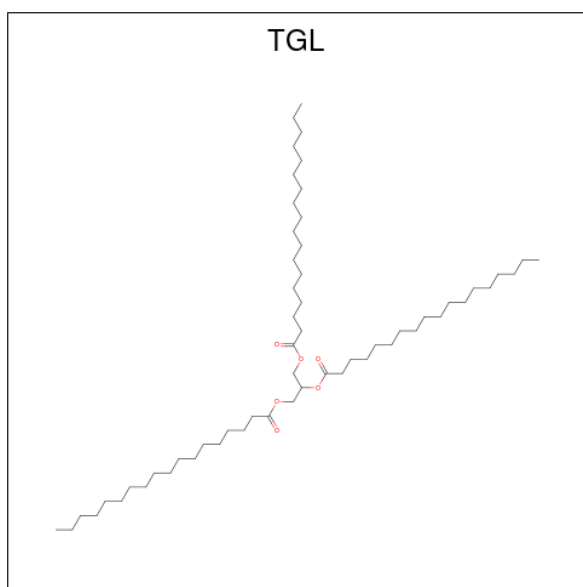
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
17	A	1	Total Na 1 1	0	0
17	N	1	Total Na 1 1	0	0

- Molecule 18 is AZIDE ION (CCD ID: AZI) (formula: N₃).



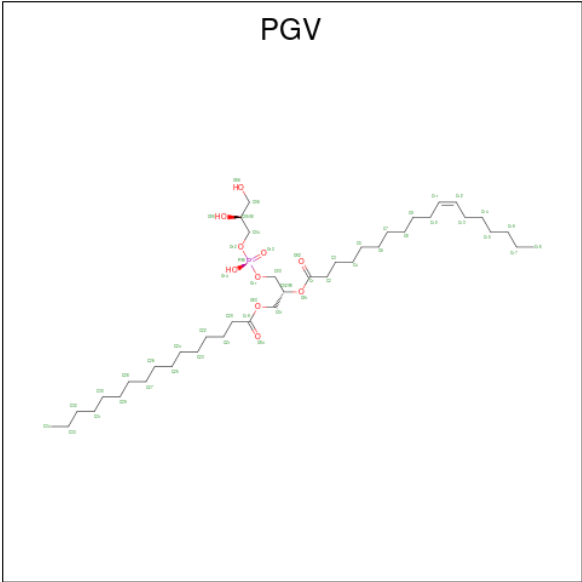
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
18	A	1	Total N 3 3	0	1
18	A	1	Total N 6 6	0	1
18	N	1	Total N 3 3	0	1
18	N	1	Total N 6 6	0	1

- Molecule 19 is TRISTEAROYLGLYCEROL (CCD ID: TGL) (formula: C₅₇H₁₁₀O₆).



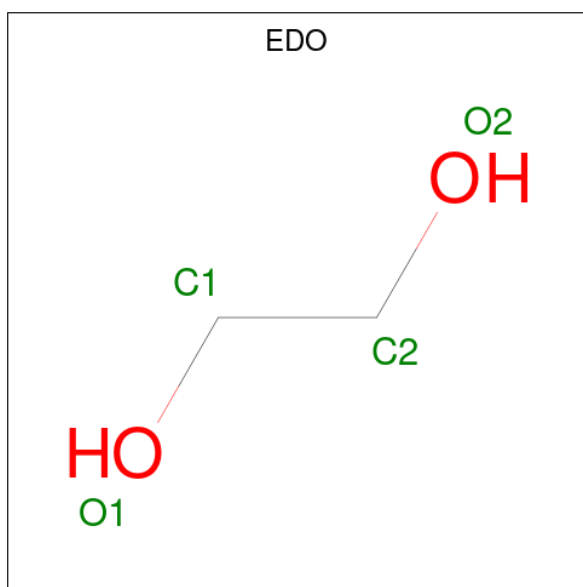
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
19	A	1	Total	C	O	0	0
			63	57	6		
19	A	1	Total	C	O	0	0
			63	57	6		
19	D	1	Total	C	O	0	0
			63	57	6		
19	N	1	Total	C	O	0	0
			63	57	6		
19	Q	1	Total	C	O	0	0
			63	57	6		
19	Y	1	Total	C	O	0	0
			63	57	6		

- Molecule 20 is (1R)-2-{{[[(2S)-2,3-DIHYDROXYPROPYL]OXY}(HYDROXY)PHOSPHORYL]OXY}-1-[(PALMITOYLOXY)METHYL]ETHYL (11E)-OCTADEC-11-ENOATE (CCD ID: PGV) (formula: C₄₀H₇₇O₁₀P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
20	A	1	Total	C	O	P	0	0
			51	40	10	1		
20	A	1	Total	C	O	P	0	0
			51	40	10	1		
20	C	1	Total	C	O	P	0	0
			51	40	10	1		
20	C	1	Total	C	O	P	0	0
			51	40	10	1		
20	N	1	Total	C	O	P	0	0
			51	40	10	1		
20	N	1	Total	C	O	P	0	0
			51	40	10	1		
20	P	1	Total	C	O	P	0	0
			51	40	10	1		
20	P	1	Total	C	O	P	0	0
			51	40	10	1		

- Molecule 21 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
21	A	1	Total	C	O	0	0
			4	2	2		
21	A	1	Total	C	O	0	0
			4	2	2		
21	A	1	Total	C	O	0	0
			4	2	2		
21	A	1	Total	C	O	0	0
			4	2	2		
21	A	1	Total	C	O	0	0
			4	2	2		
21	A	1	Total	C	O	0	0
			4	2	2		
21	A	1	Total	C	O	0	0
			4	2	2		
21	B	1	Total	C	O	0	0
			4	2	2		
21	B	1	Total	C	O	0	0
			4	2	2		
21	B	1	Total	C	O	0	0
			4	2	2		
21	B	1	Total	C	O	0	0
			4	2	2		
21	C	1	Total	C	O	0	0
			4	2	2		

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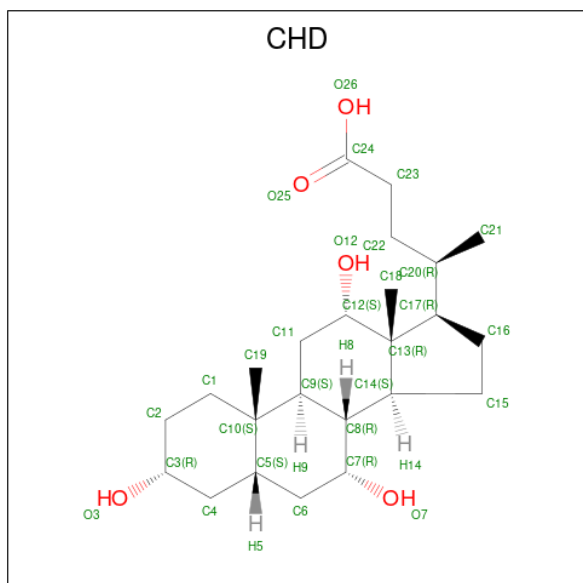
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
21	D	1	Total 4	C 2	O 2	0	0
21	D	1	Total 4	C 2	O 2	0	0
21	E	1	Total 4	C 2	O 2	0	0
21	E	1	Total 4	C 2	O 2	0	0
21	E	1	Total 4	C 2	O 2	0	0
21	F	1	Total 4	C 2	O 2	0	0
21	F	1	Total 4	C 2	O 2	0	0
21	F	1	Total 4	C 2	O 2	0	0
21	G	1	Total 4	C 2	O 2	0	0
21	G	1	Total 4	C 2	O 2	0	0
21	L	1	Total 4	C 2	O 2	0	0
21	M	1	Total 4	C 2	O 2	0	0
21	N	1	Total 4	C 2	O 2	0	0
21	N	1	Total 4	C 2	O 2	0	0
21	N	1	Total 4	C 2	O 2	0	0
21	N	1	Total 4	C 2	O 2	0	0
21	N	1	Total 4	C 2	O 2	0	0
21	N	1	Total 4	C 2	O 2	0	0
21	N	1	Total 4	C 2	O 2	0	0
21	N	1	Total 4	C 2	O 2	0	0
21	N	1	Total 4	C 2	O 2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
21	O	1	Total	C	O	0	0
			4	2	2		
21	P	1	Total	C	O	0	0
			4	2	2		
21	P	1	Total	C	O	0	0
			4	2	2		
21	R	1	Total	C	O	0	0
			4	2	2		
21	S	1	Total	C	O	0	0
			4	2	2		
21	S	1	Total	C	O	0	0
			4	2	2		
21	S	1	Total	C	O	0	0
			4	2	2		
21	T	1	Total	C	O	0	0
			4	2	2		
21	W	1	Total	C	O	0	0
			4	2	2		
21	Y	1	Total	C	O	0	0
			4	2	2		

- Molecule 22 is CHOLIC ACID (CCD ID: CHD) (formula: $C_{24}H_{40}O_5$).



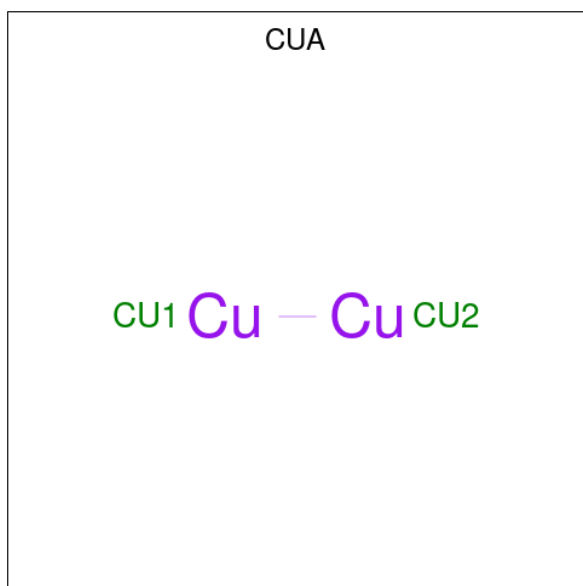
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
22	B	1	Total	C	O	0	0
			29	24	5		

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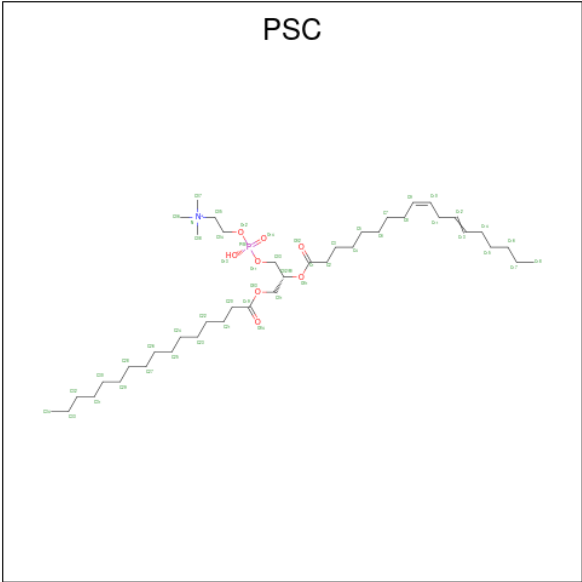
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
22	C	1	Total C O 29 24 5	0	0
22	C	1	Total C O 29 24 5	0	0
22	G	1	Total C O 29 24 5	0	0
22	J	1	Total C O 29 24 5	0	0
22	P	1	Total C O 29 24 5	0	0
22	P	1	Total C O 29 24 5	0	0
22	W	1	Total C O 29 24 5	0	0

- Molecule 23 is DINUCLEAR COPPER ION (CCD ID: CUA) (formula: Cu₂).



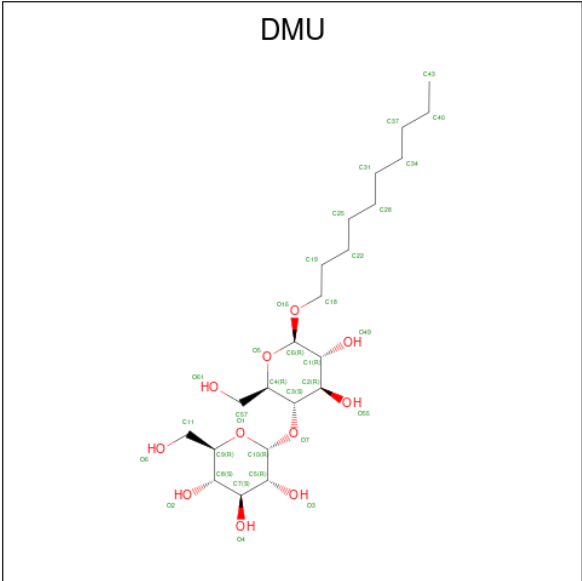
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
23	B	1	Total Cu 2 2	0	0
23	O	1	Total Cu 2 2	0	0

- Molecule 24 is (7R,17E,20E)-4-HYDROXY-N,N,N-TRIMETHYL-9-OXO-7-[(PALMITOYLOXY)METHYL]-3,5,8-TRIOXA-4-PHOSPHAHEXACOSA-17,20-DIEN-1-AMINIUM 4-OXIDE (CCD ID: PSC) (formula: C₄₂H₈₁NO₈P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
24	B	1	Total	C	N	O	P	0	0
			52	42	1	8	1		
24	N	1	Total	C	N	O	P	0	0
			52	42	1	8	1		

- Molecule 25 is DECYL-BETA-D-MALTOPYRANOSIDE (CCD ID: DMU) (formula: $C_{22}H_{42}O_{11}$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
25	C	1	Total	C	O	0	0
			33	22	11		

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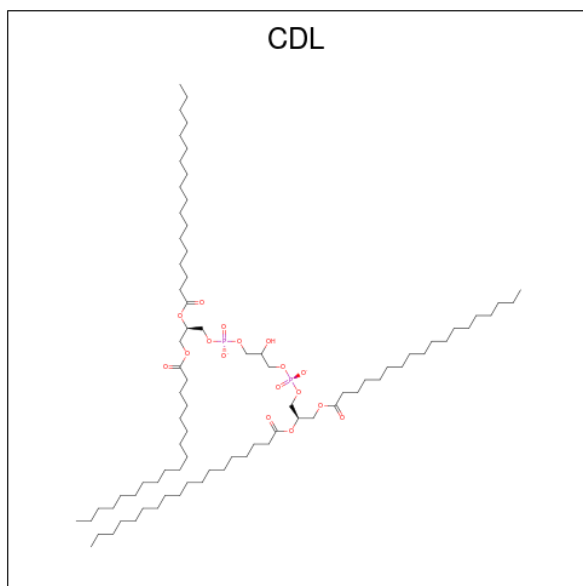
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
25	C	1	Total	C	O	0	0
			33	22	11		
25	C	1	Total	C	O	0	0
			33	22	11		
25	M	1	Total	C	O	0	0
			33	22	11		
25	P	1	Total	C	O	0	0
			33	22	11		
25	P	1	Total	C	O	0	0
			33	22	11		
25	P	1	Total	C	O	0	0
			33	22	11		
25	Z	1	Total	C	O	0	0
			33	22	11		

- Molecule 26 is UNKNOWN ATOM OR ION (CCD ID: UNX) (formula: X).

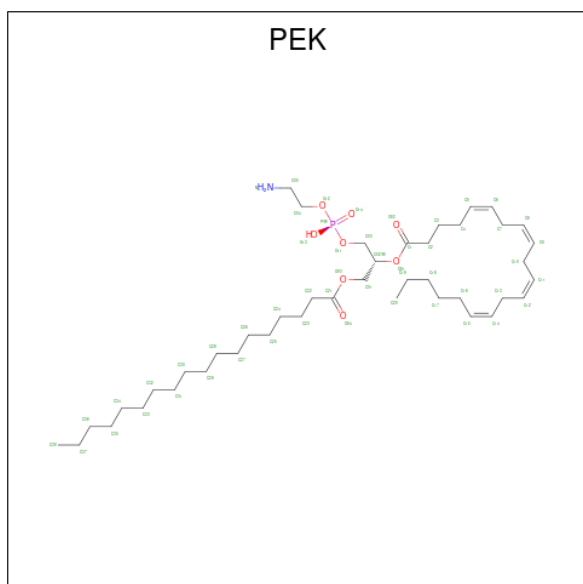
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
26	C	1	Total	X	0	0
			1	1		
26	P	1	Total	X	0	0
			1	1		

- Molecule 27 is CARDIOLIPIN (CCD ID: CDL) (formula: $C_{81}H_{156}O_{17}P_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
27	C	1	Total	C	O	P	0	0
			100	81	17	2		
27	N	1	Total	C	O	P	0	0
			100	81	17	2		
27	P	1	Total	C	O	P	0	0
			100	81	17	2		
27	T	1	Total	C	O	P	0	0
			100	81	17	2		

- Molecule 28 is (1S)-2-[[[(2-AMINOETHOXY)(HYDROXY)PHOSPHORYL]OXY}-1-[(STEAROYL)OXY]METHYL]ETHYL (5E,8E,11E,14E)-ICOSA-5,8,11,14-TETRAENOATE (CCD ID: PEK) (formula: C₄₃H₇₈NO₈P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
28	C	1	Total 53	C 43	N 1	O 8	P 1	0	0
28	C	1	Total 53	C 43	N 1	O 8	P 1	0	0
28	G	1	Total 53	C 43	N 1	O 8	P 1	0	0
28	G	1	Total 53	C 43	N 1	O 8	P 1	0	0
28	P	1	Total 53	C 43	N 1	O 8	P 1	0	0
28	T	1	Total 53	C 43	N 1	O 8	P 1	0	0

- Molecule 29 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
29	F	1	Total 1	Zn 1	0	0
29	S	1	Total 1	Zn 1	0	0

- Molecule 30 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
30	A	235	Total 235	O 235	0	0
30	B	161	Total 162	O 162	0	1
30	C	110	Total 110	O 110	0	0
30	D	118	Total 118	O 118	0	0
30	E	93	Total 93	O 93	0	0
30	F	94	Total 94	O 94	0	0
30	G	42	Total 42	O 42	0	0
30	H	42	Total 42	O 42	0	0
30	I	27	Total 27	O 27	0	0
30	J	23	Total 23	O 23	0	0
30	K	26	Total 26	O 26	0	0
30	L	38	Total 38	O 38	0	0
30	M	25	Total 25	O 25	0	0
30	N	203	Total 203	O 203	0	0
30	O	99	Total 100	O 100	0	1
30	P	95	Total 95	O 95	0	0
30	Q	29	Total 29	O 29	0	0
30	R	40	Total 40	O 40	0	0

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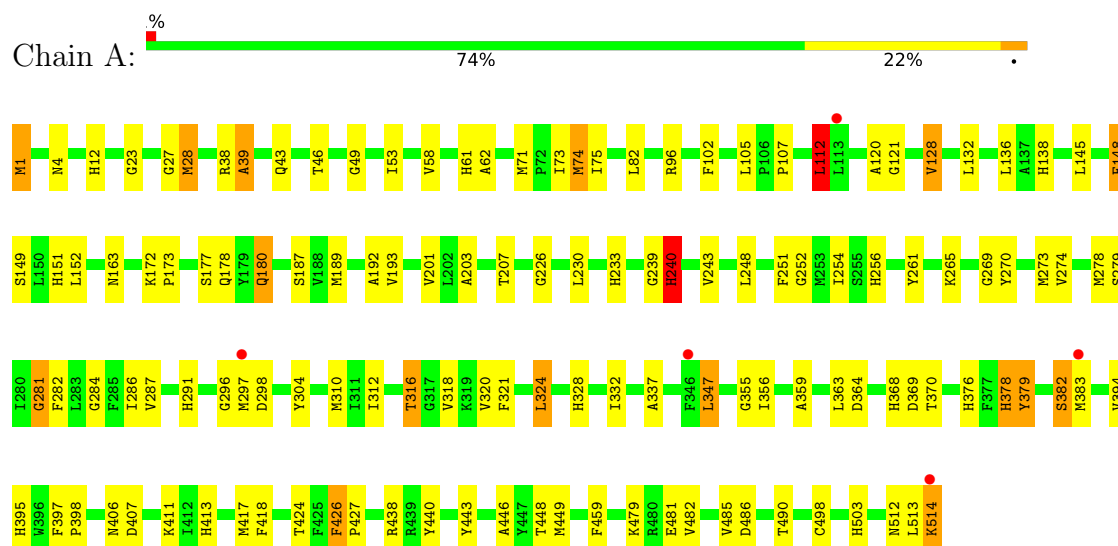
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
30	S	45	Total 45	O 45	0	0
30	T	35	Total 35	O 35	0	0
30	U	29	Total 29	O 29	0	0
30	V	16	Total 16	O 16	0	0
30	W	7	Total 7	O 7	0	0
30	X	11	Total 11	O 11	0	0
30	Y	12	Total 12	O 12	0	0
30	Z	12	Total 12	O 12	0	0

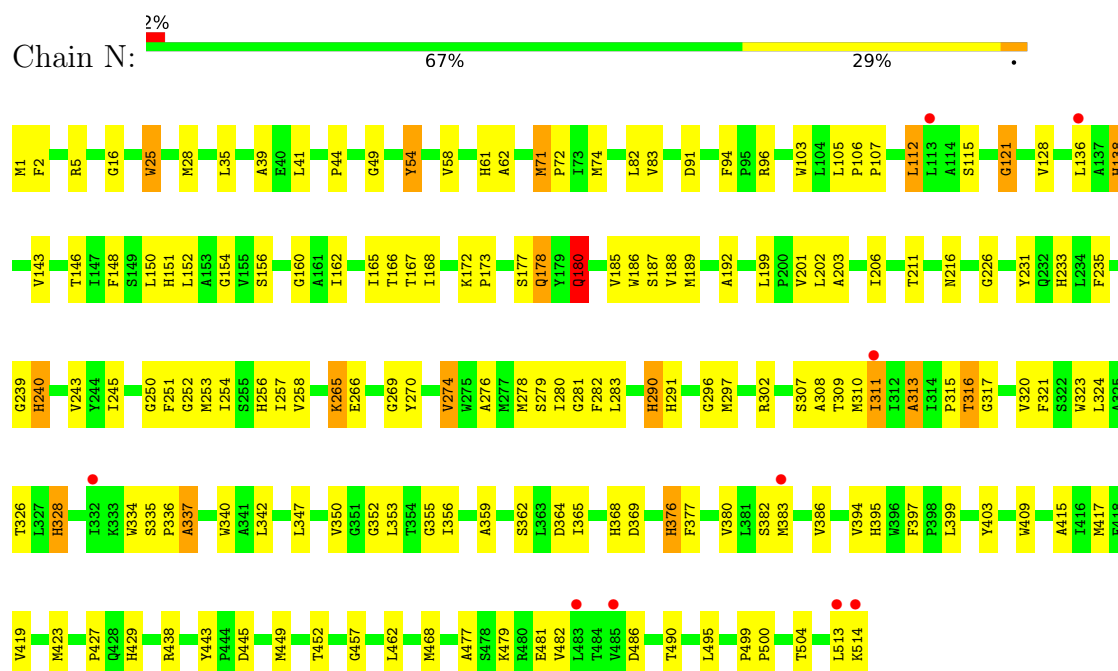
3 Residue-property plots

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 electron 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 dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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.

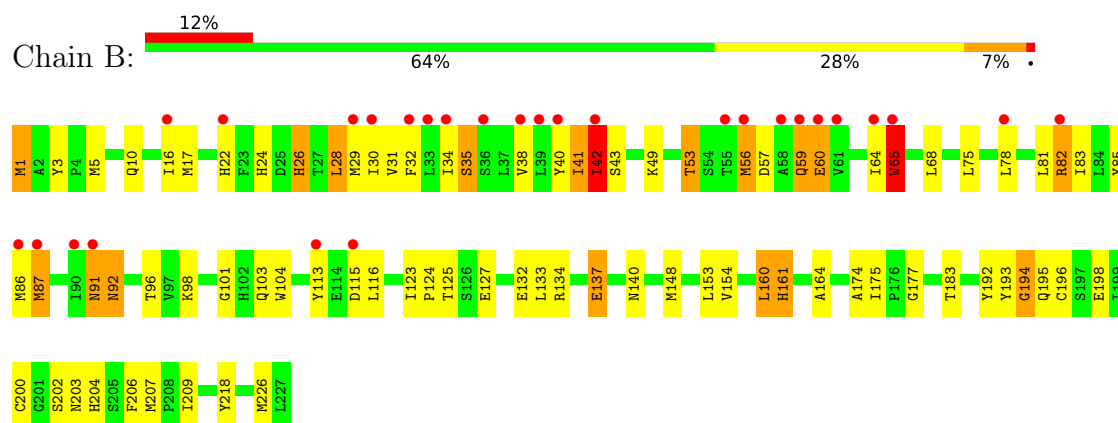
• Molecule 1: Cytochrome c oxidase subunit 1



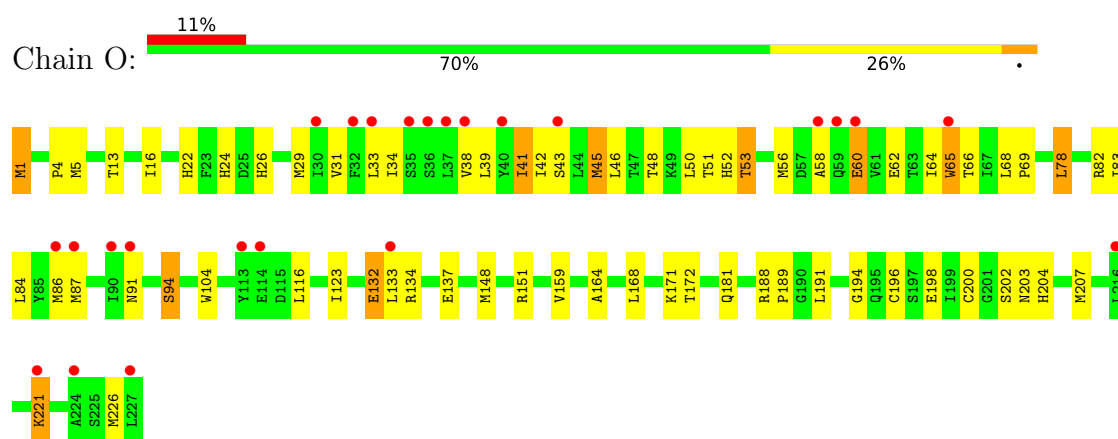
• Molecule 1: Cytochrome c oxidase subunit 1



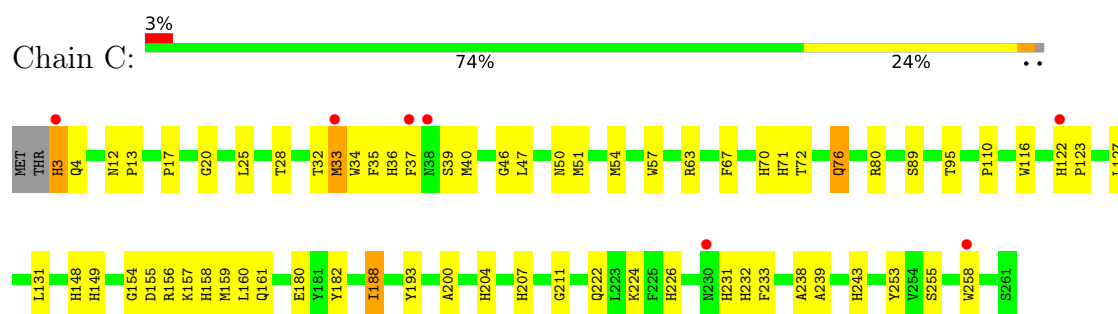
- Molecule 2: Cytochrome c oxidase subunit 2



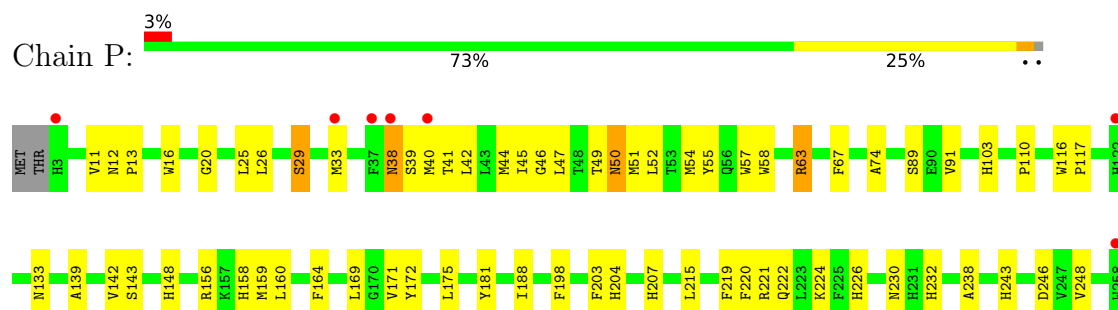
- Molecule 2: Cytochrome c oxidase subunit 2



- Molecule 3: Cytochrome c oxidase subunit 3

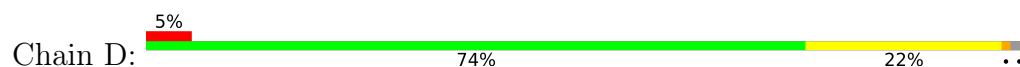


- Molecule 3: Cytochrome c oxidase subunit 3

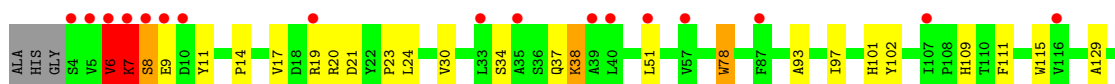
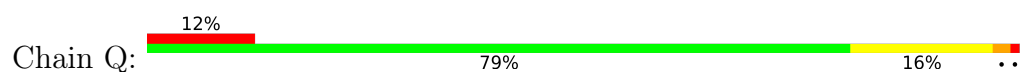




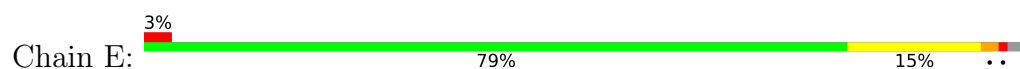
- Molecule 4: Cytochrome c oxidase subunit 4 isoform 1, mitochondrial



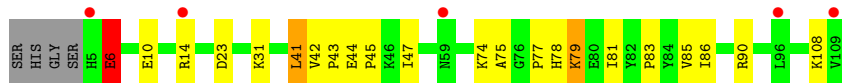
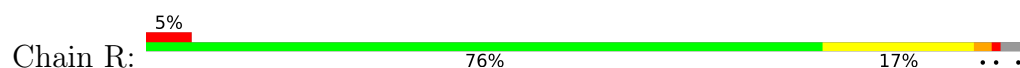
- Molecule 4: Cytochrome c oxidase subunit 4 isoform 1, mitochondrial



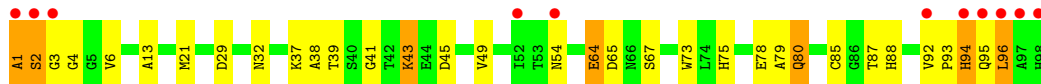
- Molecule 5: Cytochrome c oxidase subunit 5A, mitochondrial



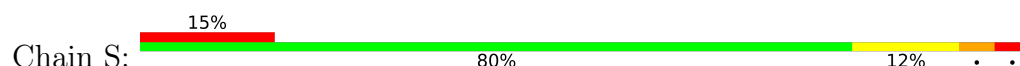
- Molecule 5: Cytochrome c oxidase subunit 5A, mitochondrial



- Molecule 6: Cytochrome c oxidase subunit 5B, mitochondrial

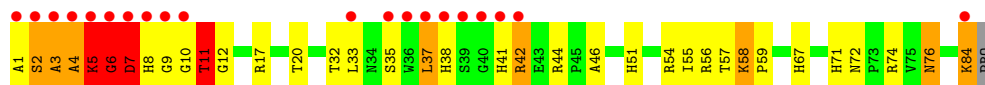


- Molecule 6: Cytochrome c oxidase subunit 5B, mitochondrial

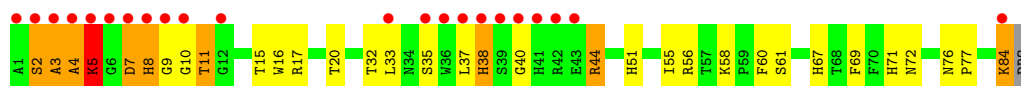




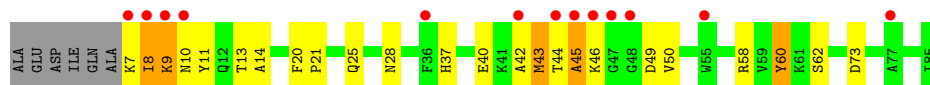
- Molecule 7: Cytochrome c oxidase subunit 6A2, mitochondrial



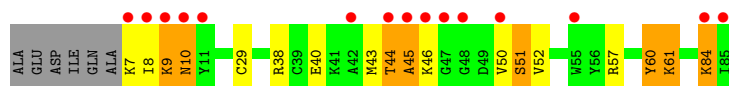
- Molecule 7: Cytochrome c oxidase subunit 6A2, mitochondrial



- Molecule 8: Cytochrome c oxidase subunit 6B1



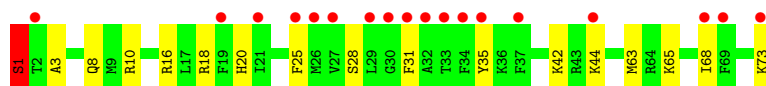
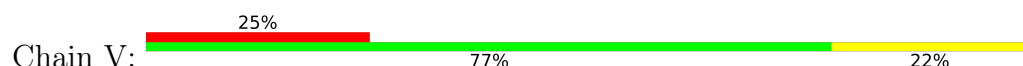
- Molecule 8: Cytochrome c oxidase subunit 6B1



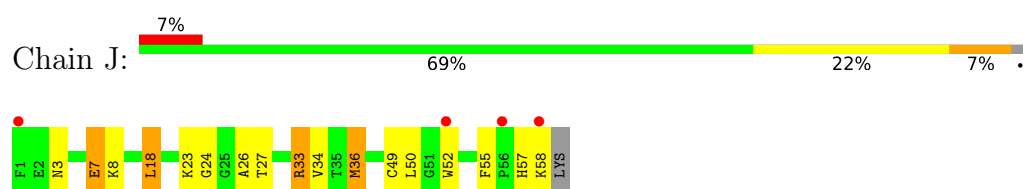
- Molecule 9: Cytochrome c oxidase subunit 6C



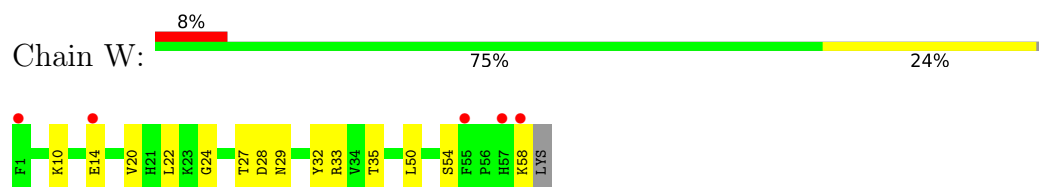
- Molecule 9: Cytochrome c oxidase subunit 6C



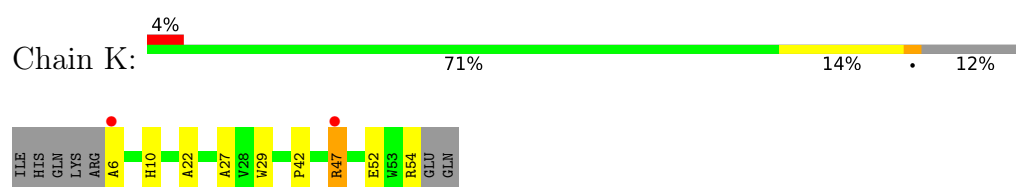
- Molecule 10: Cytochrome c oxidase subunit 7A1, mitochondrial



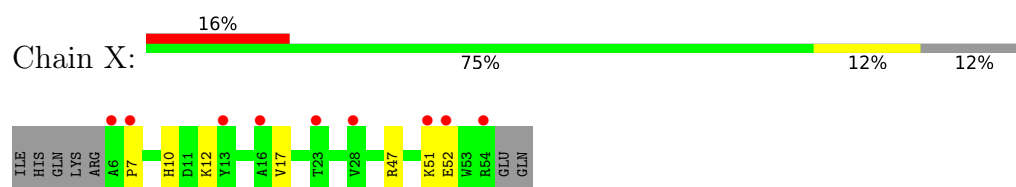
- Molecule 10: Cytochrome c oxidase subunit 7A1, mitochondrial



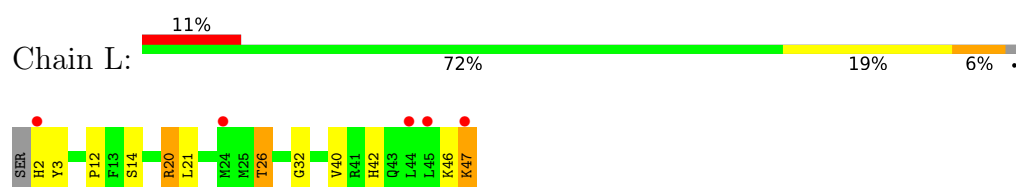
- Molecule 11: Cytochrome c oxidase subunit 7B, mitochondrial



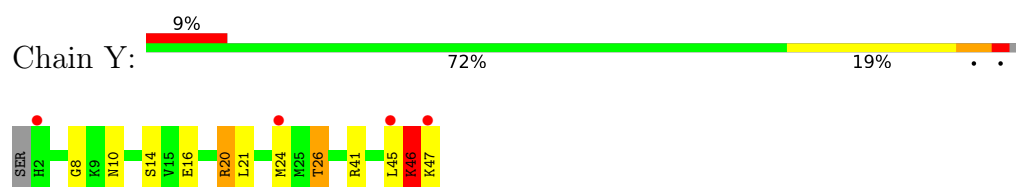
- Molecule 11: Cytochrome c oxidase subunit 7B, mitochondrial



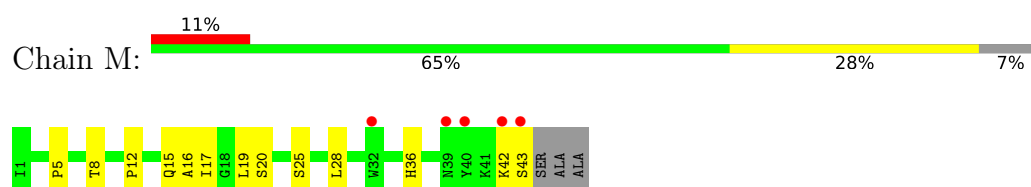
- Molecule 12: Cytochrome c oxidase subunit 7C, mitochondrial



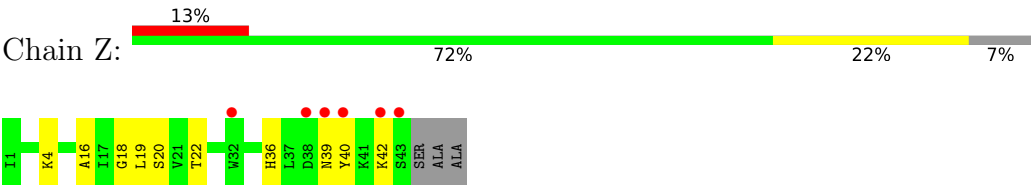
- Molecule 12: Cytochrome c oxidase subunit 7C, mitochondrial



- Molecule 13: Cytochrome c oxidase subunit 8B, mitochondrial



● Molecule 13: Cytochrome c oxidase subunit 8B, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	183.32Å 206.17Å 177.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 1.65 40.00 – 1.65	Depositor EDS
% Data completeness (in resolution range)	99.5 (40.00-1.65) 99.5 (40.00-1.65)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.96 (at 1.65Å)	Xtriage
Refinement program	REFMAC 5.8.0048	Depositor
R, R_{free}	0.175 , 0.197 0.177 , 0.198	Depositor DCC
R_{free} test set	39973 reflections (3.14%)	wwPDB-VP
Wilson B-factor (Å ²)	27.4	Xtriage
Anisotropy	0.590	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.005 for l,-k,h	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	33609	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.51% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, FME, EDO, DMU, CDL, PSC, SAC, CU, NA, MG, UNX, AZI, TGL, PEK, CUA, HEA, CHD, TPO, PGV

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	A	1.96	71/4322 (1.6%)	1.63	38/5897 (0.6%)
1	N	1.97	101/4308 (2.3%)	1.62	52/5878 (0.9%)
2	B	1.89	38/1937 (2.0%)	1.52	22/2637 (0.8%)
2	O	1.64	22/1908 (1.2%)	1.34	8/2597 (0.3%)
3	C	1.85	42/2272 (1.8%)	1.43	6/3102 (0.2%)
3	P	1.93	46/2272 (2.0%)	1.49	11/3102 (0.4%)
4	D	1.78	21/1268 (1.7%)	1.46	6/1709 (0.4%)
4	Q	1.41	8/1259 (0.6%)	1.25	5/1698 (0.3%)
5	E	1.85	13/871 (1.5%)	1.55	8/1182 (0.7%)
5	R	1.67	12/882 (1.4%)	1.39	2/1196 (0.2%)
6	F	1.80	13/795 (1.6%)	1.51	9/1079 (0.8%)
6	S	1.67	5/780 (0.6%)	1.45	2/1058 (0.2%)
7	G	1.95	14/702 (2.0%)	1.48	8/953 (0.8%)
7	T	1.95	19/702 (2.7%)	1.51	11/953 (1.2%)
8	H	1.59	7/682 (1.0%)	1.31	2/921 (0.2%)
8	U	1.29	1/682 (0.1%)	1.24	1/921 (0.1%)
9	I	1.57	5/605 (0.8%)	1.43	1/802 (0.1%)
9	V	1.33	1/605 (0.2%)	1.34	0/802
10	J	1.80	7/471 (1.5%)	1.44	3/636 (0.5%)
10	W	1.68	6/480 (1.2%)	1.38	2/648 (0.3%)
11	K	1.64	4/398 (1.0%)	1.29	1/546 (0.2%)
11	X	1.44	2/405 (0.5%)	1.20	0/556
12	L	1.77	4/393 (1.0%)	1.44	2/526 (0.4%)
12	Y	1.63	2/401 (0.5%)	1.29	2/536 (0.4%)
13	M	1.86	7/345 (2.0%)	1.37	0/470
13	Z	1.60	3/345 (0.9%)	1.28	2/470 (0.4%)
All	All	1.81	474/30090 (1.6%)	1.48	204/40875 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected

by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4
1	N	0	2
6	F	0	1
6	S	0	2
7	G	0	1
9	V	0	1
All	All	0	11

The worst 5 of 474 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	N	49	GLY	C-O	15.40	1.34	1.23
1	A	310	MET	SD-CE	-10.49	1.53	1.79
5	R	78	HIS	CE1-NE2	10.49	1.43	1.32
1	N	173	PRO	CA-CB	10.26	1.64	1.54
1	A	376	HIS	ND1-CE1	10.18	1.42	1.32

The worst 5 of 204 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	N	71	MET	CG-SD-CE	-12.09	74.31	100.90
5	E	90	ARG	NE-CZ-NH1	11.79	133.29	121.50
1	A	71	MET	CG-SD-CE	-11.38	75.86	100.90
1	A	112	LEU	CD1-CG-CD2	-10.41	87.89	110.80
7	G	7	ASP	N-CA-C	10.17	123.45	108.60

There are no chirality outliers.

5 of 11 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	240	HIS	Sidechain
1	A	296	GLY	Mainchain
1	A	379	TYR	Mainchain
1	A	38	ARG	Sidechain
6	F	93	PRO	Peptide

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	A	4193	0	4162	98	0
1	N	4179	0	4154	95	0
2	B	1899	0	1898	69	0
2	O	1870	0	1868	51	0
3	C	2185	0	2097	48	0
3	P	2185	0	2097	38	0
4	D	1233	0	1223	37	0
4	Q	1224	0	1211	21	0
5	E	852	0	845	4	0
5	R	863	0	857	6	0
6	F	778	0	754	26	0
6	S	763	0	742	24	0
7	G	686	0	652	32	0
7	T	686	0	651	21	0
8	H	662	0	623	19	0
8	U	662	0	623	12	0
9	I	601	0	613	20	0
9	V	601	0	613	15	0
10	J	460	0	459	9	0
10	W	469	0	464	4	0
11	K	384	0	366	3	0
11	X	391	0	374	3	0
12	L	380	0	380	16	0
12	Y	388	0	388	14	0
13	M	335	0	352	8	0
13	Z	335	0	352	8	0
14	A	180	0	162	18	0
14	N	180	0	162	22	0
15	A	1	0	0	1	0
15	N	1	0	0	0	0
16	A	1	0	0	0	0
16	N	1	0	0	0	0
17	A	1	0	0	0	0
17	N	1	0	0	0	0
18	A	9	0	0	7	0
18	N	9	0	0	2	0
19	A	126	0	220	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
19	D	63	0	110	13	0
19	N	63	0	110	7	0
19	Q	63	0	110	12	0
19	Y	63	0	110	20	0
20	A	102	0	152	13	0
20	C	102	0	152	5	0
20	N	102	0	152	10	0
20	P	102	0	152	3	0
21	A	36	0	52	8	0
21	B	16	0	24	2	0
21	C	4	0	6	0	0
21	D	8	0	12	9	0
21	E	12	0	18	0	0
21	F	12	0	18	0	0
21	G	8	0	12	3	0
21	L	4	0	6	0	0
21	M	4	0	6	0	0
21	N	36	0	54	1	0
21	O	4	0	6	0	0
21	P	8	0	12	0	0
21	R	4	0	6	0	0
21	S	12	0	18	0	0
21	T	4	0	6	0	0
21	W	4	0	6	0	0
21	Y	4	0	6	0	0
22	B	29	0	39	0	0
22	C	58	0	78	6	0
22	G	29	0	39	1	0
22	J	29	0	38	1	0
22	P	58	0	78	6	0
22	W	29	0	38	1	0
23	B	2	0	0	0	0
23	O	2	0	0	0	0
24	B	52	0	80	15	0
24	N	52	0	80	16	0
25	C	99	0	126	15	0
25	M	33	0	42	0	0
25	P	99	0	126	15	0
25	Z	33	0	42	0	0
26	C	1	0	0	0	0
26	P	1	0	0	1	0
27	C	100	0	156	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
27	N	100	0	156	29	0
27	P	100	0	156	24	0
27	T	100	0	156	16	0
28	C	106	0	154	16	0
28	G	106	0	154	8	0
28	P	53	0	77	6	0
28	T	53	0	77	3	0
29	F	1	0	0	0	0
29	S	1	0	0	0	0
30	A	235	0	0	32	0
30	B	162	0	0	25	2
30	C	110	0	0	3	0
30	D	118	0	0	12	1
30	E	93	0	0	1	0
30	F	94	0	0	5	0
30	G	42	0	0	5	0
30	H	42	0	0	1	0
30	I	27	0	0	6	1
30	J	23	0	0	2	0
30	K	26	0	0	1	0
30	L	38	0	0	4	1
30	M	25	0	0	7	1
30	N	203	0	0	13	0
30	O	100	0	0	1	0
30	P	95	0	0	3	0
30	Q	29	0	0	5	0
30	R	40	0	0	1	0
30	S	45	0	0	3	0
30	T	35	0	0	3	0
30	U	29	0	0	0	0
30	V	16	0	0	3	0
30	W	7	0	0	0	0
30	X	11	0	0	0	0
30	Y	12	0	0	0	0
30	Z	12	0	0	0	0
All	All	33609	0	32570	796	3

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 12.

The worst 5 of 796 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)
1:A:74:MET:CG	1:A:74:MET:CB	1.75	1.61
6:S:43:LYS:H	6:S:43:LYS:CD	1.08	1.49
20:N:609:PGV:H011	20:N:609:PGV:C2	1.38	1.33
24:B:303:PSC:O02	9:I:14:ALA:HB2	1.16	1.31
30:A:709:HOH:O	19:D:201:TGL:HG11	1.23	1.29

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:B:491:HOH:O	30:D:341:HOH:O[2_584]	1.22	0.98
30:I:126:HOH:O	30:M:216:HOH:O[2_584]	1.92	0.28
30:B:541:HOH:O	30:L:228:HOH:O[2_584]	2.14	0.06

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 X-ray entries followed by that with respect to entries of similar resolution.

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	A	534/514 (104%)	516 (97%)	18 (3%)	0	100	100
1	N	532/514 (104%)	515 (97%)	17 (3%)	0	100	100
2	B	234/227 (103%)	227 (97%)	5 (2%)	2 (1%)	14	3
2	O	230/227 (101%)	223 (97%)	7 (3%)	0	100	100
3	C	266/261 (102%)	262 (98%)	4 (2%)	0	100	100
3	P	266/261 (102%)	261 (98%)	5 (2%)	0	100	100
4	D	146/147 (99%)	142 (97%)	4 (3%)	0	100	100
4	Q	145/147 (99%)	138 (95%)	4 (3%)	3 (2%)	5	0
5	E	103/109 (94%)	103 (100%)	0	0	100	100
5	R	104/109 (95%)	103 (99%)	0	1 (1%)	12	2
6	F	100/98 (102%)	96 (96%)	1 (1%)	3 (3%)	3	0

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
6	S	98/98 (100%)	91 (93%)	3 (3%)	4 (4%)	2	0
7	G	82/85 (96%)	68 (83%)	9 (11%)	5 (6%)	1	0
7	T	82/85 (96%)	70 (85%)	8 (10%)	4 (5%)	1	0
8	H	77/85 (91%)	70 (91%)	4 (5%)	3 (4%)	2	0
8	U	77/85 (91%)	68 (88%)	5 (6%)	4 (5%)	1	0
9	I	71/73 (97%)	70 (99%)	1 (1%)	0	100	100
9	V	71/73 (97%)	70 (99%)	1 (1%)	0	100	100
10	J	56/59 (95%)	56 (100%)	0	0	100	100
10	W	57/59 (97%)	56 (98%)	1 (2%)	0	100	100
11	K	47/56 (84%)	45 (96%)	2 (4%)	0	100	100
11	X	48/56 (86%)	46 (96%)	2 (4%)	0	100	100
12	L	44/47 (94%)	41 (93%)	3 (7%)	0	100	100
12	Y	45/47 (96%)	43 (96%)	1 (2%)	1 (2%)	5	0
13	M	41/46 (89%)	38 (93%)	2 (5%)	1 (2%)	4	0
13	Z	41/46 (89%)	39 (95%)	2 (5%)	0	100	100
All	All	3597/3614 (100%)	3457 (96%)	109 (3%)	31 (1%)	14	3

5 of 31 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
6	F	94	HIS
7	G	2	SER
7	G	3	ALA
7	G	4	ALA
7	G	5	LYS

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 X-ray entries followed by that with respect to entries of similar resolution.

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	A	447/426 (105%)	438 (98%)	9 (2%)	48	26

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	N	445/426 (104%)	434 (98%)	11 (2%)	42	18
2	B	219/210 (104%)	208 (95%)	11 (5%)	22	5
2	O	215/210 (102%)	207 (96%)	8 (4%)	30	8
3	C	233/226 (103%)	228 (98%)	5 (2%)	47	24
3	P	233/226 (103%)	230 (99%)	3 (1%)	61	42
4	D	132/129 (102%)	129 (98%)	3 (2%)	44	21
4	Q	131/129 (102%)	126 (96%)	5 (4%)	29	7
5	E	92/95 (97%)	89 (97%)	3 (3%)	33	11
5	R	93/95 (98%)	88 (95%)	5 (5%)	20	4
6	F	85/81 (105%)	81 (95%)	4 (5%)	23	5
6	S	83/81 (102%)	73 (88%)	10 (12%)	5	1
7	G	68/68 (100%)	62 (91%)	6 (9%)	9	1
7	T	68/68 (100%)	63 (93%)	5 (7%)	13	2
8	H	71/75 (95%)	68 (96%)	3 (4%)	26	6
8	U	71/75 (95%)	65 (92%)	6 (8%)	10	1
9	I	57/57 (100%)	55 (96%)	2 (4%)	32	10
9	V	57/57 (100%)	54 (95%)	3 (5%)	20	4
10	J	49/50 (98%)	48 (98%)	1 (2%)	48	26
10	W	50/50 (100%)	48 (96%)	2 (4%)	28	7
11	K	39/46 (85%)	38 (97%)	1 (3%)	40	17
11	X	40/46 (87%)	38 (95%)	2 (5%)	22	5
12	L	39/40 (98%)	38 (97%)	1 (3%)	40	17
12	Y	40/40 (100%)	38 (95%)	2 (5%)	22	5
13	M	37/38 (97%)	37 (100%)	0	100	100
13	Z	37/38 (97%)	37 (100%)	0	100	100
All	All	3131/3082 (102%)	3020 (96%)	111 (4%)	32	10

5 of 111 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	N	369	ASP
12	Y	26	THR
4	Q	6	VAL

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Mol	Chain	Res	Type
12	Y	16	GLU
8	U	60	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 57 such sidechains are listed below:

Mol	Chain	Res	Type
1	N	503	HIS
11	X	35	GLN
4	Q	37	GLN
10	W	57	HIS
8	U	25	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

8 non-standard protein/DNA/RNA residues are 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$
1	FME	A	1	1	8,9,10	0.94	1 (12%)	8,9,11	1.60	3 (37%)
9	SAC	V	1	9	7,8,9	1.56	1 (14%)	7,9,11	1.00	1 (14%)
7	TPO	T	11	7	8,10,11	1.80	2 (25%)	10,14,16	1.29	1 (10%)
2	FME	O	1	2	8,9,10	1.23	1 (12%)	8,9,11	1.78	2 (25%)
9	SAC	I	1	9	7,8,9	1.43	1 (14%)	7,9,11	2.11	3 (42%)
1	FME	N	1	1	8,9,10	1.15	1 (12%)	8,9,11	1.23	1 (12%)
2	FME	B	1	2	8,9,10	1.92	4 (50%)	8,9,11	1.64	2 (25%)
7	TPO	G	11	7	8,10,11	2.12	3 (37%)	10,14,16	0.92	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
1	FME	A	1	1	-	2/7/9/11	-
9	SAC	V	1	9	-	7/7/8/10	-
7	TPO	T	11	7	-	6/9/11/13	-
2	FME	O	1	2	-	0/7/9/11	-
9	SAC	I	1	9	-	4/7/8/10	-
1	FME	N	1	1	-	3/7/9/11	-
2	FME	B	1	2	-	1/7/9/11	-
7	TPO	G	11	7	-	4/9/11/13	-

The worst 5 of 14 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	V	1	SAC	CA-N	3.96	1.52	1.46
7	G	11	TPO	P-O1P	3.44	1.61	1.50
9	I	1	SAC	CA-N	3.24	1.51	1.46
7	T	11	TPO	P-O1P	3.22	1.60	1.50
2	B	1	FME	CB-CA	2.99	1.59	1.53

The worst 5 of 13 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	O	1	FME	CA-N-CN	3.36	127.98	122.82
9	I	1	SAC	C-CA-N	3.35	115.97	109.50
9	I	1	SAC	OG-CB-CA	-3.33	101.96	111.13
2	B	1	FME	C-CA-N	-2.87	103.96	109.50
2	B	1	FME	CG-CB-CA	-2.72	104.55	112.87

There are no chirality outliers.

5 of 27 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	1	FME	N-CA-CB-CG
7	G	11	TPO	N-CA-CB-OG1
7	G	11	TPO	O-C-CA-CB
7	G	11	TPO	CA-CB-OG1-P
9	I	1	SAC	C-CA-CB-OG

There are no ring outliers.

6 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	1	FME	1	0
9	V	1	SAC	1	0
7	T	11	TPO	1	0
2	O	1	FME	2	0
2	B	1	FME	4	0
7	G	11	TPO	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 111 ligands modelled in this entry, 8 are monoatomic and 2 are unknown - leaving 101 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	$\# Z > 2$
21	EDO	M	102	-	3,3,3	0.22	0	2,2,2	1.05	0
20	PGV	C	308	-	50,50,50	1.28	2 (4%)	53,56,56	1.45	7 (13%)
21	EDO	N	613	-	3,3,3	1.62	1 (33%)	2,2,2	0.54	0
21	EDO	D	202	-	3,3,3	0.94	0	2,2,2	0.39	0
22	CHD	W	101	-	32,32,32	1.28	3 (9%)	51,51,51	2.63	27 (52%)
21	EDO	E	203	-	3,3,3	0.74	0	2,2,2	0.55	0
21	EDO	A	617	-	3,3,3	1.36	0	2,2,2	0.40	0
21	EDO	B	307	-	3,3,3	0.61	0	2,2,2	0.05	0
21	EDO	B	306	-	3,3,3	0.55	0	2,2,2	0.34	0
25	DMU	P	310	-	34,34,34	1.11	2 (5%)	45,45,45	1.83	11 (24%)
20	PGV	N	610	-	50,50,50	1.18	6 (12%)	53,56,56	1.51	9 (16%)
25	DMU	Z	101	-	34,34,34	0.80	1 (2%)	45,45,45	1.22	3 (6%)
21	EDO	N	615	-	3,3,3	0.69	0	2,2,2	1.25	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
14	HEA	A	601	1	67,67,67	1.91	22 (32%)	81,103,103	2.60	32 (39%)
23	CUA	B	302	2,30	0,1,1	-	-	-		
21	EDO	B	305	-	3,3,3	1.55	0	2,2,2	0.23	0
22	CHD	G	102	-	32,32,32	2.01	10 (31%)	51,51,51	1.98	14 (27%)
21	EDO	S	102	-	3,3,3	0.99	0	2,2,2	0.43	0
24	PSC	N	612	-	51,51,51	1.29	3 (5%)	57,59,59	1.20	4 (7%)
14	HEA	N	603[A]	18,1	67,67,67	1.98	25 (37%)	81,103,103	2.04	30 (37%)
21	EDO	A	615	-	3,3,3	0.63	0	2,2,2	1.79	1 (50%)
21	EDO	A	612	-	3,3,3	0.39	0	2,2,2	1.41	0
18	AZI	A	606[B]	14	2,2,2	1.27	0	0,1,1	-	-
21	EDO	F	104	-	3,3,3	0.67	0	2,2,2	0.38	0
25	DMU	C	302	-	34,34,34	0.77	1 (2%)	45,45,45	1.48	8 (17%)
21	EDO	N	620	-	3,3,3	0.55	0	2,2,2	0.11	0
21	EDO	N	616	-	3,3,3	1.00	0	2,2,2	0.23	0
14	HEA	N	603[B]	18,1	67,67,67	1.65	14 (20%)	81,103,103	2.23	29 (35%)
28	PEK	G	101	-	52,52,52	1.12	6 (11%)	55,57,57	1.46	8 (14%)
21	EDO	A	620	-	3,3,3	0.45	0	2,2,2	0.85	0
18	AZI	A	607[A]	14,15	2,2,2	3.47	2 (100%)	0,1,1	-	-
21	EDO	O	302	-	3,3,3	0.61	0	2,2,2	0.61	0
20	PGV	C	304	-	50,50,50	1.01	3 (6%)	53,56,56	1.13	4 (7%)
21	EDO	P	312	-	3,3,3	0.73	0	2,2,2	1.60	1 (50%)
18	AZI	N	607[B]	14	2,2,2	1.89	1 (50%)	0,1,1	-	-
22	CHD	B	301	-	32,32,32	1.82	10 (31%)	51,51,51	2.26	22 (43%)
21	EDO	A	613	-	3,3,3	0.55	0	2,2,2	1.29	0
18	AZI	A	607[B]	15	2,2,2	7.06	2 (100%)	0,1,1	-	-
21	EDO	P	311	-	3,3,3	0.81	0	2,2,2	0.48	0
21	EDO	B	304	-	3,3,3	0.16	0	2,2,2	1.38	0
22	CHD	C	301	-	32,32,32	1.81	7 (21%)	51,51,51	2.29	19 (37%)
21	EDO	N	618	-	3,3,3	0.62	0	2,2,2	0.68	0
23	CUA	O	301	2	0,1,1	-	-	-		
21	EDO	S	103	-	3,3,3	1.23	0	2,2,2	0.53	0
21	EDO	N	614	-	3,3,3	0.44	0	2,2,2	1.01	0
21	EDO	E	202	-	3,3,3	0.92	0	2,2,2	0.61	0
25	DMU	C	311	-	34,34,34	1.05	1 (2%)	45,45,45	2.38	13 (28%)
14	HEA	N	602	1	67,67,67	1.87	25 (37%)	81,103,103	2.20	26 (32%)
25	DMU	P	307	-	34,34,34	0.95	1 (2%)	45,45,45	1.49	9 (20%)
20	PGV	N	609	-	50,50,50	1.21	2 (4%)	53,56,56	1.25	6 (11%)
22	CHD	P	306	-	32,32,32	1.52	6 (18%)	51,51,51	3.43	22 (43%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
22	CHD	J	101	-	32,32,32	0.92	0	51,51,51	2.28	19 (37%)
21	EDO	T	103	-	3,3,3	0.94	0	2,2,2	0.82	0
21	EDO	A	619	-	3,3,3	0.24	0	2,2,2	0.80	0
21	EDO	D	203	-	3,3,3	0.54	0	2,2,2	0.35	0
25	DMU	P	309	-	34,34,34	0.81	1 (2%)	45,45,45	2.18	12 (26%)
20	PGV	A	609	-	50,50,50	1.22	6 (12%)	53,56,56	1.33	6 (11%)
21	EDO	C	312	-	3,3,3	0.59	0	2,2,2	0.60	0
19	TGL	N	611	-	62,62,62	1.13	4 (6%)	65,65,65	1.72	10 (15%)
22	CHD	P	301	-	32,32,32	1.62	9 (28%)	51,51,51	2.04	15 (29%)
21	EDO	A	614	-	3,3,3	1.60	1 (33%)	2,2,2	0.80	0
21	EDO	W	102	-	3,3,3	0.45	0	2,2,2	0.67	0
21	EDO	A	616	-	3,3,3	2.67	1 (33%)	2,2,2	5.31	1 (50%)
21	EDO	E	201	-	3,3,3	0.68	0	2,2,2	0.78	0
19	TGL	A	608	-	62,62,62	1.33	6 (9%)	65,65,65	2.26	10 (15%)
21	EDO	Y	102	-	3,3,3	0.49	0	2,2,2	0.64	0
21	EDO	F	103	-	3,3,3	1.06	0	2,2,2	0.16	0
20	PGV	P	304	-	50,50,50	1.10	4 (8%)	53,56,56	1.36	6 (11%)
21	EDO	N	619	-	3,3,3	1.22	0	2,2,2	0.24	0
19	TGL	A	611	-	62,62,62	1.29	3 (4%)	65,65,65	1.84	13 (20%)
21	EDO	G	105	-	3,3,3	0.56	0	2,2,2	0.24	0
20	PGV	A	610	-	50,50,50	1.42	5 (10%)	53,56,56	1.92	7 (13%)
28	PEK	C	307	-	52,52,52	1.50	4 (7%)	55,57,57	1.52	9 (16%)
14	HEA	A	602[A]	18,1	67,67,67	1.93	22 (32%)	81,103,103	2.19	30 (37%)
28	PEK	G	103	-	52,52,52	1.11	2 (3%)	55,57,57	1.20	4 (7%)
19	TGL	Q	201	-	62,62,62	1.52	4 (6%)	65,65,65	1.50	7 (10%)
25	DMU	C	310	-	34,34,34	0.91	1 (2%)	45,45,45	2.55	15 (33%)
21	EDO	G	104	-	3,3,3	0.92	0	2,2,2	0.64	0
21	EDO	S	104	-	3,3,3	0.61	0	2,2,2	1.63	1 (50%)
21	EDO	F	102	-	3,3,3	0.91	0	2,2,2	0.23	0
27	CDL	N	601	-	99,99,99	1.61	16 (16%)	105,111,111	1.53	18 (17%)
24	PSC	B	303	-	51,51,51	1.29	4 (7%)	57,59,59	1.39	5 (8%)
14	HEA	A	602[B]	18,1	67,67,67	1.87	18 (26%)	81,103,103	2.09	23 (28%)
18	AZI	N	608[A]	14,15	2,2,2	4.14	2 (100%)	0,1,1	-	-
21	EDO	A	618	-	3,3,3	0.40	0	2,2,2	1.03	0
25	DMU	M	101	-	34,34,34	0.91	1 (2%)	45,45,45	1.45	6 (13%)
28	PEK	T	101	-	52,52,52	1.39	7 (13%)	55,57,57	2.16	8 (14%)
21	EDO	L	101	-	3,3,3	0.73	0	2,2,2	0.86	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
28	PEK	C	309	-	52,52,52	1.19	2 (3%)	55,57,57	1.36	6 (10%)
27	CDL	P	305	-	99,99,99	1.61	18 (18%)	105,111,111	1.65	18 (17%)
21	EDO	R	201	-	3,3,3	0.84	0	2,2,2	0.65	0
27	CDL	C	305	-	99,99,99	1.58	18 (18%)	105,111,111	1.54	16 (15%)
18	AZI	N	608[B]	15	2,2,2	3.59	2 (100%)	0,1,1	-	-
19	TGL	Y	101	-	62,62,62	1.50	5 (8%)	65,65,65	1.83	10 (15%)
28	PEK	P	308	-	52,52,52	1.26	2 (3%)	55,57,57	1.38	5 (9%)
20	PGV	P	302	-	50,50,50	1.17	2 (4%)	53,56,56	1.35	5 (9%)
27	CDL	T	102	-	99,99,99	1.52	13 (13%)	105,111,111	1.38	17 (16%)
22	CHD	C	306	-	32,32,32	1.33	5 (15%)	51,51,51	3.26	18 (35%)
21	EDO	N	621	-	3,3,3	0.75	0	2,2,2	1.20	0
19	TGL	D	201	-	62,62,62	2.01	5 (8%)	65,65,65	2.60	11 (16%)
21	EDO	N	617	-	3,3,3	0.61	0	2,2,2	0.82	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
21	EDO	M	102	-	-	0/1/1/1	-
20	PGV	C	308	-	-	23/55/55/55	-
21	EDO	N	613	-	-	0/1/1/1	-
21	EDO	D	202	-	-	1/1/1/1	-
22	CHD	W	101	-	-	8/9/74/74	0/4/4/4
21	EDO	E	203	-	-	0/1/1/1	-
21	EDO	A	617	-	-	0/1/1/1	-
21	EDO	B	307	-	-	0/1/1/1	-
21	EDO	B	306	-	-	0/1/1/1	-
25	DMU	P	310	-	-	10/19/59/59	0/2/2/2
20	PGV	N	610	-	-	11/55/55/55	-
25	DMU	Z	101	-	-	7/19/59/59	0/2/2/2
21	EDO	N	615	-	-	0/1/1/1	-
14	HEA	A	601	1	-	4/36/76/76	-
21	EDO	B	305	-	-	1/1/1/1	-
22	CHD	G	102	-	-	2/9/74/74	0/4/4/4
21	EDO	S	102	-	-	0/1/1/1	-
24	PSC	N	612	-	-	27/55/55/55	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
14	HEA	N	603[A]	18,1	-	6/36/76/76	-
21	EDO	A	615	-	-	1/1/1/1	-
21	EDO	A	612	-	-	1/1/1/1	-
21	EDO	F	104	-	-	0/1/1/1	-
25	DMU	C	302	-	-	9/19/59/59	0/2/2/2
21	EDO	N	620	-	-	0/1/1/1	-
21	EDO	N	616	-	-	0/1/1/1	-
14	HEA	N	603[B]	18,1	-	7/36/76/76	-
28	PEK	G	101	-	-	18/56/56/56	-
21	EDO	A	620	-	-	0/1/1/1	-
21	EDO	O	302	-	-	0/1/1/1	-
20	PGV	C	304	-	-	17/55/55/55	-
21	EDO	P	312	-	-	0/1/1/1	-
22	CHD	B	301	-	-	2/9/74/74	0/4/4/4
21	EDO	A	613	-	-	1/1/1/1	-
21	EDO	P	311	-	-	0/1/1/1	-
21	EDO	B	304	-	-	1/1/1/1	-
22	CHD	C	301	-	-	2/9/74/74	0/4/4/4
21	EDO	N	618	-	-	1/1/1/1	-
21	EDO	S	103	-	-	0/1/1/1	-
21	EDO	N	614	-	-	0/1/1/1	-
21	EDO	E	202	-	-	1/1/1/1	-
25	DMU	C	311	-	-	8/19/59/59	0/2/2/2
14	HEA	N	602	1	-	4/36/76/76	-
25	DMU	P	307	-	-	5/19/59/59	0/2/2/2
20	PGV	N	609	-	-	34/55/55/55	-
22	CHD	P	306	-	-	5/9/74/74	0/4/4/4
22	CHD	J	101	-	-	5/9/74/74	0/4/4/4
21	EDO	T	103	-	-	0/1/1/1	-
21	EDO	A	619	-	-	1/1/1/1	-
21	EDO	D	203	-	-	0/1/1/1	-
25	DMU	P	309	-	-	7/19/59/59	0/2/2/2
20	PGV	A	609	-	-	8/55/55/55	-
21	EDO	C	312	-	-	0/1/1/1	-
19	TGL	N	611	-	-	38/65/65/65	-
22	CHD	P	301	-	-	3/9/74/74	0/4/4/4
21	EDO	A	614	-	-	0/1/1/1	-
21	EDO	W	102	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
21	EDO	A	616	-	-	1/1/1/1	-
21	EDO	E	201	-	-	0/1/1/1	-
19	TGL	A	608	-	-	36/65/65/65	-
21	EDO	Y	102	-	-	1/1/1/1	-
21	EDO	F	103	-	-	0/1/1/1	-
20	PGV	P	304	-	-	13/55/55/55	-
21	EDO	N	619	-	-	0/1/1/1	-
19	TGL	A	611	-	-	38/65/65/65	-
21	EDO	G	105	-	-	0/1/1/1	-
20	PGV	A	610	-	-	29/55/55/55	-
28	PEK	C	307	-	-	30/56/56/56	-
14	HEA	A	602[A]	18,1	-	7/36/76/76	-
28	PEK	G	103	-	-	29/56/56/56	-
19	TGL	Q	201	-	-	31/65/65/65	-
25	DMU	C	310	-	-	7/19/59/59	0/2/2/2
21	EDO	G	104	-	-	0/1/1/1	-
21	EDO	S	104	-	-	0/1/1/1	-
21	EDO	F	102	-	-	0/1/1/1	-
27	CDL	N	601	-	-	58/110/110/110	-
24	PSC	B	303	-	-	33/55/55/55	-
14	HEA	A	602[B]	18,1	-	6/36/76/76	-
21	EDO	A	618	-	-	1/1/1/1	-
25	DMU	M	101	-	-	4/19/59/59	0/2/2/2
28	PEK	T	101	-	-	22/56/56/56	-
21	EDO	L	101	-	-	1/1/1/1	-
28	PEK	C	309	-	-	25/56/56/56	-
27	CDL	P	305	-	-	54/110/110/110	-
21	EDO	R	201	-	-	0/1/1/1	-
27	CDL	C	305	-	-	53/110/110/110	-
19	TGL	Y	101	-	-	34/65/65/65	-
28	PEK	P	308	-	-	25/56/56/56	-
20	PGV	P	302	-	-	23/55/55/55	-
27	CDL	T	102	-	-	48/110/110/110	-
22	CHD	C	306	-	-	5/9/74/74	0/4/4/4
21	EDO	N	621	-	-	1/1/1/1	-
19	TGL	D	201	-	-	35/65/65/65	-
21	EDO	N	617	-	-	0/1/1/1	-

The worst 5 of 349 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
19	D	201	TGL	OB1-CB1	9.96	1.52	1.22
18	A	607[B]	AZI	N1-N2	7.95	1.39	1.23
19	D	201	TGL	OG2-CB1	7.59	1.55	1.34
19	Y	101	TGL	OG2-CB1	6.72	1.53	1.34
19	Q	201	TGL	OG2-CB1	6.51	1.52	1.34

The worst 5 of 636 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
22	C	306	CHD	C23-C22-C20	-17.52	81.74	114.46
22	P	306	CHD	C23-C22-C20	-17.39	81.97	114.46
28	T	101	PEK	C2-C3-C4	12.47	140.59	113.35
19	D	201	TGL	OG2-CB1-CB2	-12.27	84.94	111.48
25	C	311	DMU	O16-C6-C1	11.41	125.59	108.27

There are no chirality outliers.

5 of 929 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
14	A	602[B]	HEA	C2A-C3A-CMA-OMA
14	A	602[B]	HEA	C4D-C3D-CAD-CBD
14	N	603[A]	HEA	C2A-C3A-CMA-OMA
19	A	611	TGL	CB2-CB1-OG2-CG2
19	A	611	TGL	OB1-CB1-OG2-CG2

There are no ring outliers.

56 monomers are involved in 349 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
20	C	308	PGV	1	0
21	D	202	EDO	9	0
22	W	101	CHD	1	0
25	P	310	DMU	2	0
20	N	610	PGV	1	0
14	A	601	HEA	10	0
22	G	102	CHD	1	0
24	N	612	PSC	16	0
14	N	603[A]	HEA	5	0
25	C	302	DMU	12	0
14	N	603[B]	HEA	6	0

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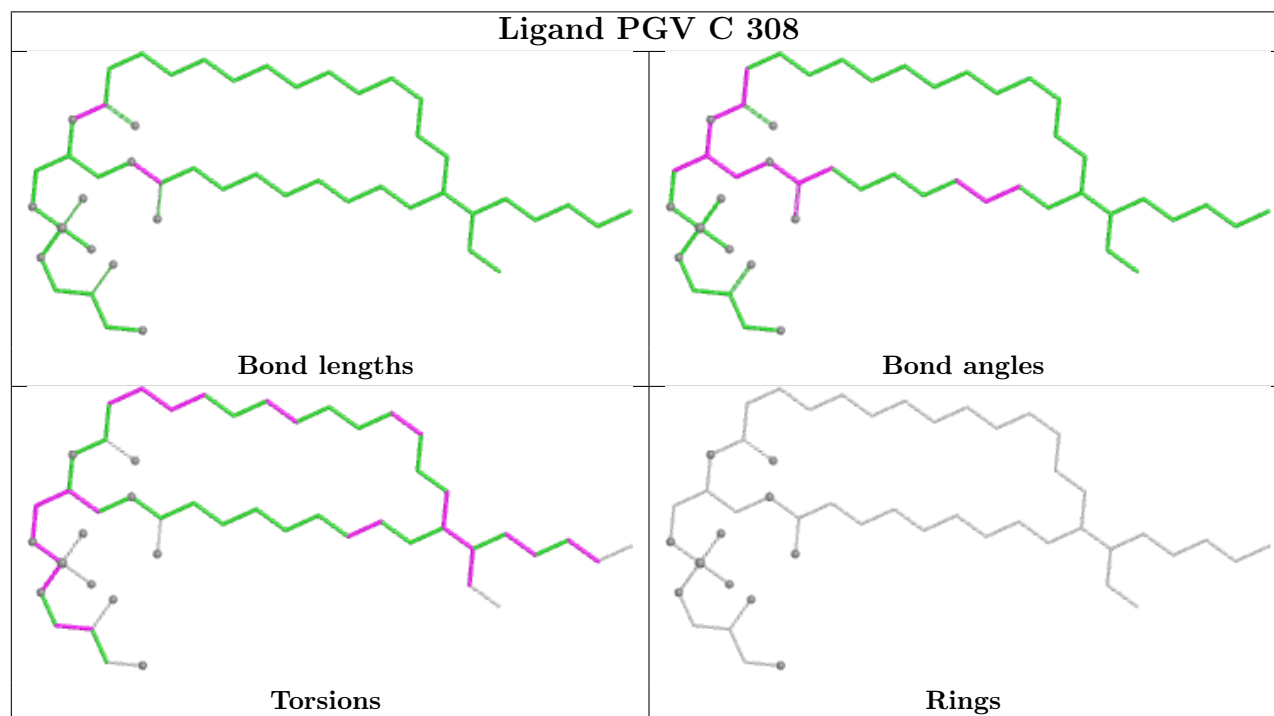
Mol	Chain	Res	Type	Clashes	Symm-Clashes
28	G	101	PEK	7	0
21	A	620	EDO	1	0
18	A	607[A]	AZI	2	0
20	C	304	PGV	4	0
18	A	607[B]	AZI	5	0
21	B	304	EDO	2	0
22	C	301	CHD	1	0
14	N	602	HEA	11	0
25	P	307	DMU	11	0
20	N	609	PGV	9	0
22	P	306	CHD	5	0
22	J	101	CHD	1	0
21	A	619	EDO	3	0
25	P	309	DMU	3	0
20	A	609	PGV	4	0
19	N	611	TGL	7	0
22	P	301	CHD	1	0
21	A	616	EDO	2	0
19	A	608	TGL	5	0
20	P	304	PGV	2	0
19	A	611	TGL	6	0
21	G	105	EDO	3	0
20	A	610	PGV	9	0
28	C	307	PEK	12	0
14	A	602[A]	HEA	3	0
28	G	103	PEK	1	0
19	Q	201	TGL	12	0
25	C	310	DMU	3	0
27	N	601	CDL	29	0
24	B	303	PSC	15	0
14	A	602[B]	HEA	5	0
18	N	608[A]	AZI	1	0
21	A	618	EDO	2	0
28	T	101	PEK	3	0
28	C	309	PEK	4	0
27	P	305	CDL	24	0
27	C	305	CDL	16	0
18	N	608[B]	AZI	1	0
19	Y	101	TGL	20	0
28	P	308	PEK	6	0
20	P	302	PGV	1	0
27	T	102	CDL	16	0

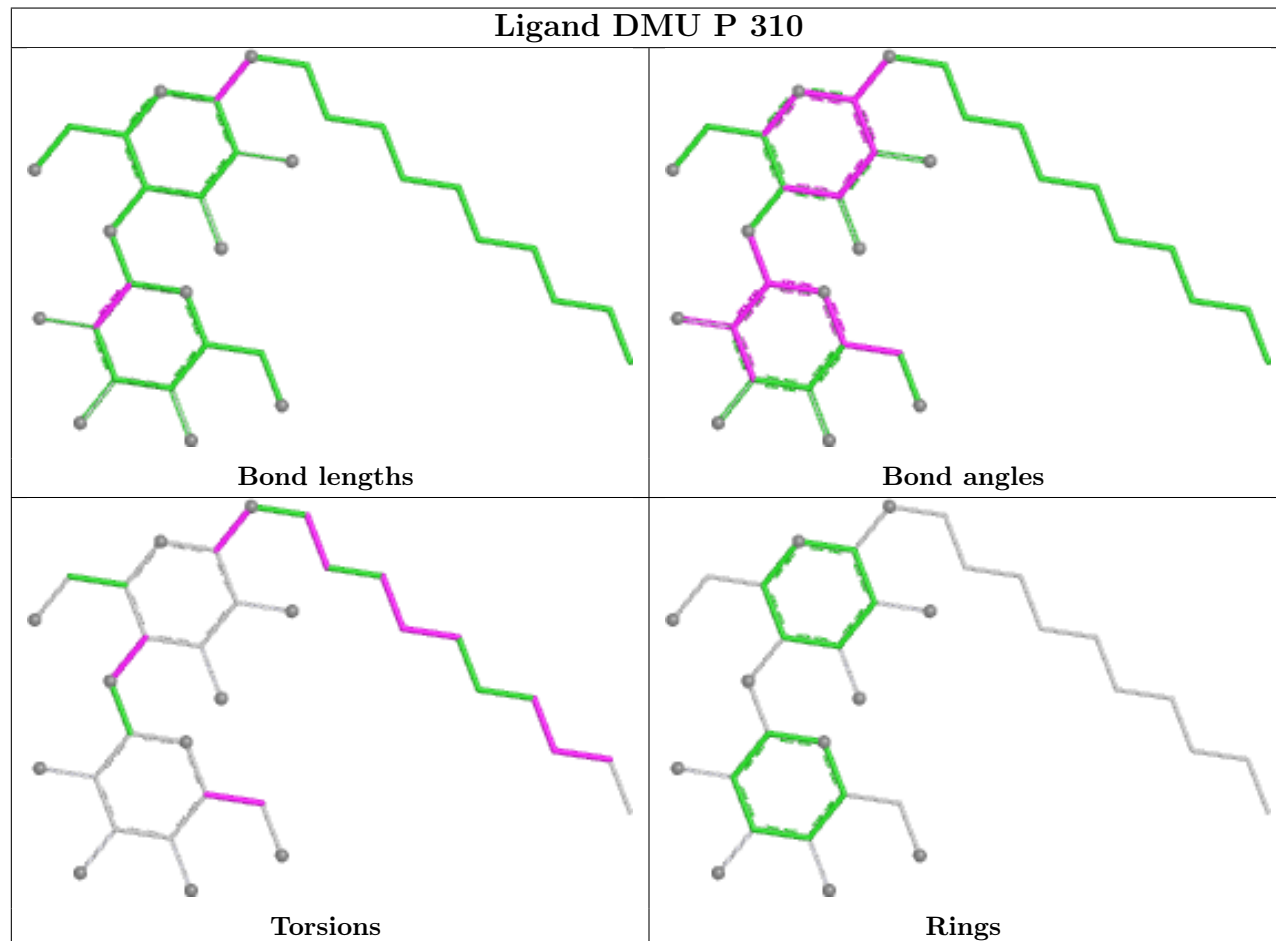
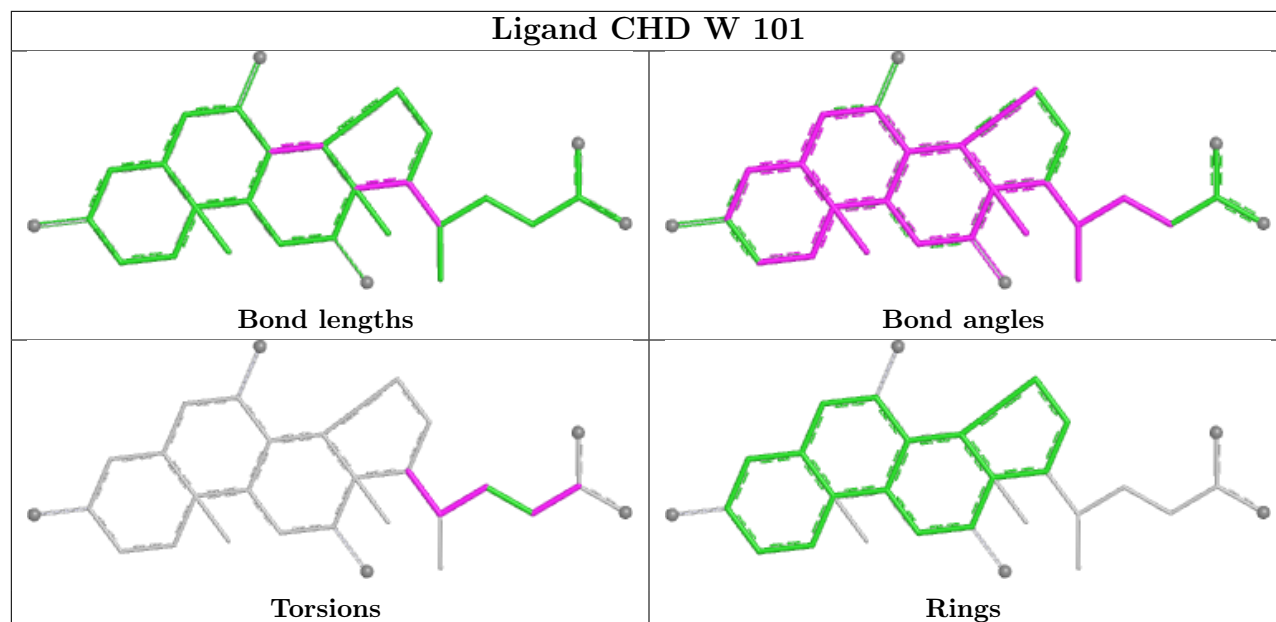
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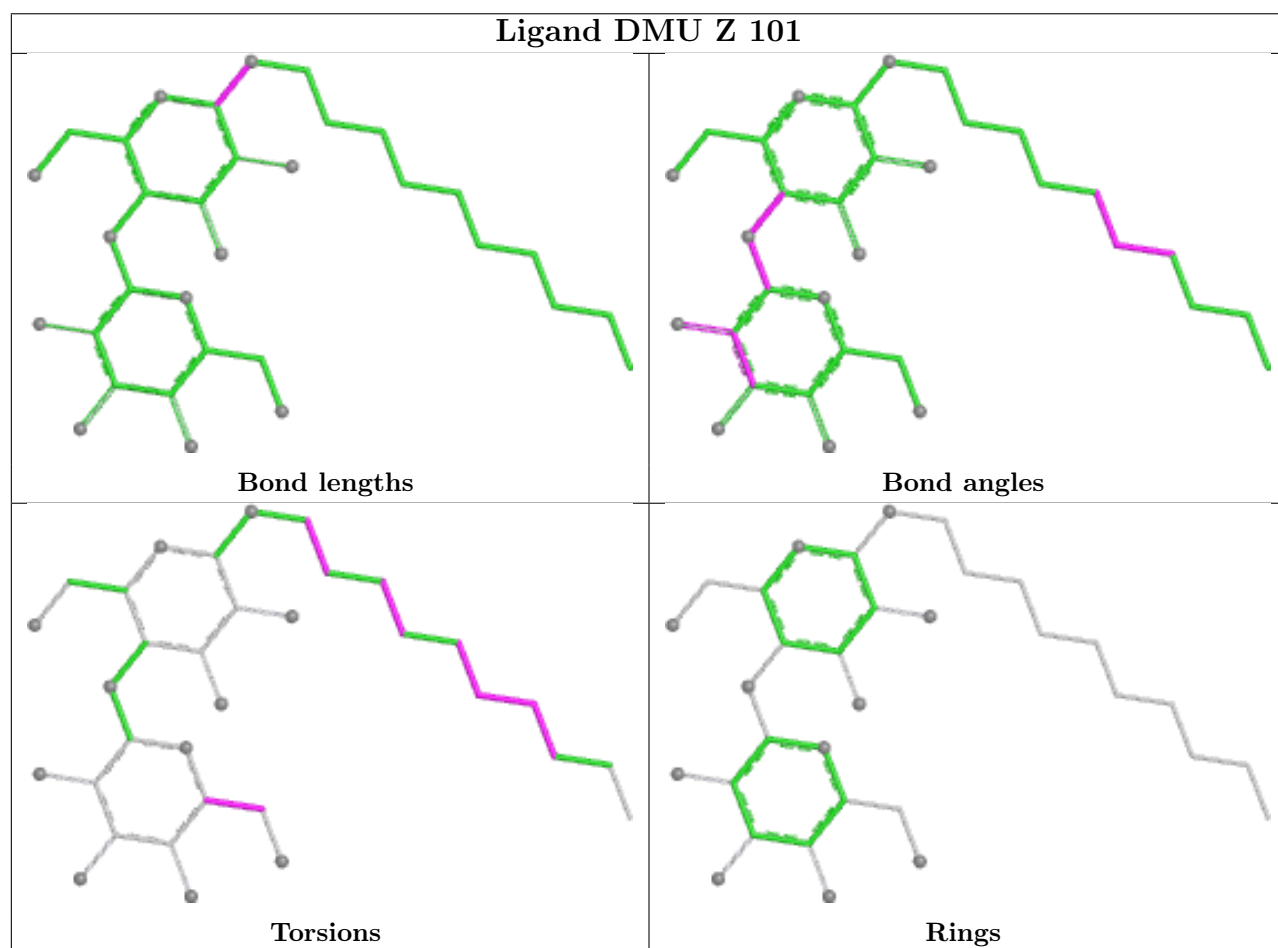
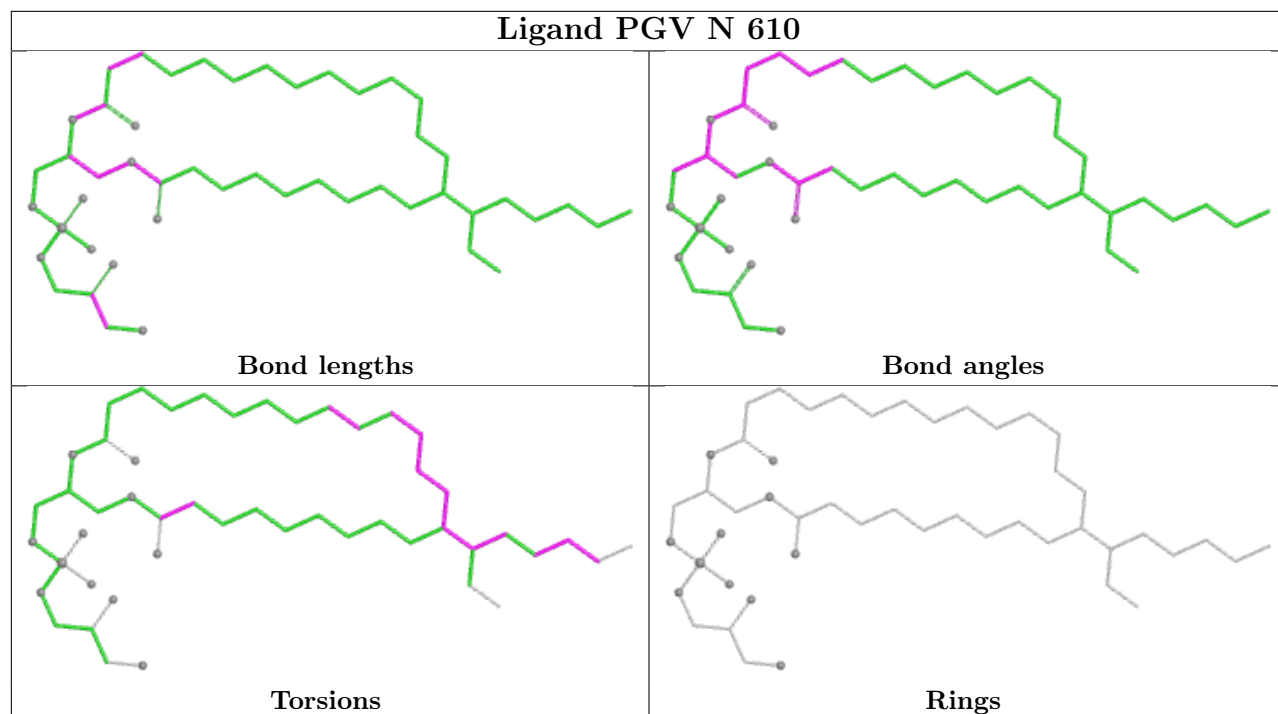
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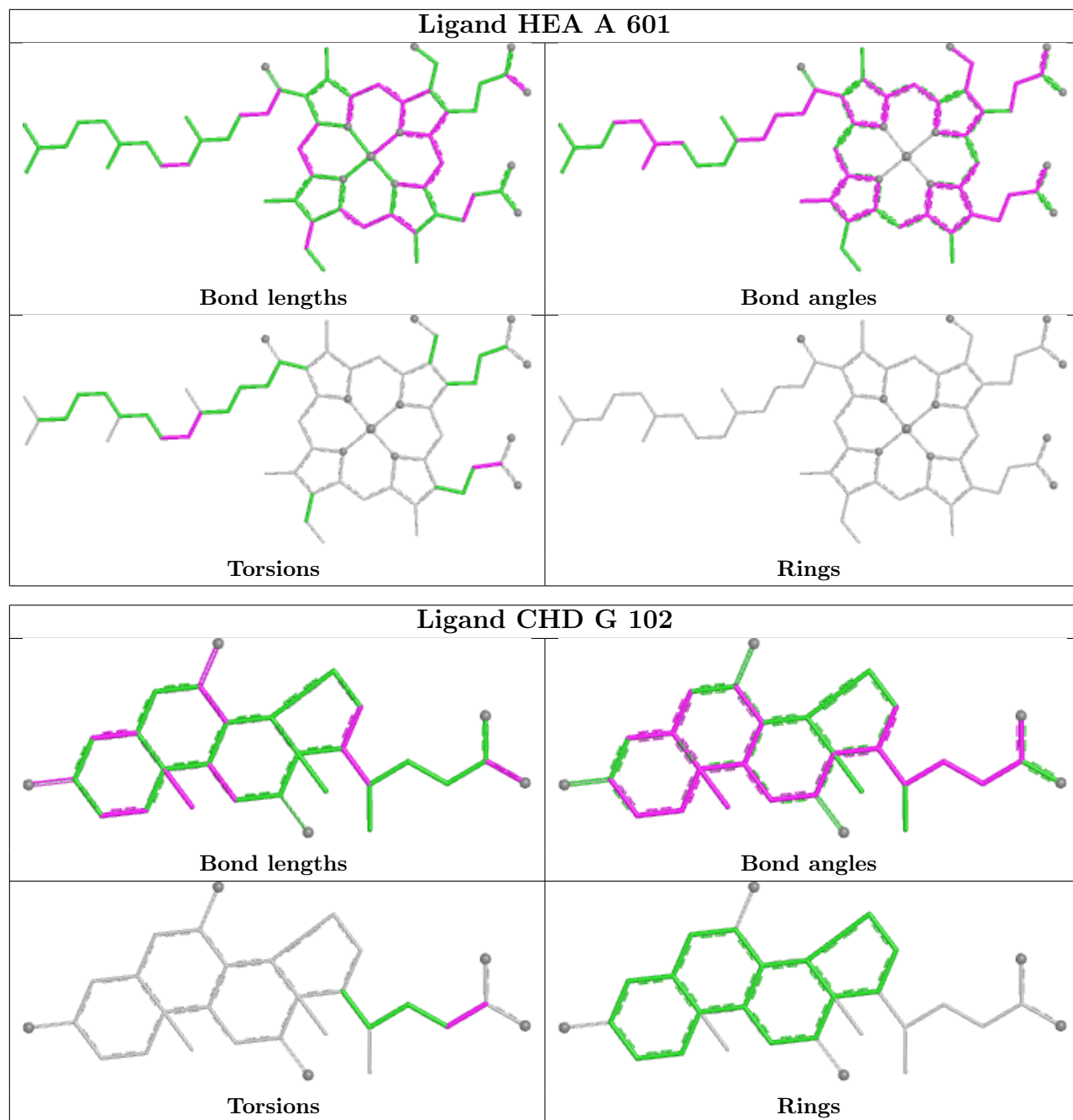
Mol	Chain	Res	Type	Clashes	Symm-Clashes
22	C	306	CHD	5	0
21	N	621	EDO	1	0
19	D	201	TGL	13	0

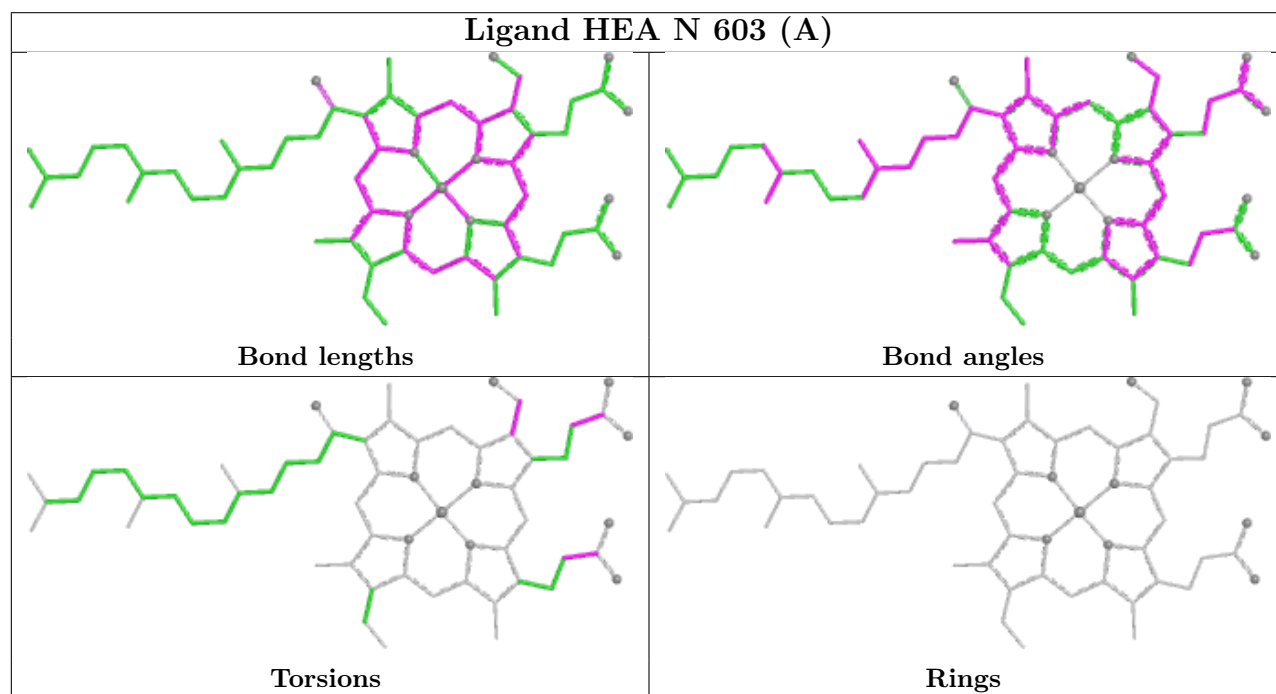
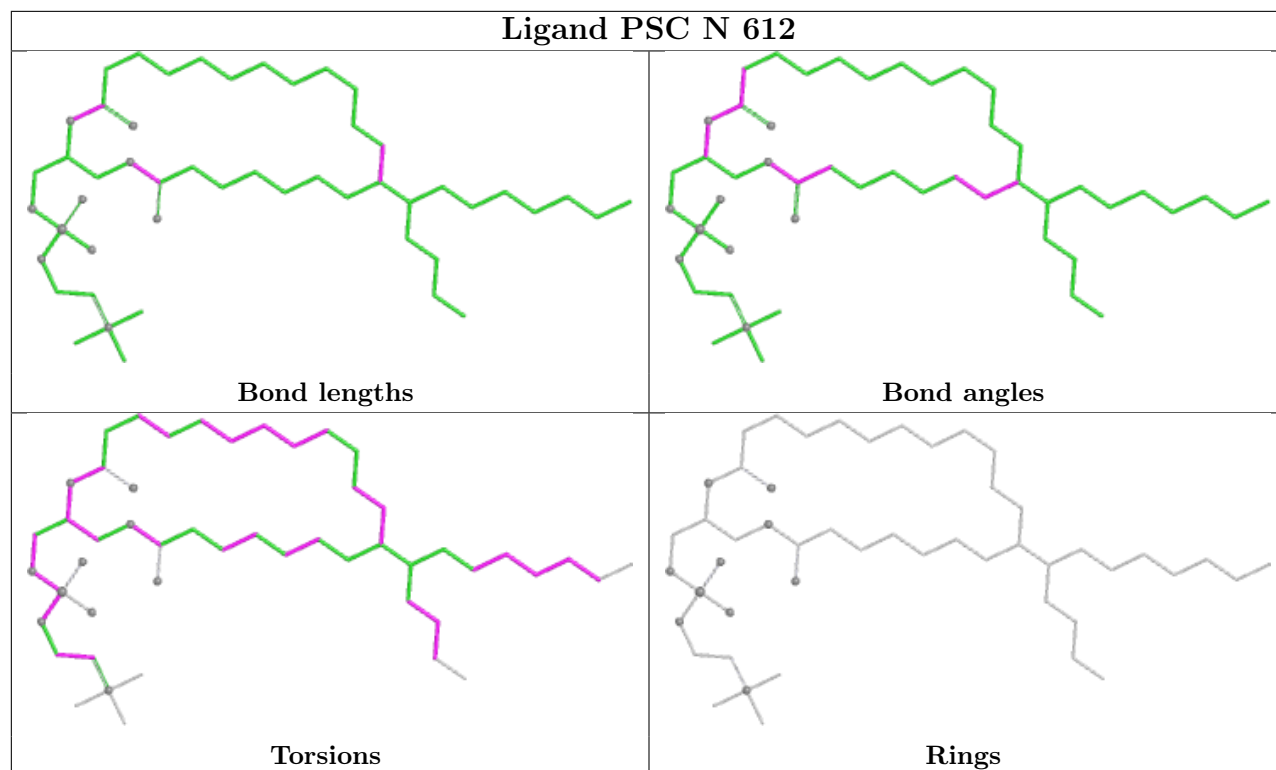
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

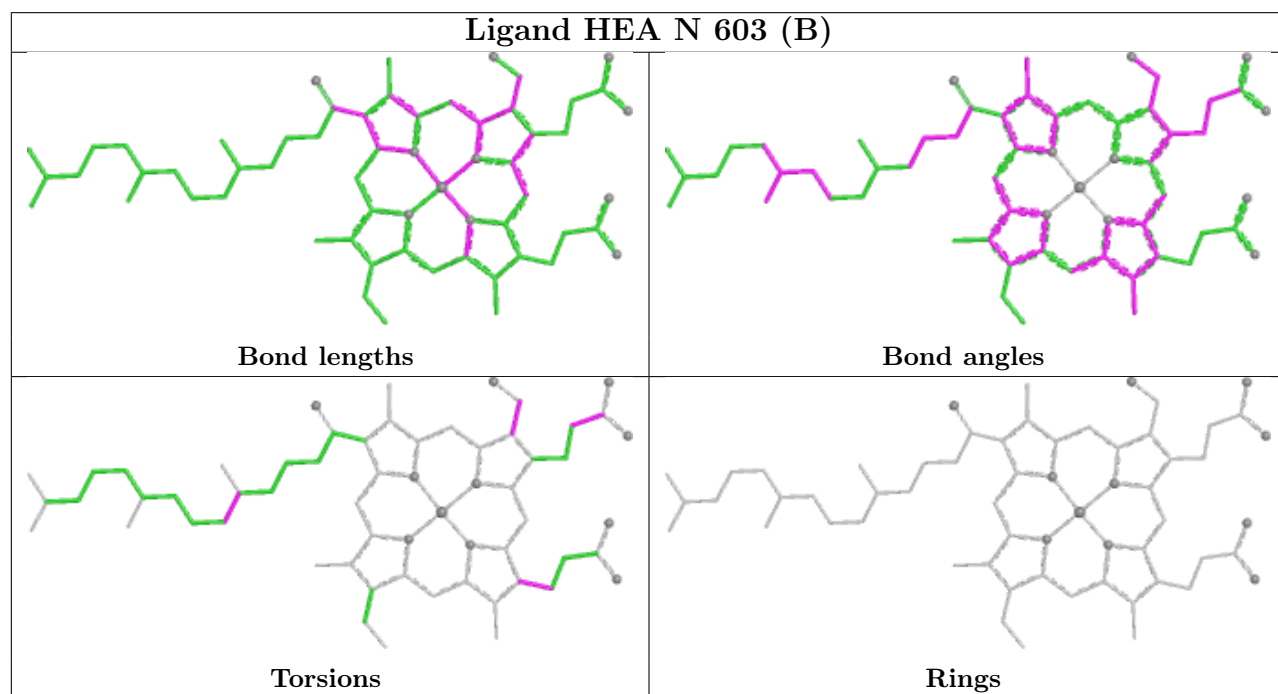
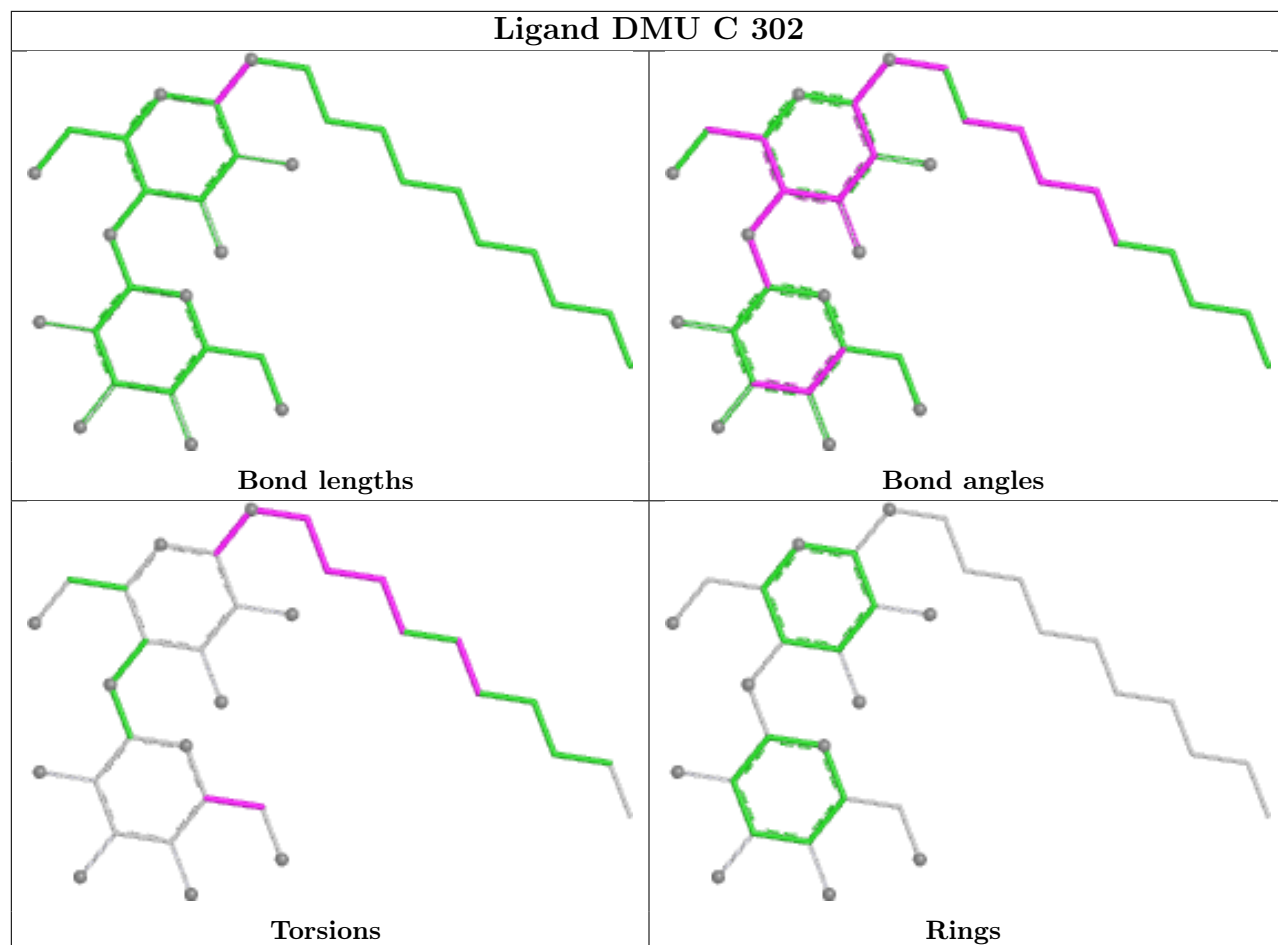


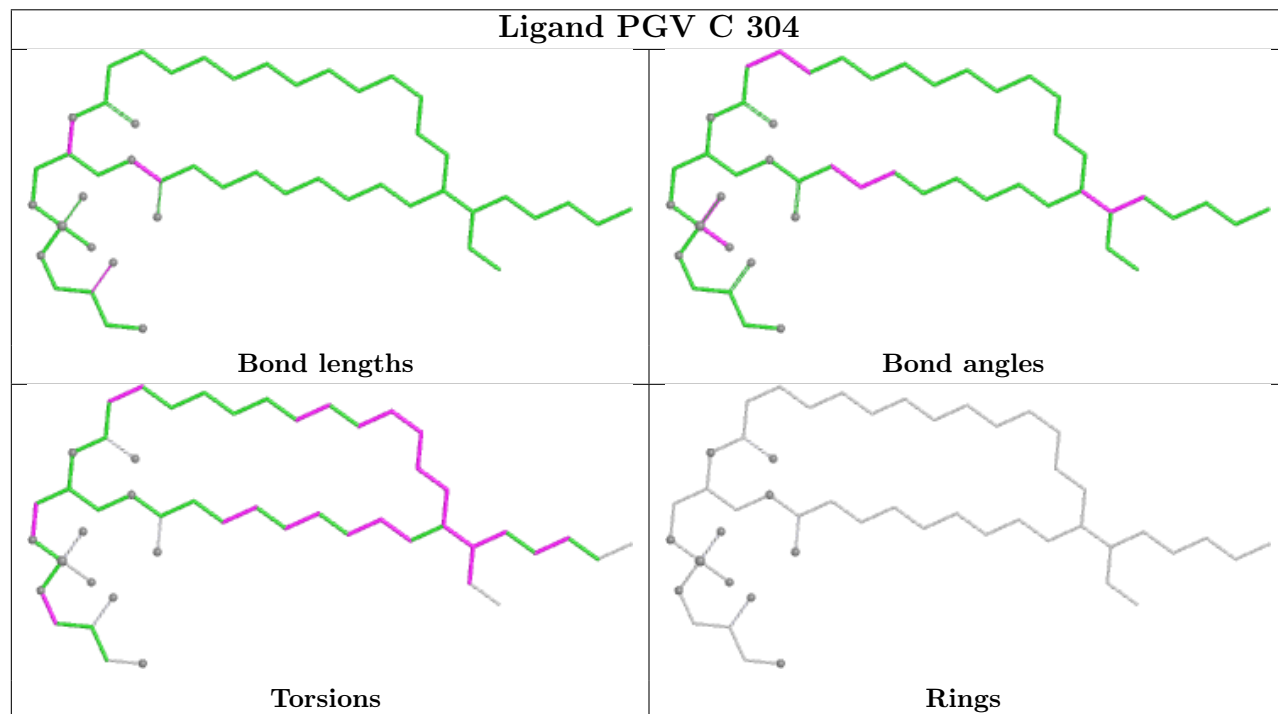
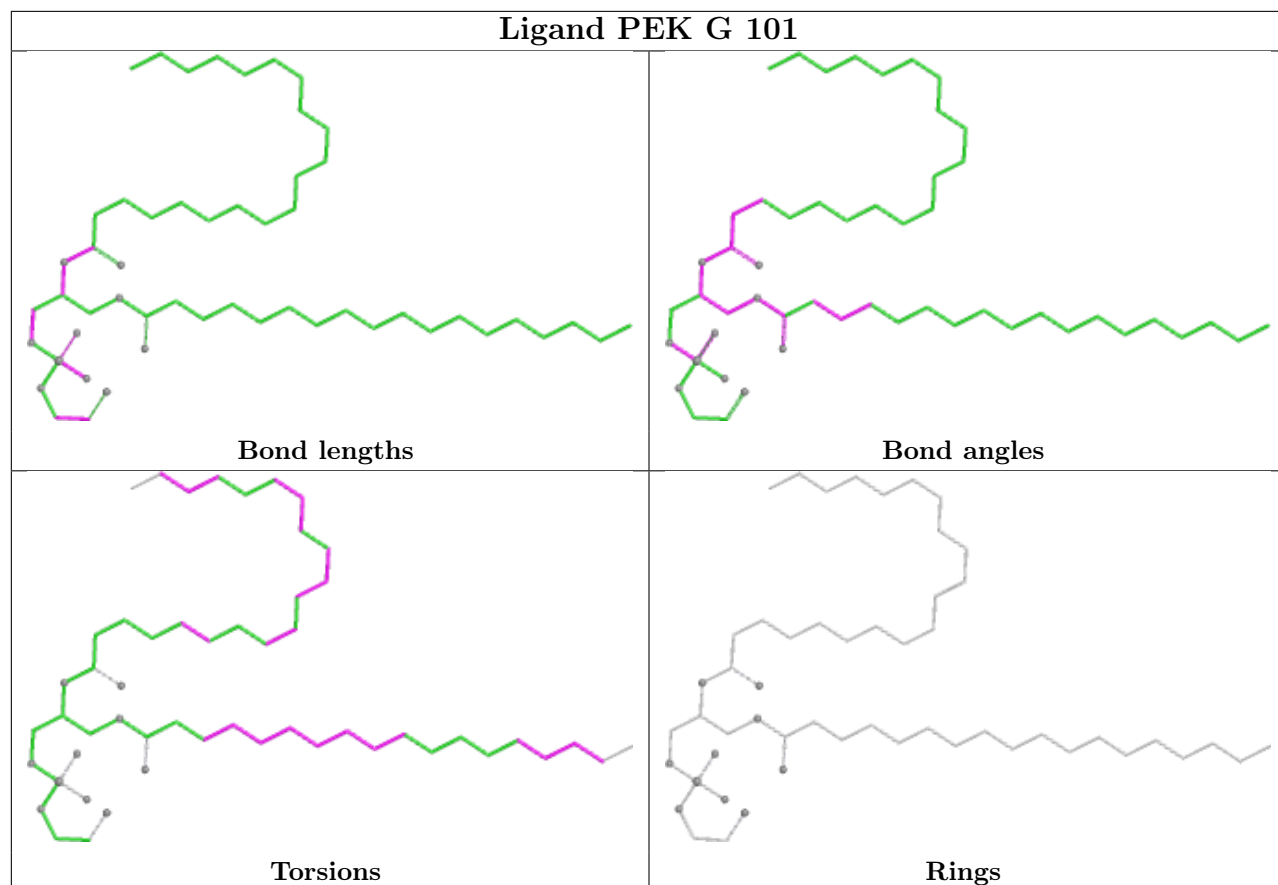


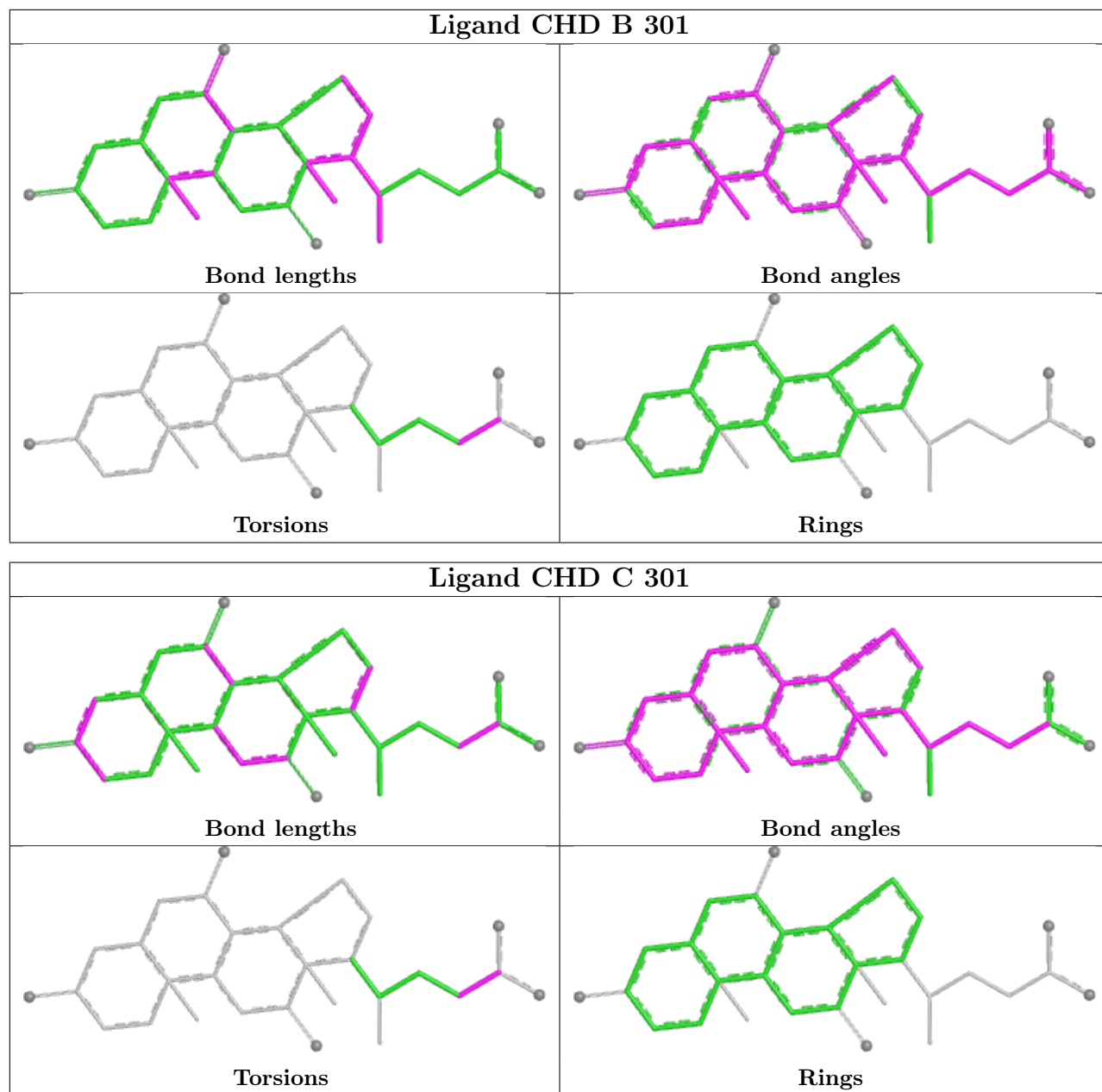


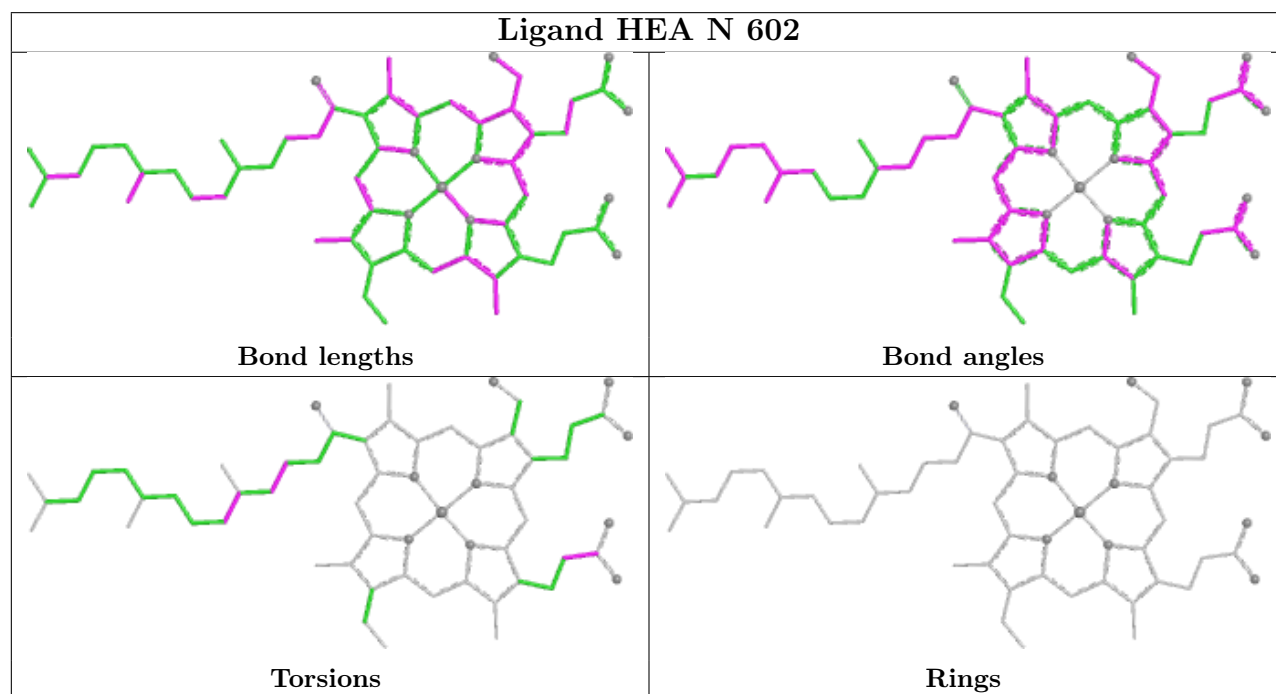
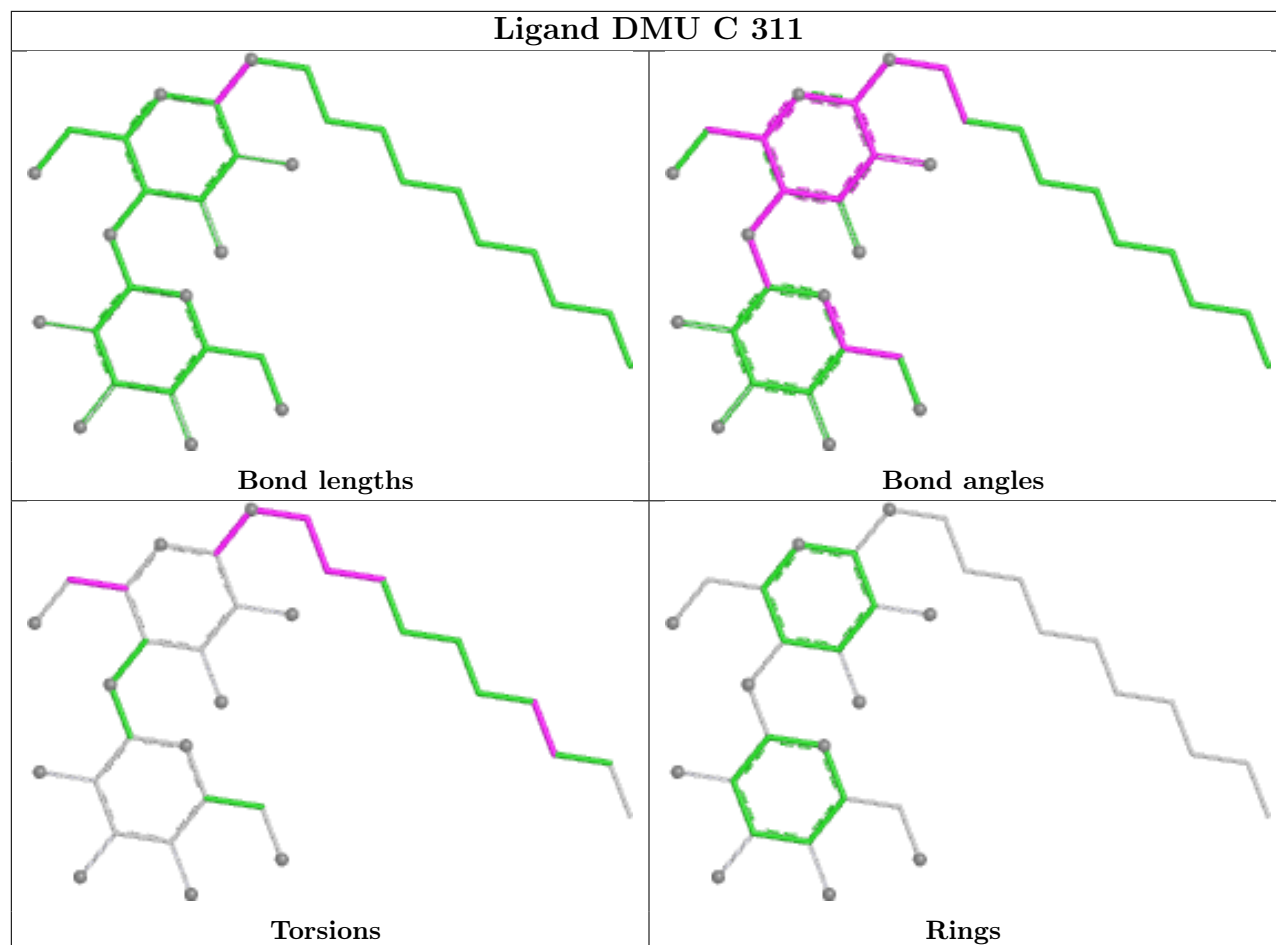


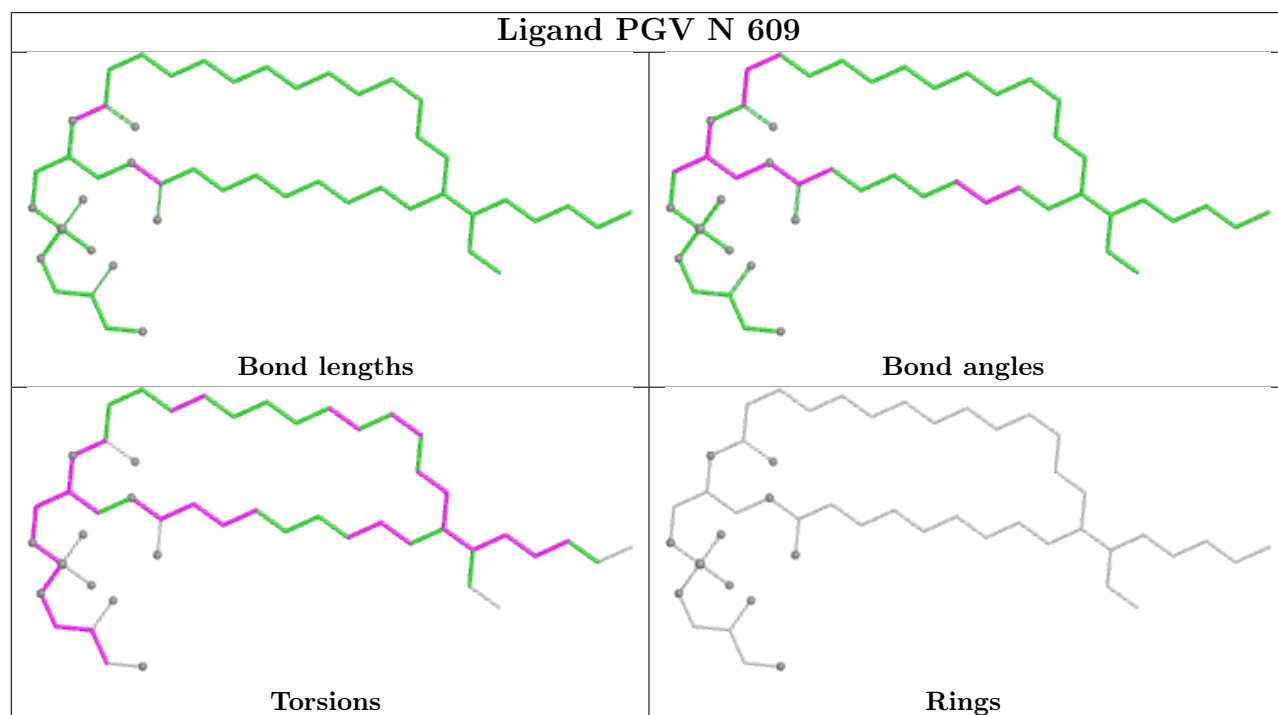
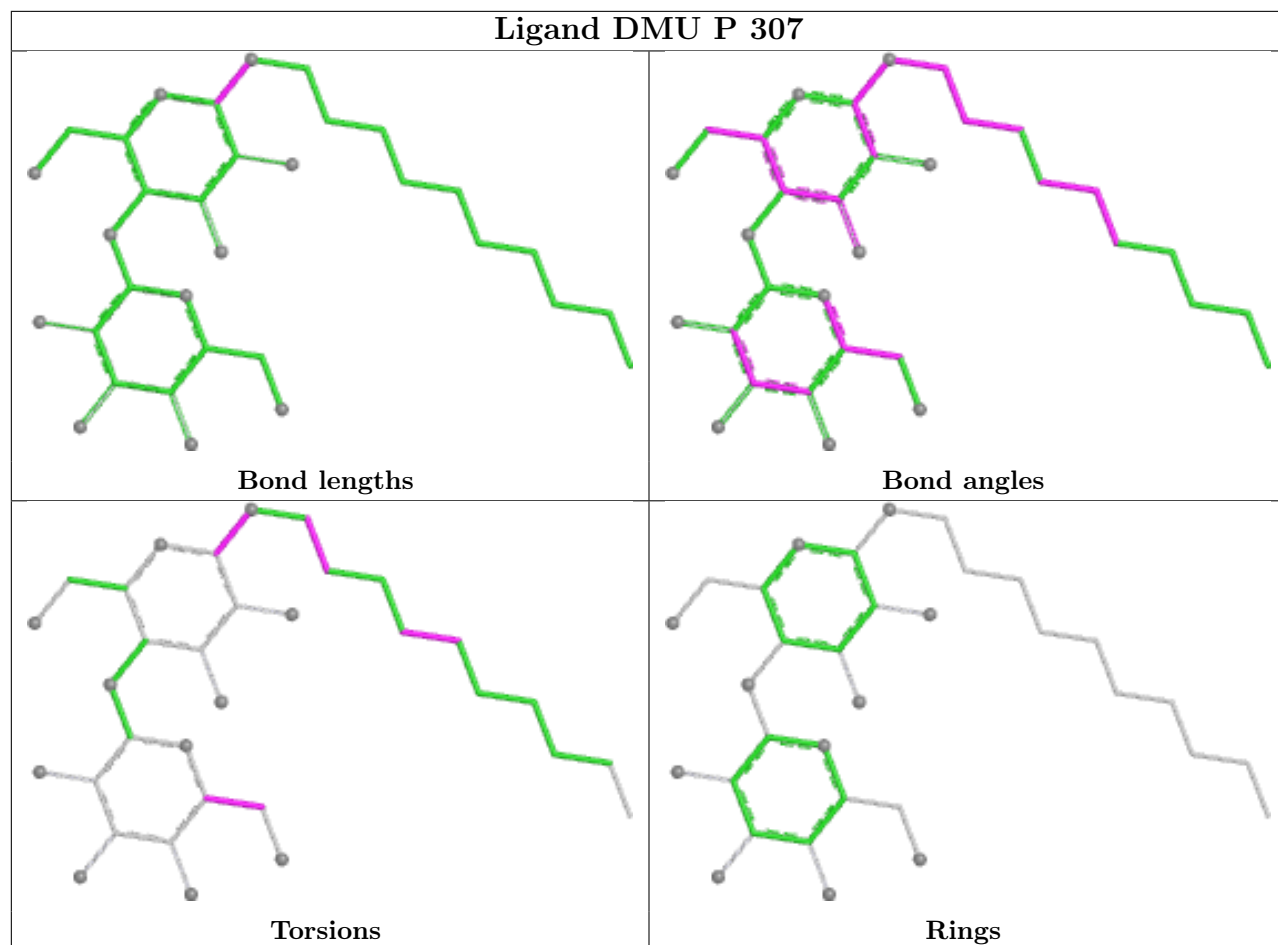


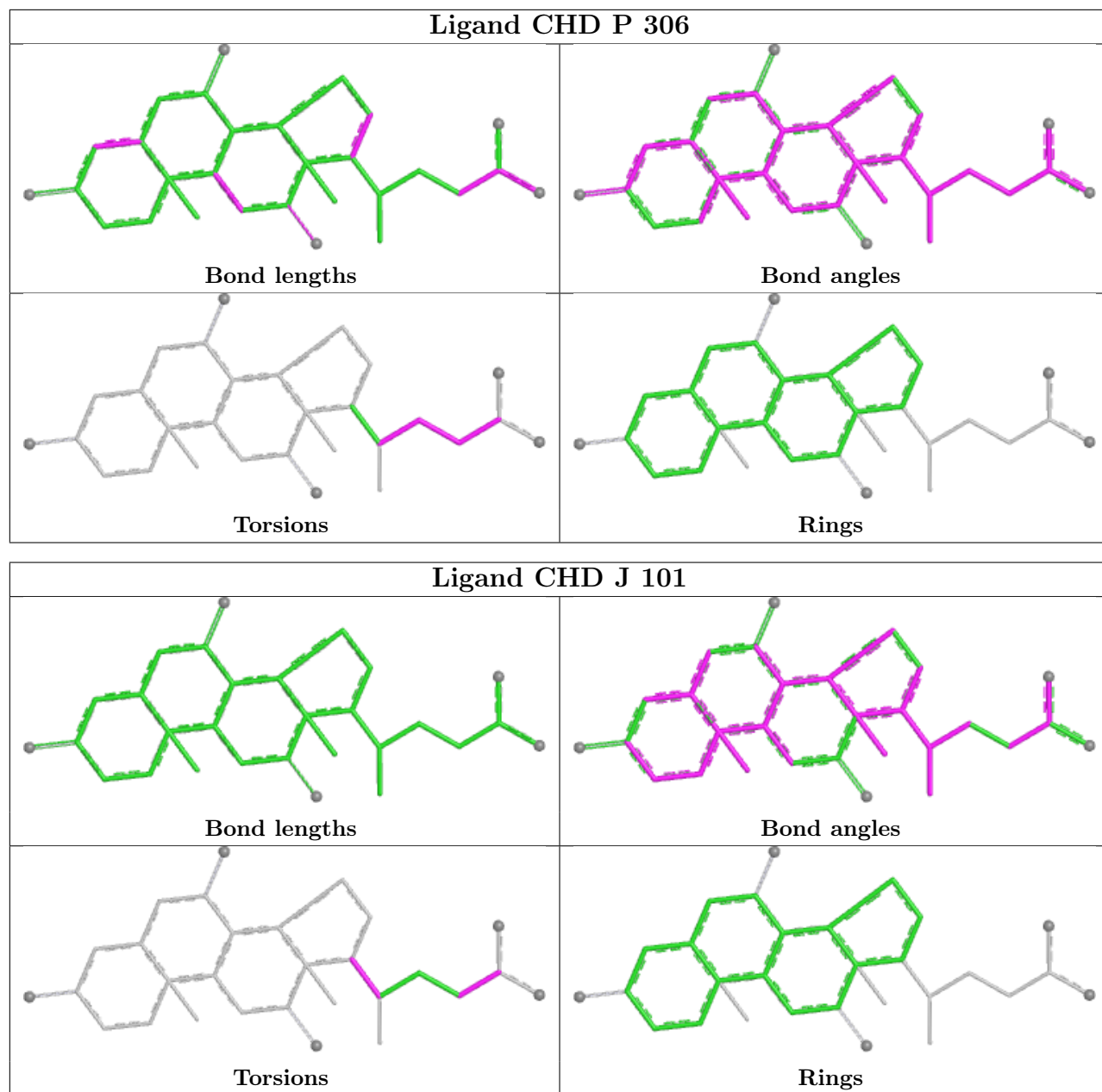


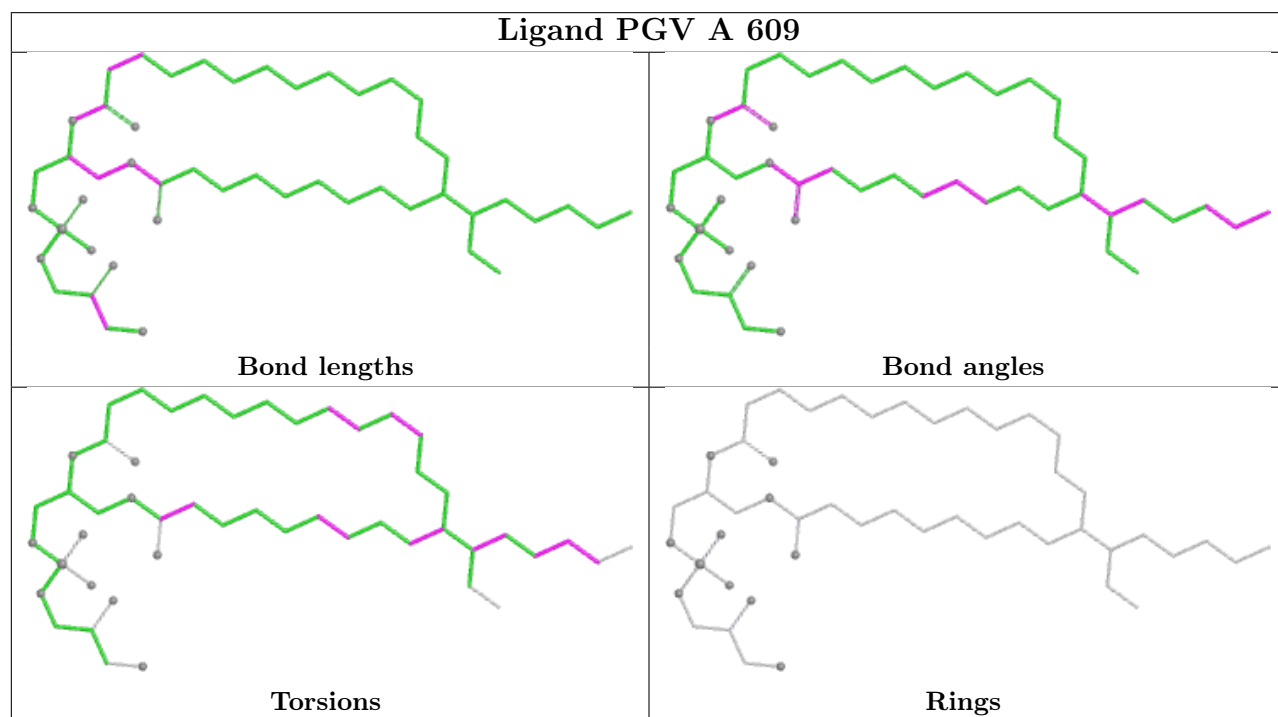
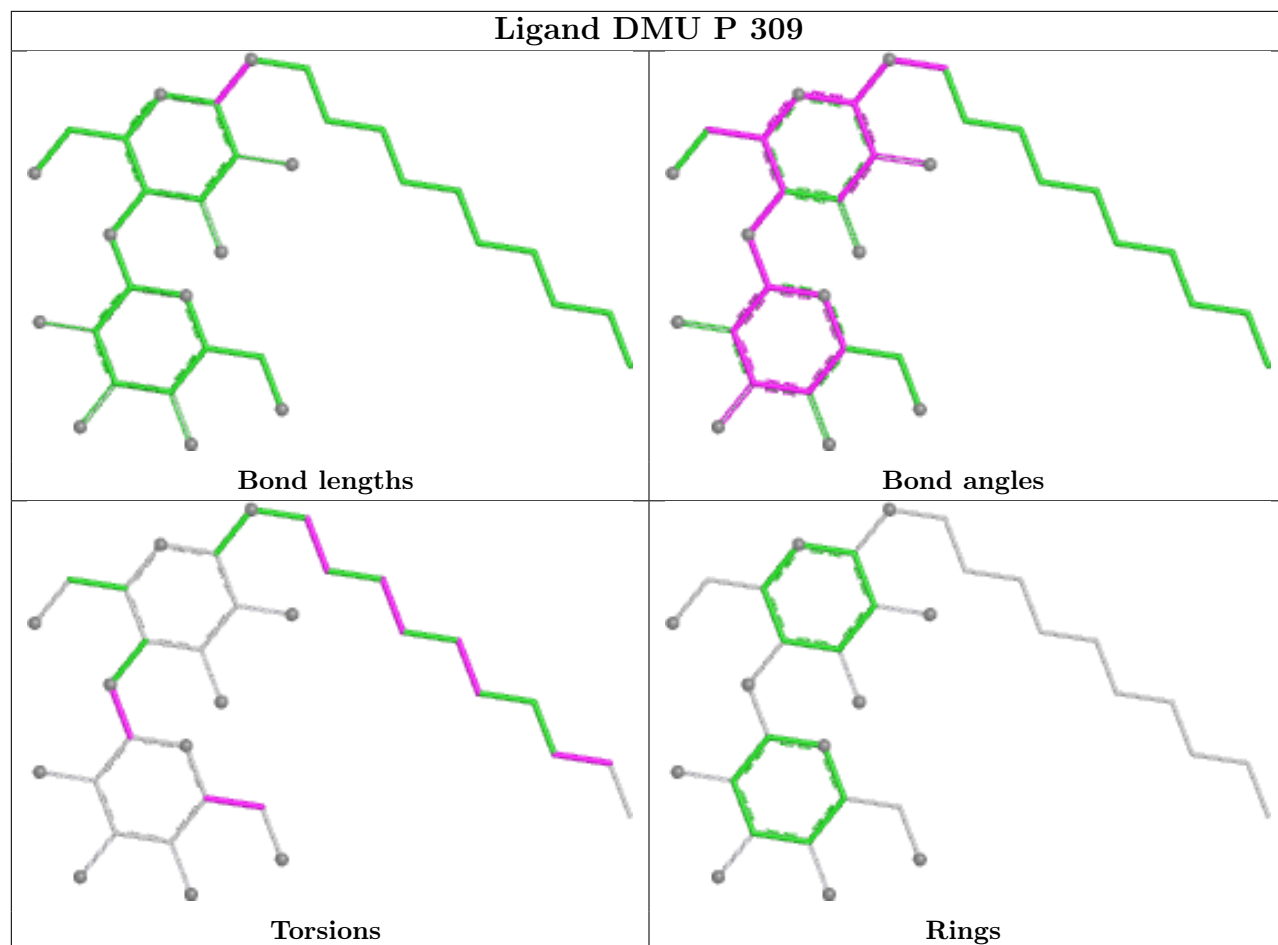


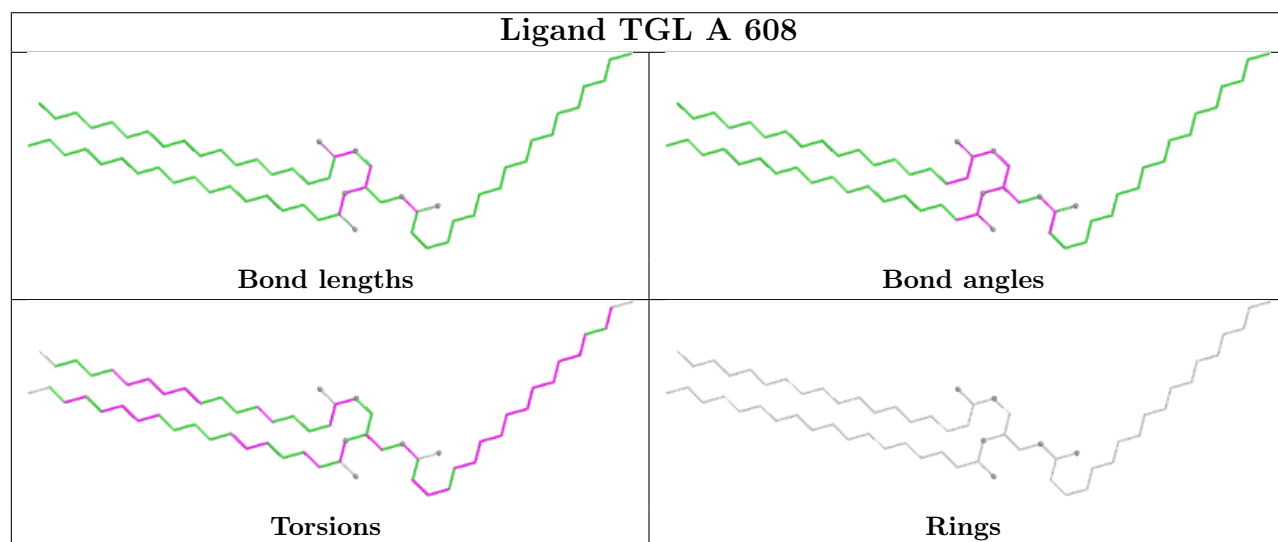
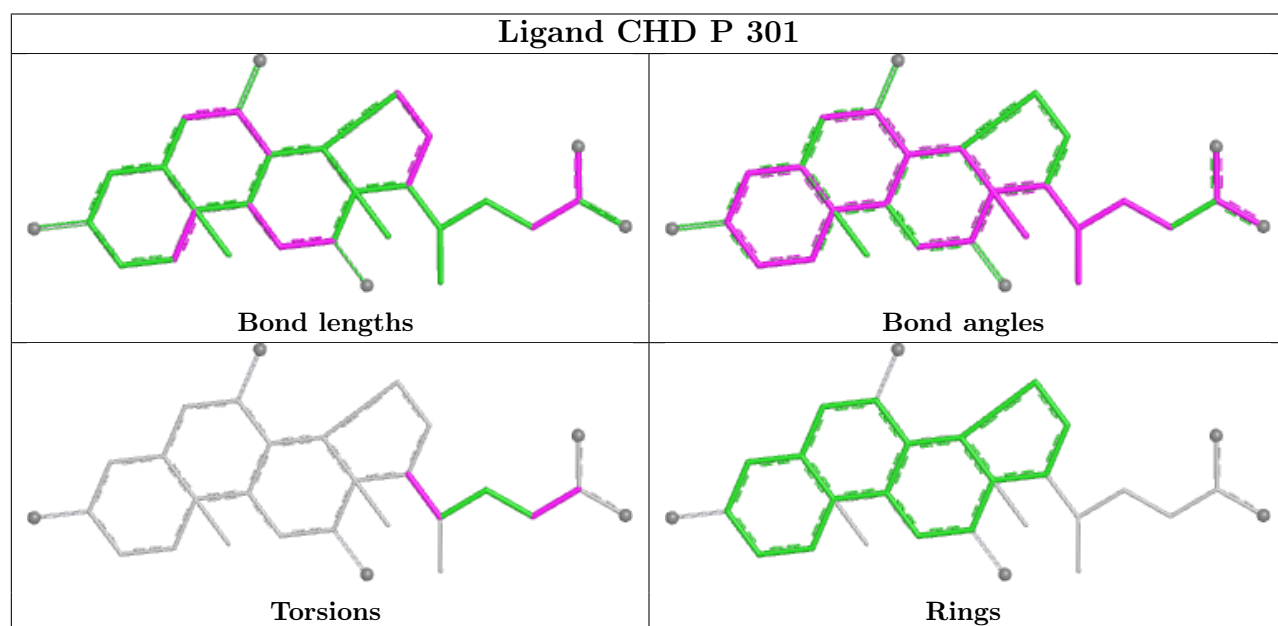
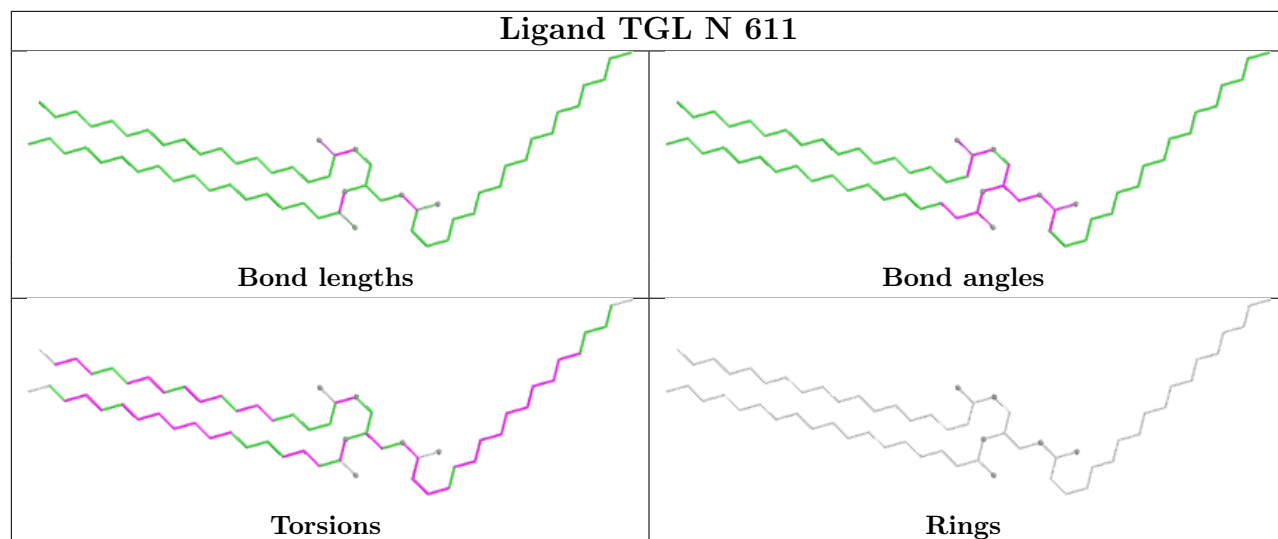


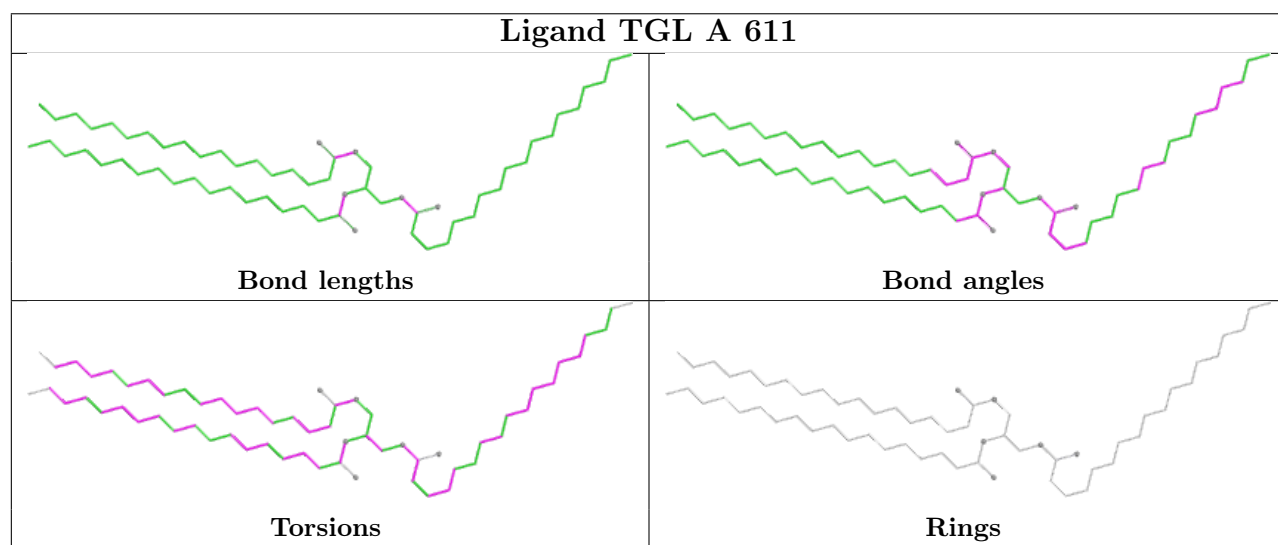
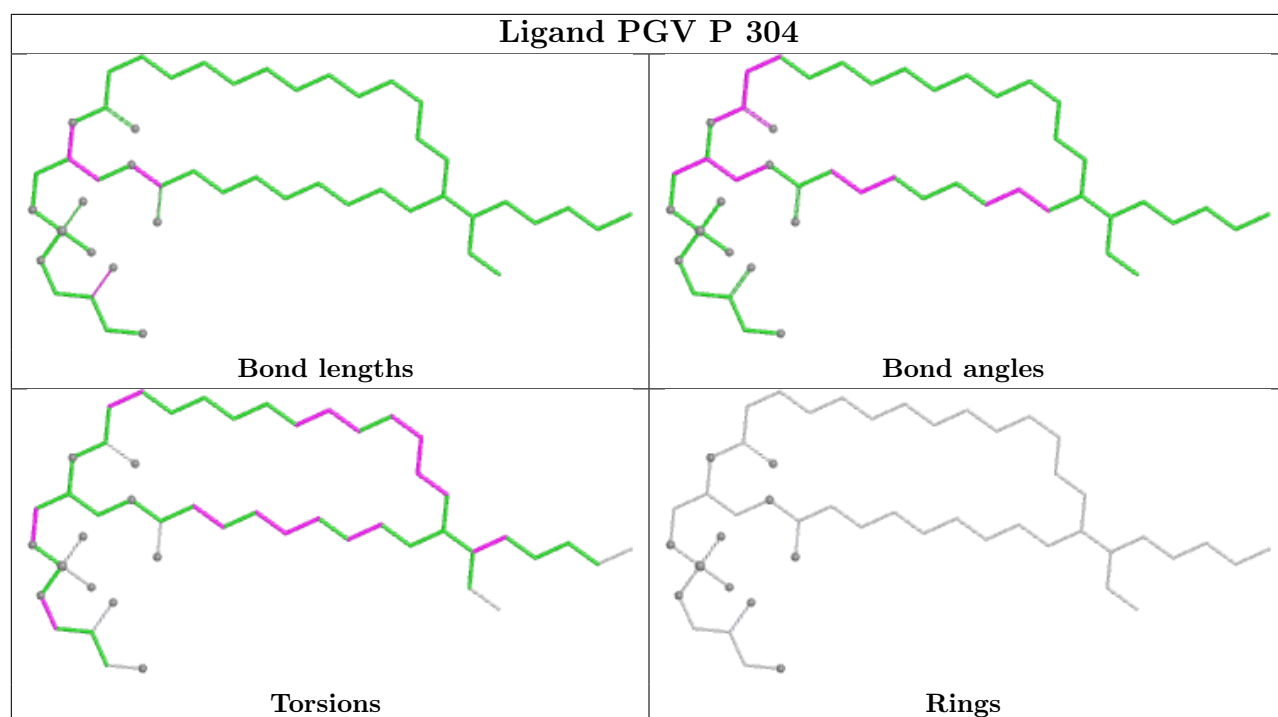


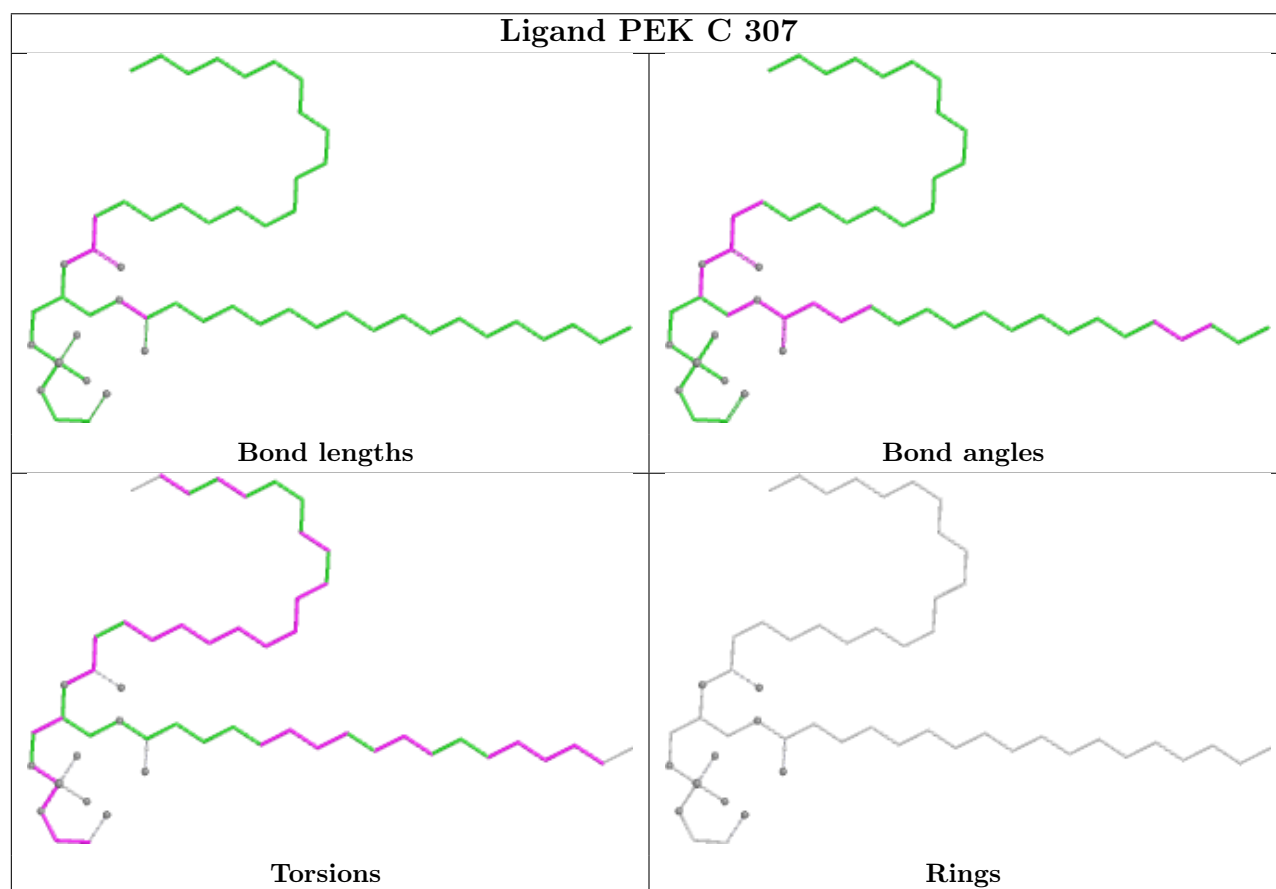
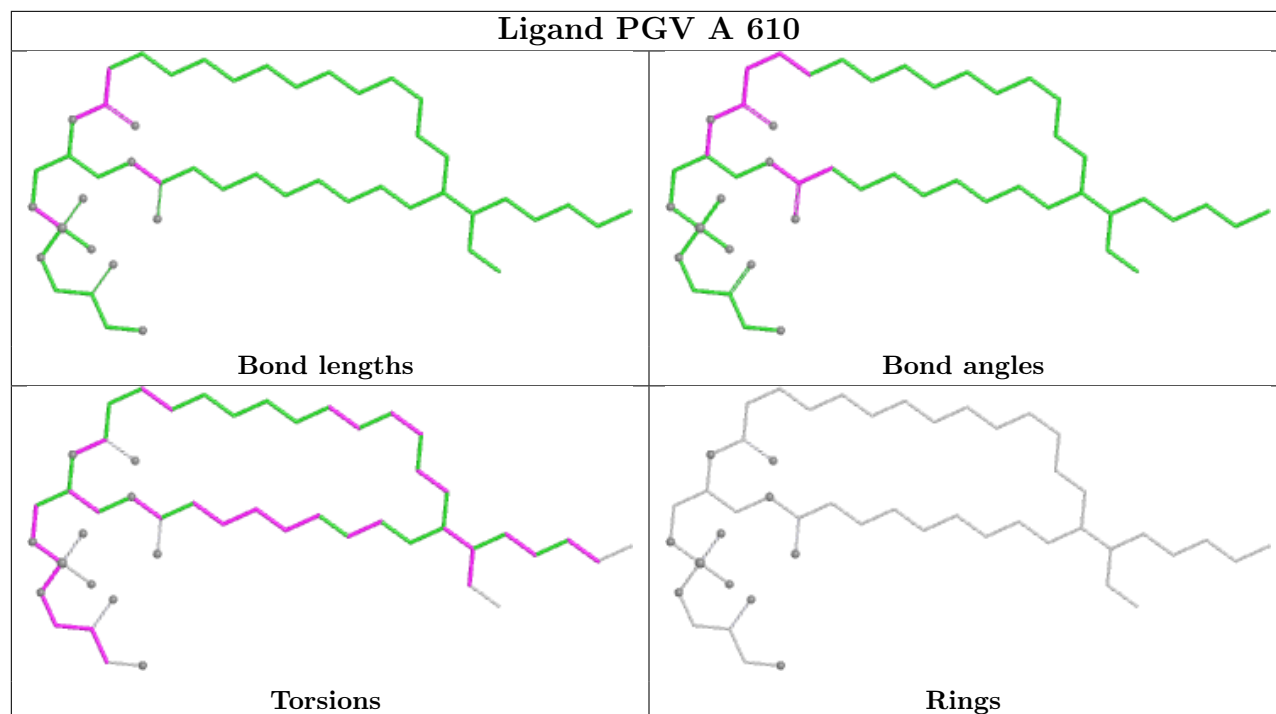


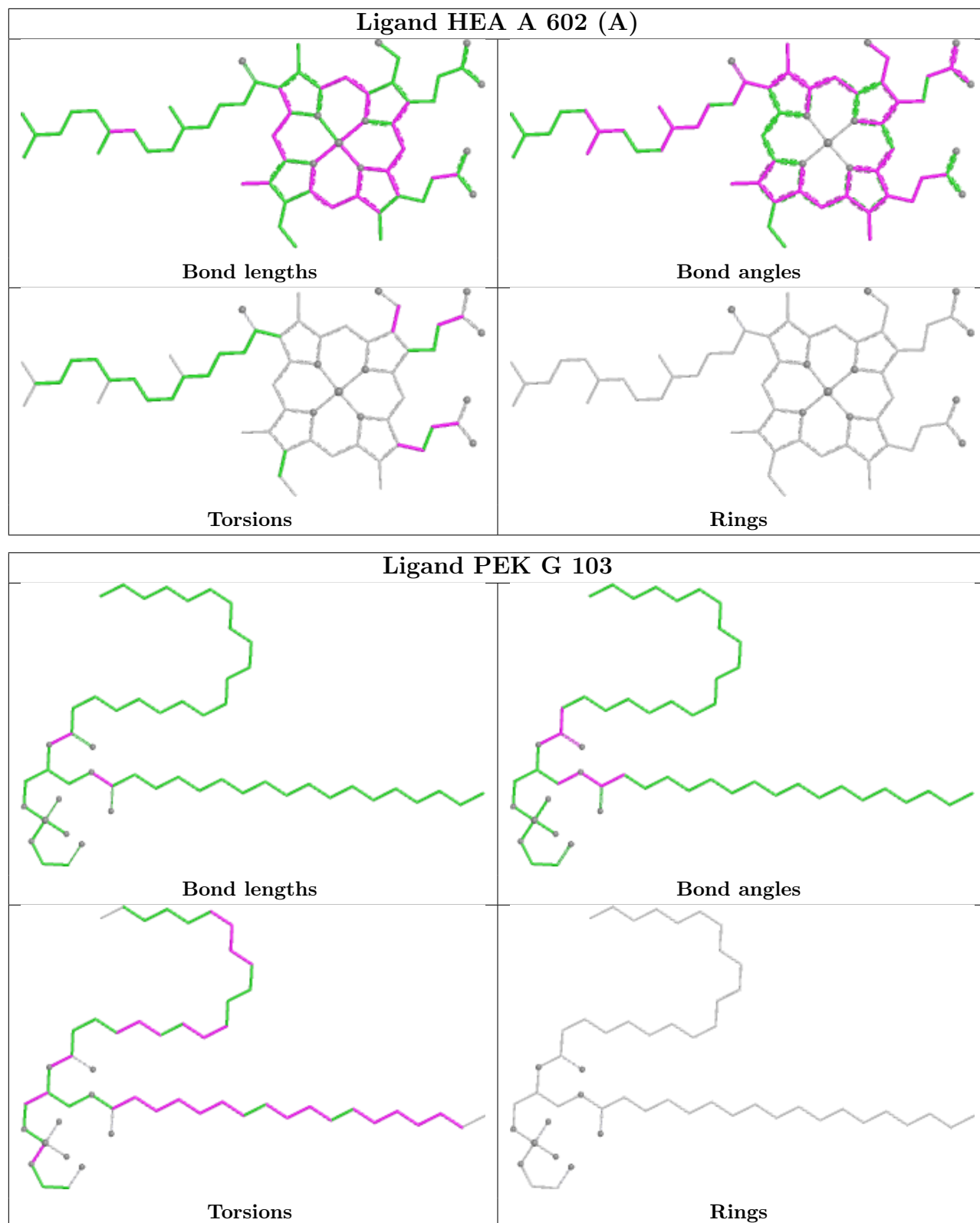


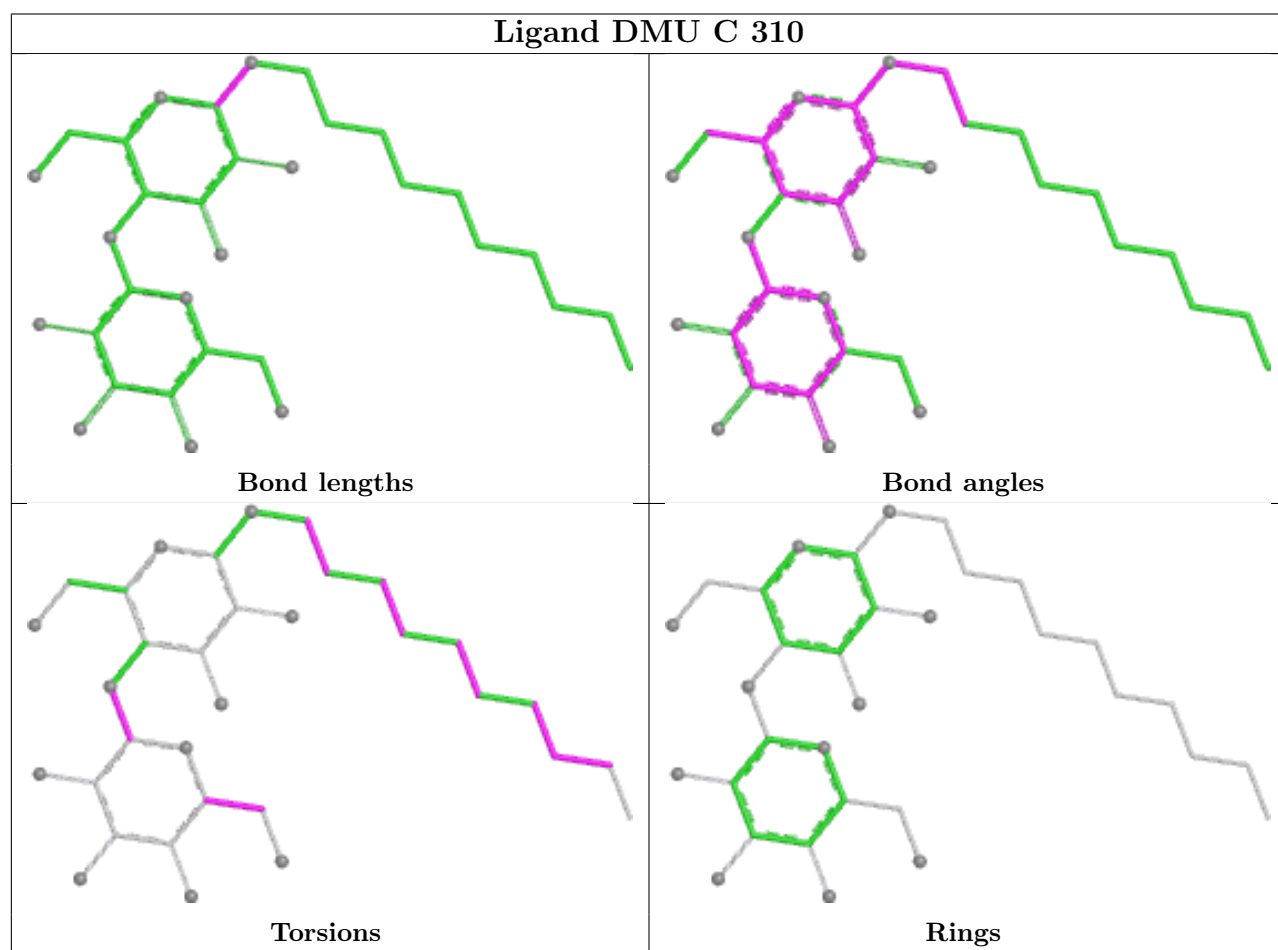
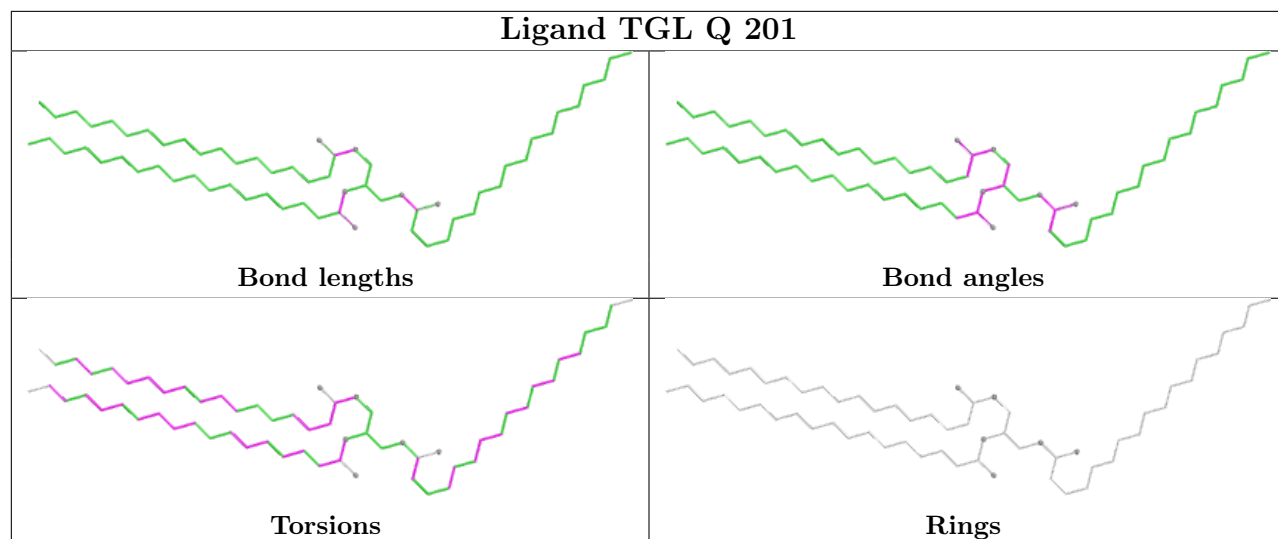


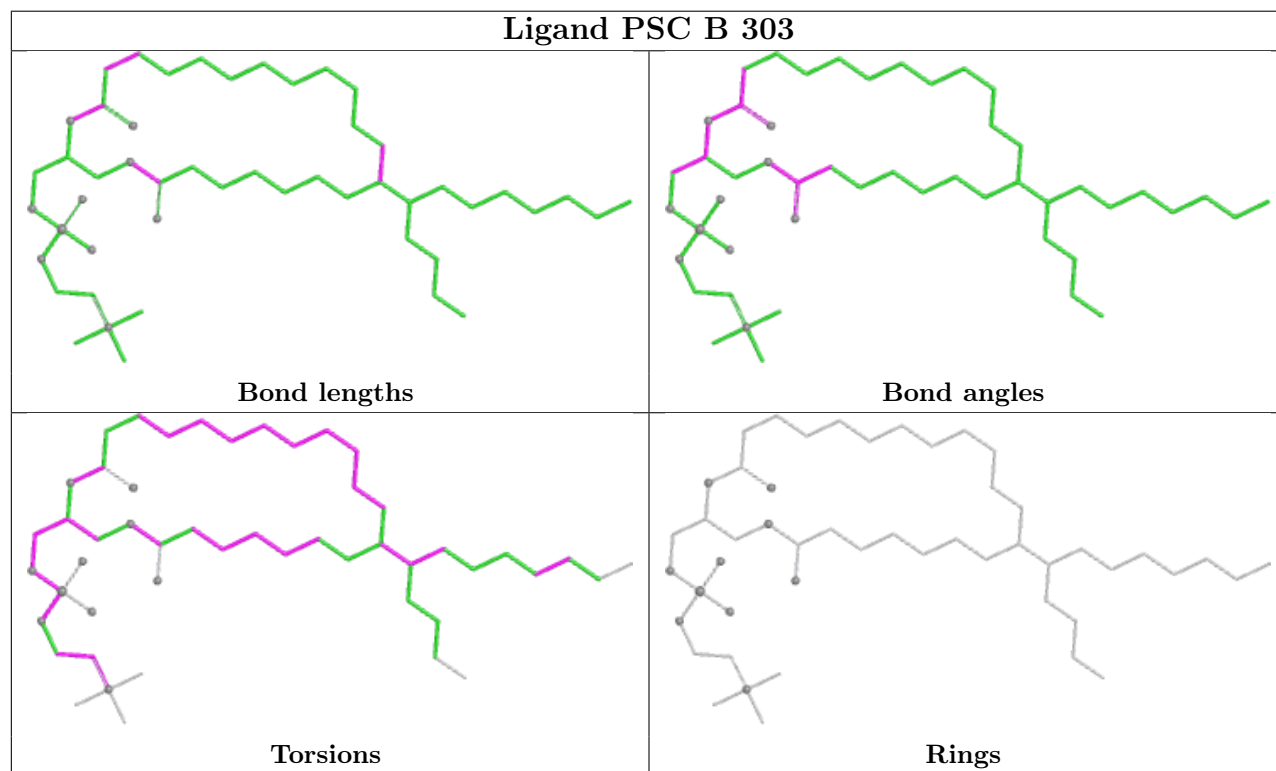
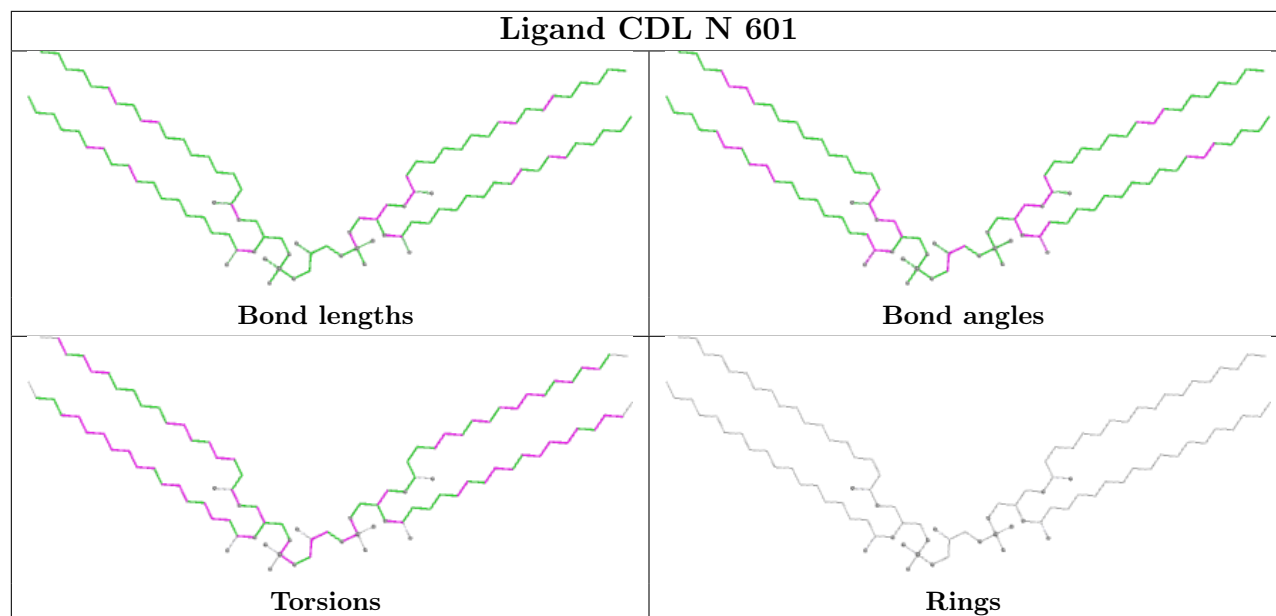


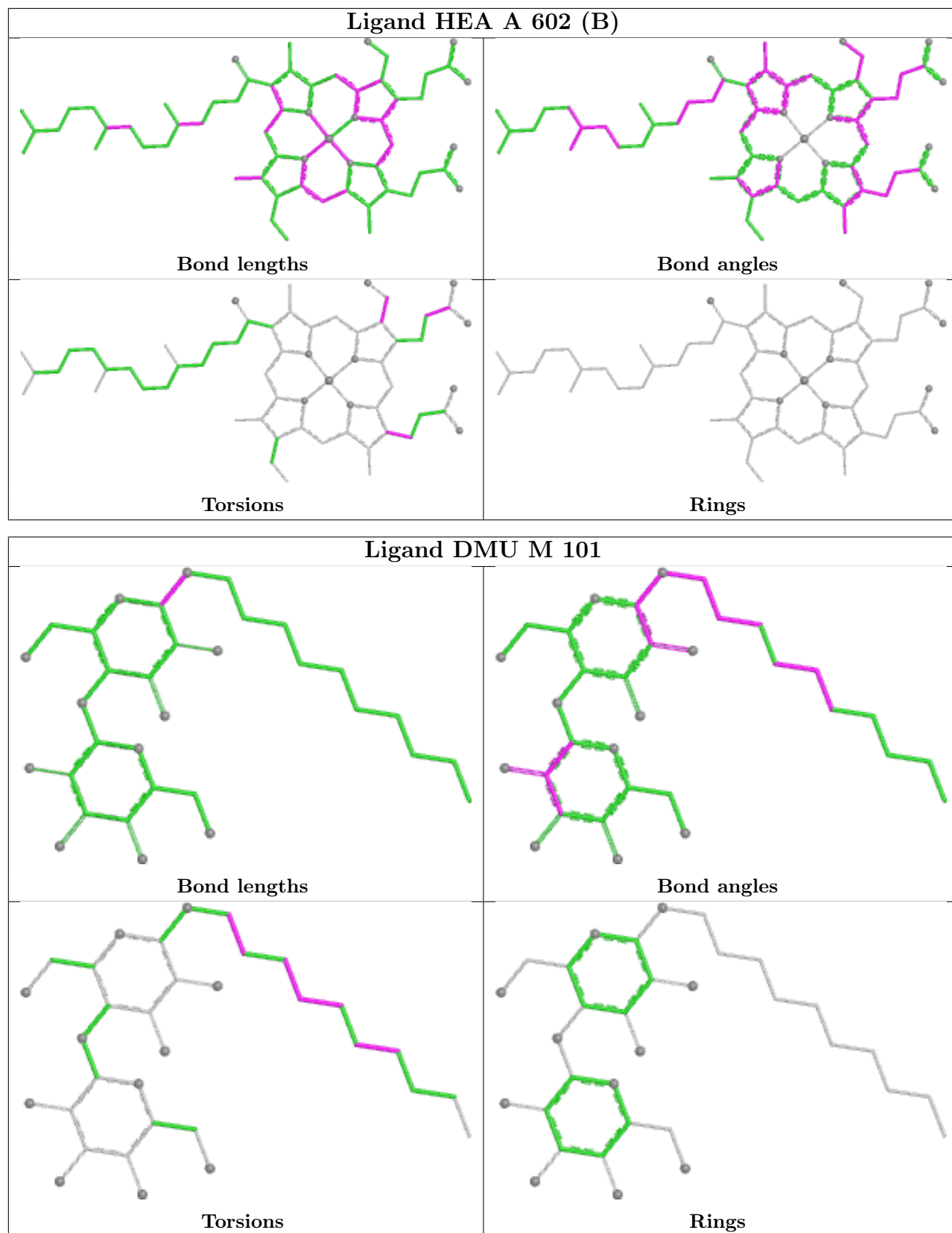




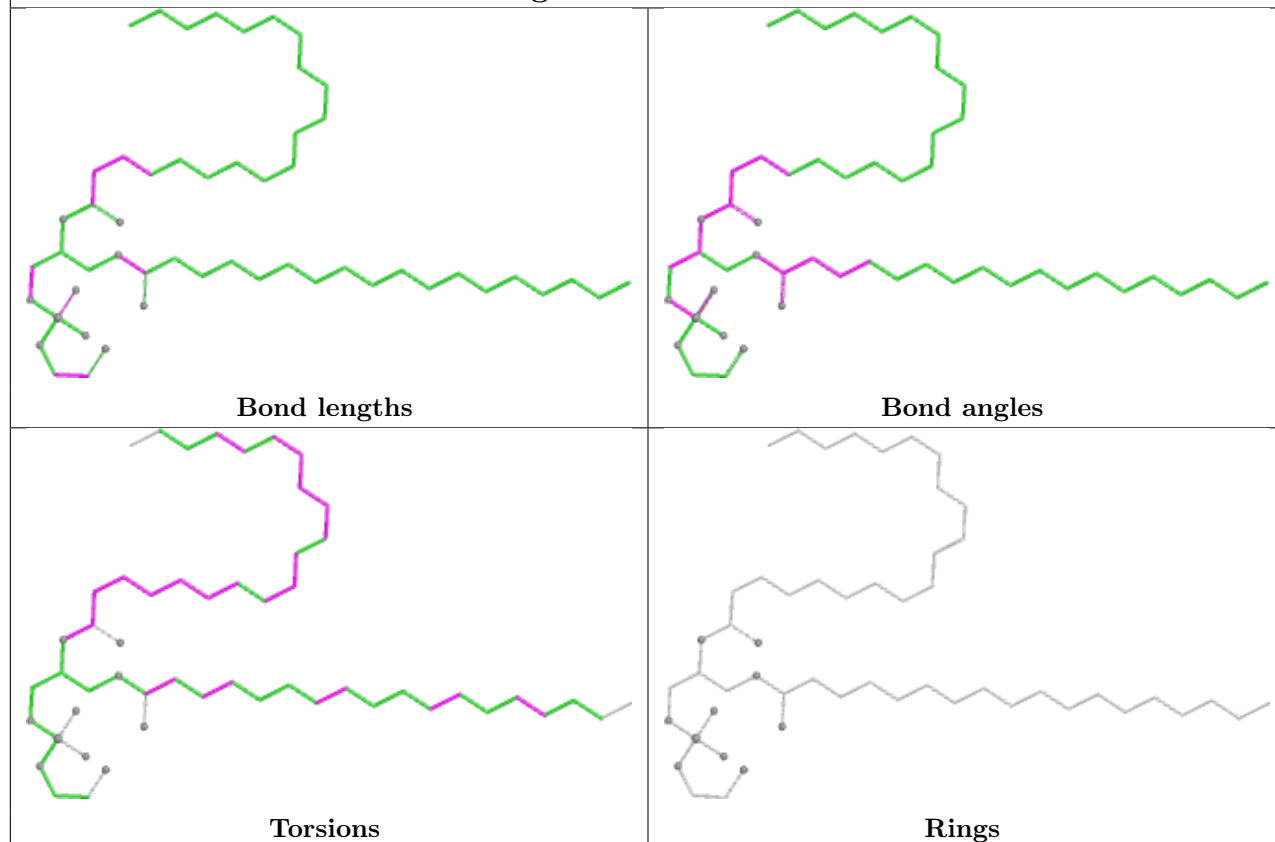




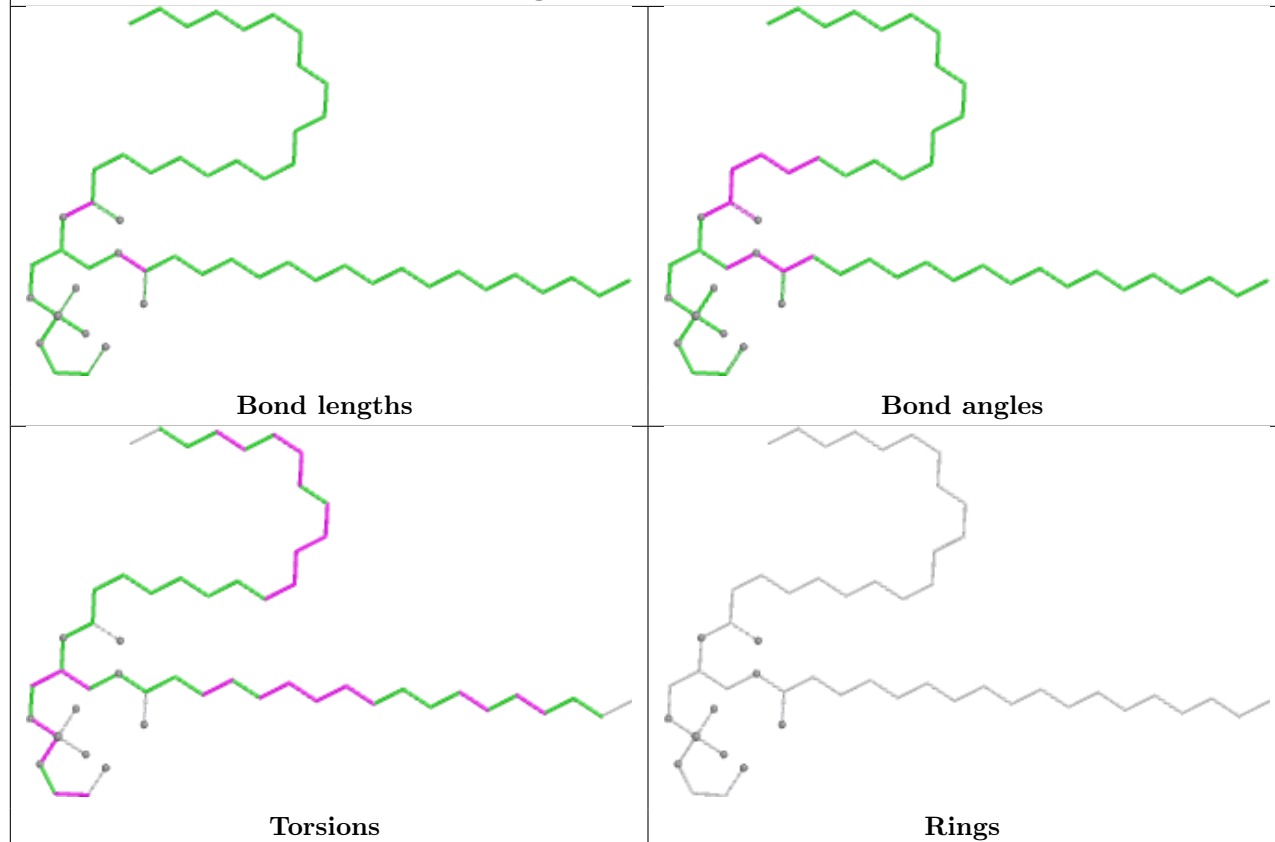


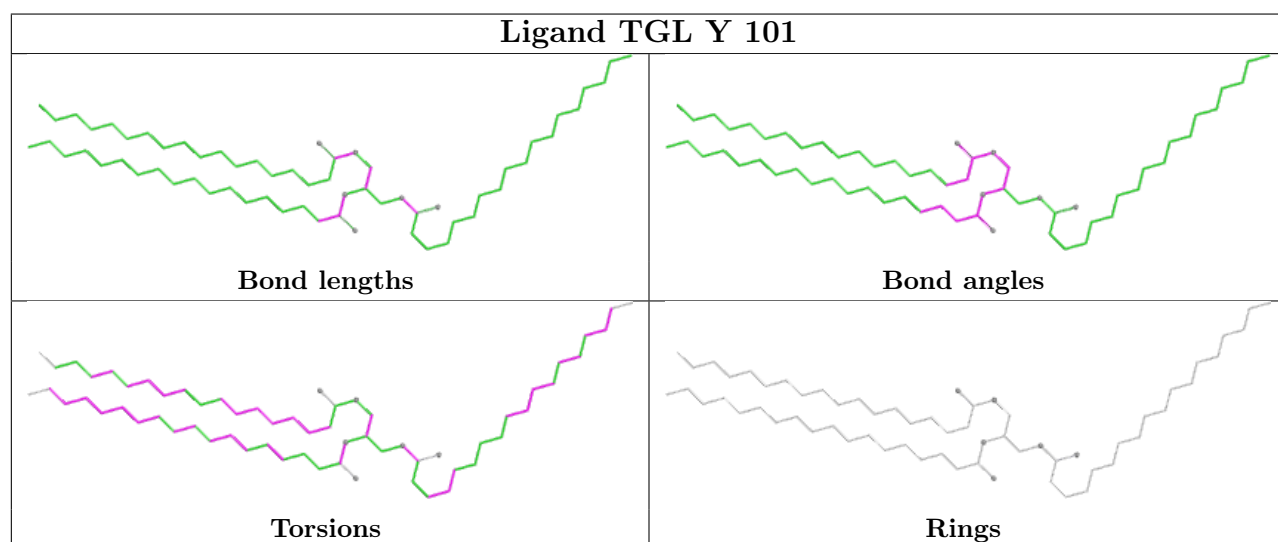
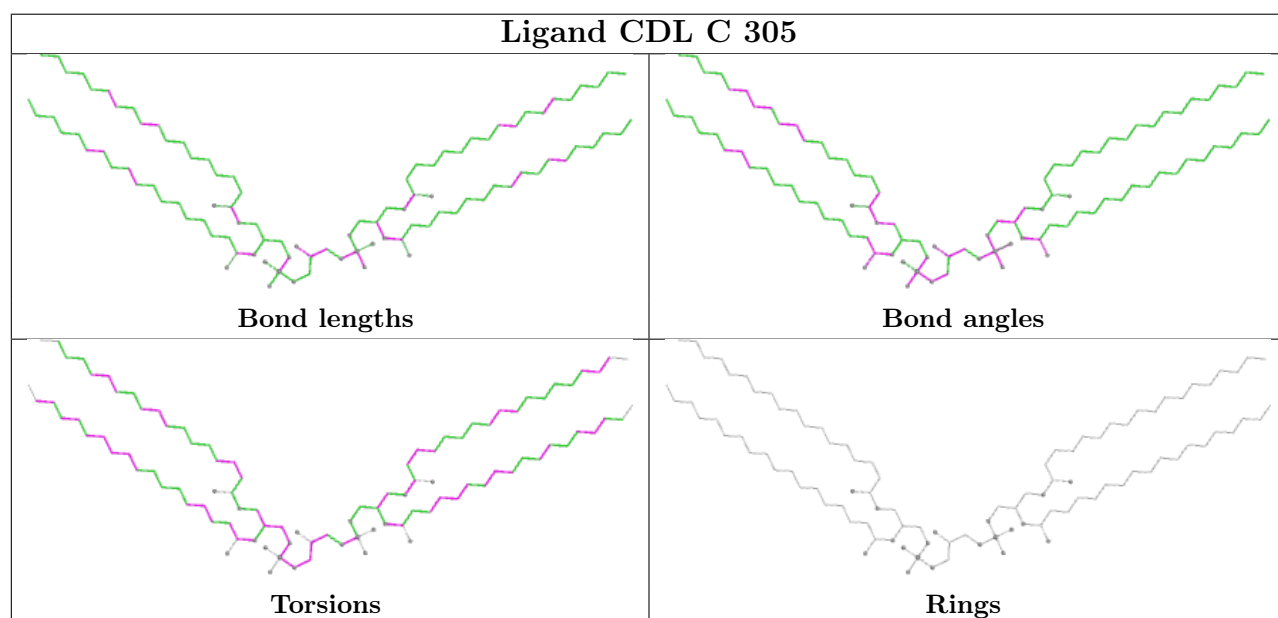
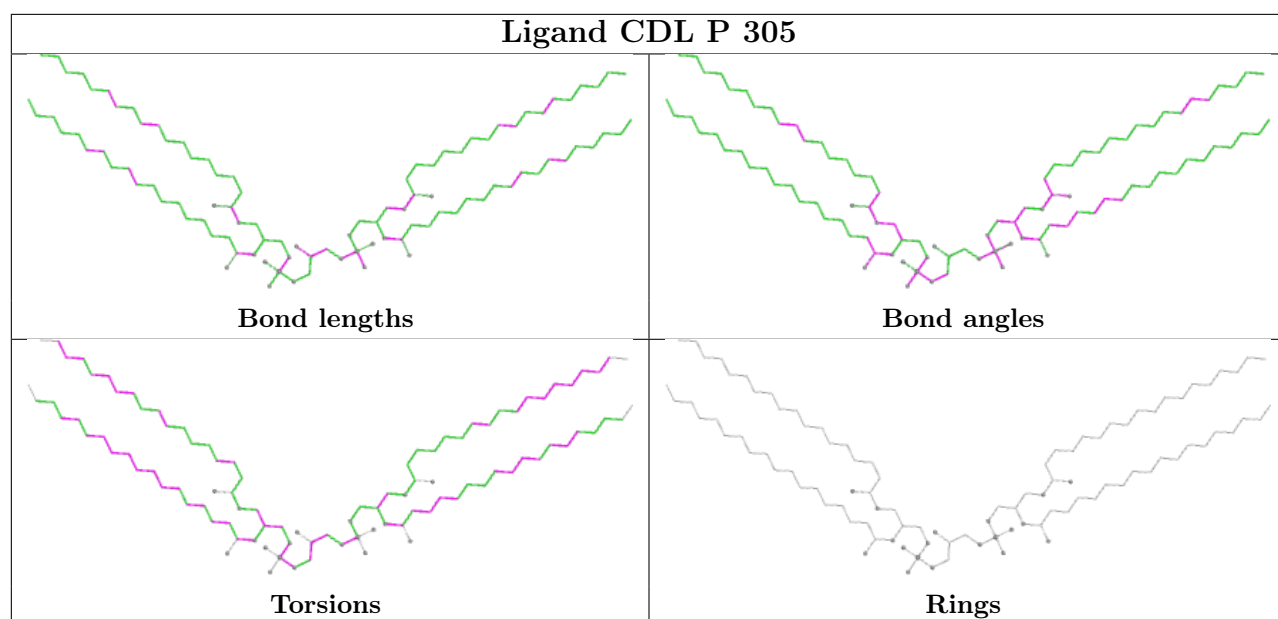


Ligand PEK T 101

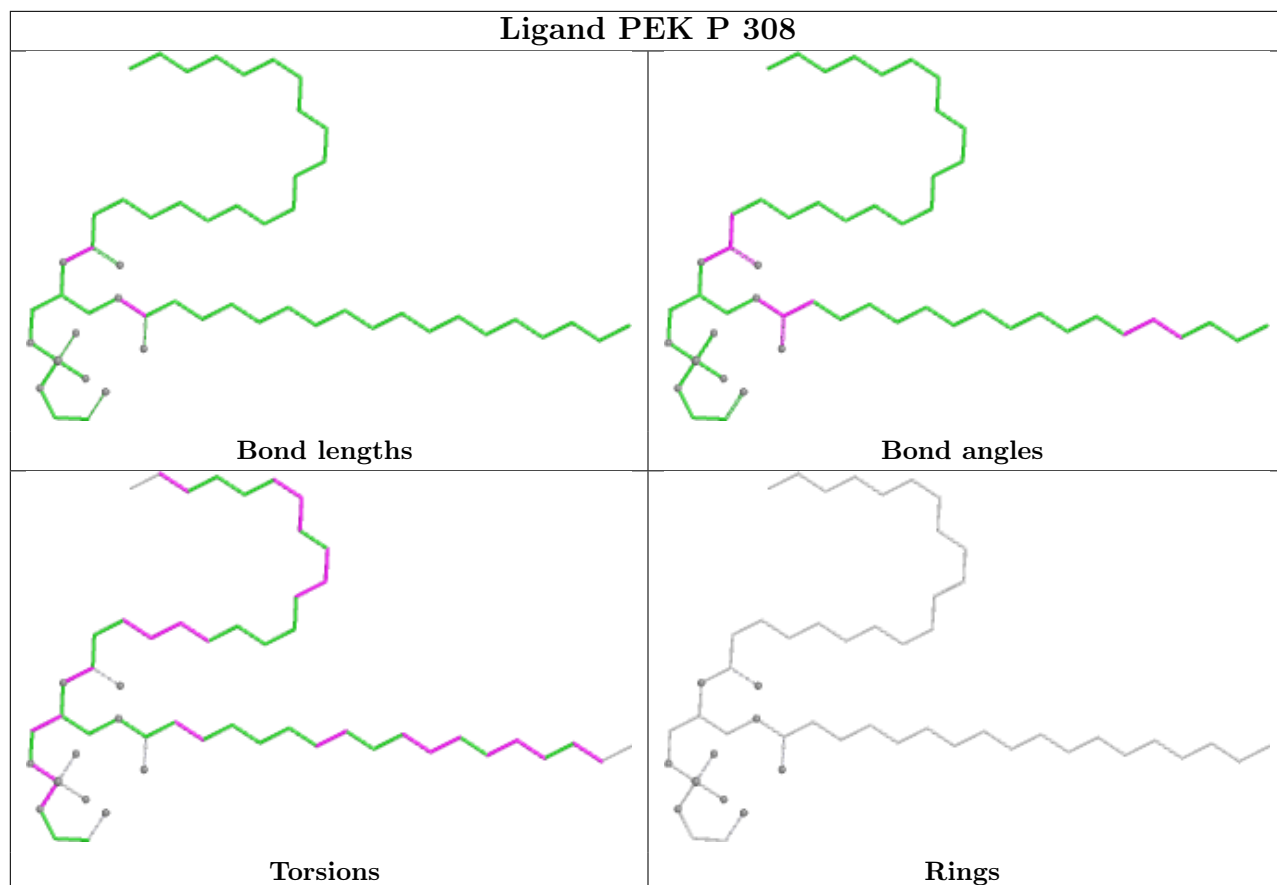


Ligand PEK C 309

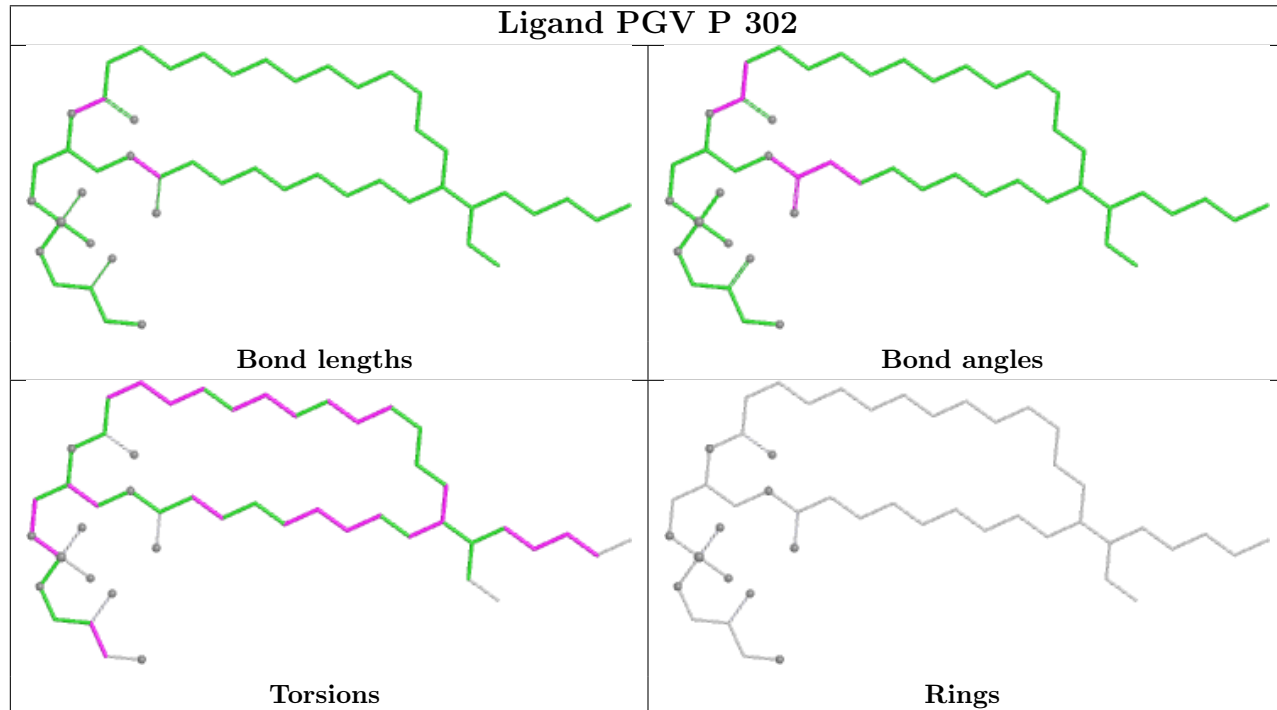


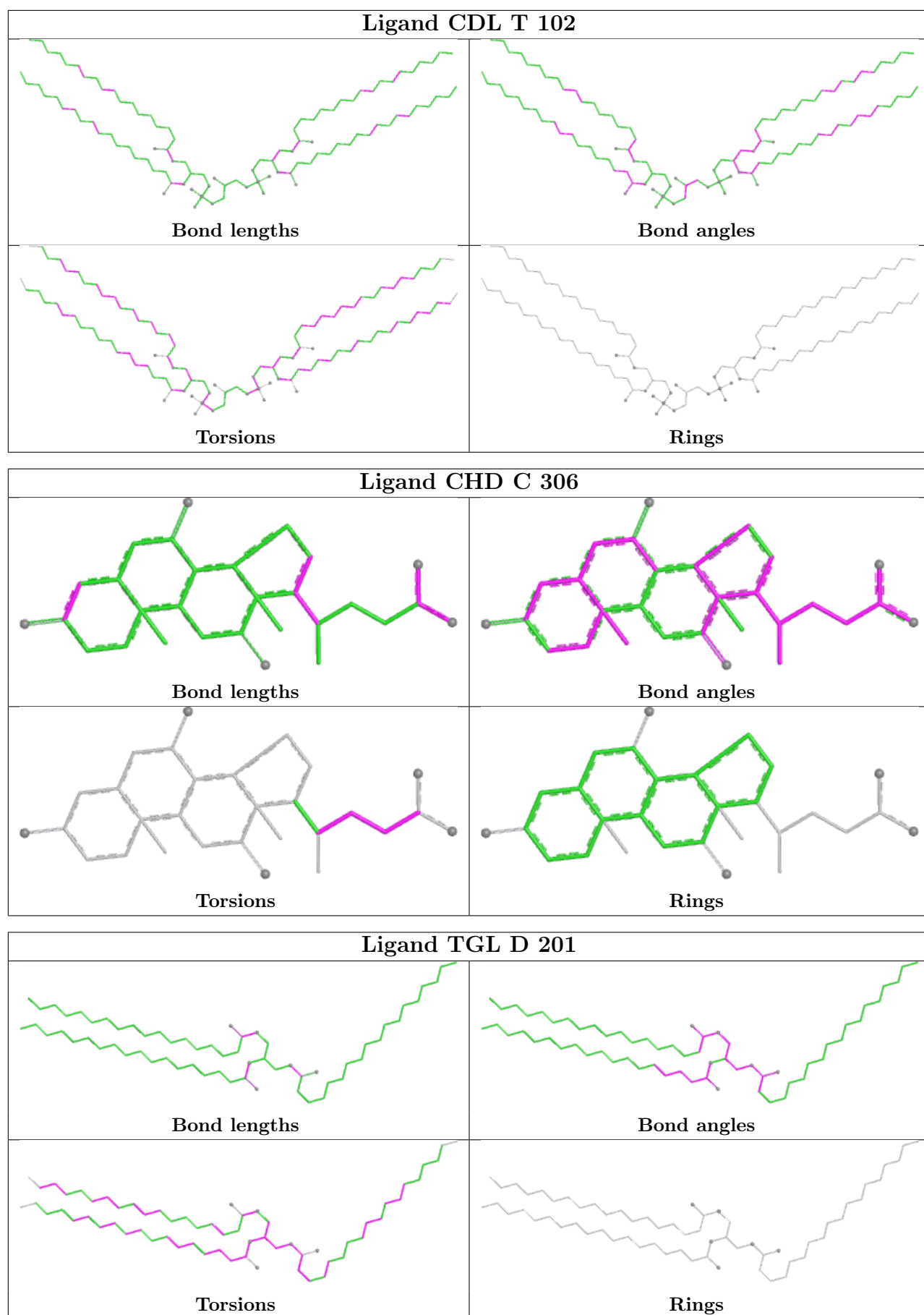


Ligand PEK P 308



Ligand PGV P 302





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 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	513/514 (99%)	-0.21	5 (0%) 79 84	12, 27, 34, 70	22 (4%)
1	N	513/514 (99%)	0.10	9 (1%) 67 73	14, 31, 39, 74	20 (3%)
2	B	226/227 (99%)	0.57	28 (12%) 8 9	19, 34, 54, 68	9 (3%)
2	O	226/227 (99%)	0.78	24 (10%) 11 12	21, 41, 65, 88	5 (2%)
3	C	259/261 (99%)	0.04	7 (2%) 56 61	13, 30, 40, 75	9 (3%)
3	P	259/261 (99%)	0.06	7 (2%) 56 61	12, 31, 41, 80	9 (3%)
4	D	144/147 (97%)	0.45	8 (5%) 30 34	16, 35, 57, 79	4 (2%)
4	Q	144/147 (97%)	1.25	17 (11%) 9 9	19, 48, 74, 136	3 (2%)
5	E	105/109 (96%)	0.33	3 (2%) 53 58	27, 34, 57, 118	0
5	R	105/109 (96%)	0.63	5 (4%) 35 39	24, 40, 58, 121	1 (0%)
6	F	98/98 (100%)	0.86	11 (11%) 10 11	17, 37, 90, 148	4 (4%)
6	S	98/98 (100%)	1.22	15 (15%) 5 5	18, 37, 99, 143	2 (2%)
7	G	83/85 (97%)	1.47	20 (24%) 2 2	17, 39, 108, 141	1 (1%)
7	T	83/85 (97%)	1.46	22 (26%) 1 1	19, 39, 101, 131	1 (1%)
8	H	79/85 (92%)	1.15	13 (16%) 4 4	33, 44, 91, 101	0
8	U	79/85 (92%)	1.18	15 (18%) 3 3	37, 47, 103, 123	0
9	I	72/73 (98%)	1.47	18 (25%) 2 1	33, 44, 83, 90	0
9	V	72/73 (98%)	1.50	18 (25%) 2 1	32, 54, 80, 104	0
10	J	58/59 (98%)	0.78	4 (6%) 23 26	30, 39, 63, 121	0
10	W	58/59 (98%)	0.92	5 (8%) 16 18	22, 42, 69, 127	1 (1%)
11	K	49/56 (87%)	0.74	2 (4%) 41 45	34, 41, 54, 61	0
11	X	49/56 (87%)	1.23	9 (18%) 3 4	23, 50, 70, 81	1 (2%)
12	L	46/47 (97%)	0.44	5 (10%) 10 11	28, 33, 52, 91	0
12	Y	46/47 (97%)	0.92	4 (8%) 16 17	23, 40, 65, 114	1 (2%)

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	M	43/46 (93%)	0.58	5 (11%) 9 10	30, 33, 68, 118	0
13	Z	43/46 (93%)	1.05	6 (13%) 6 7	39, 44, 78, 108	0
All	All	3550/3614 (98%)	0.51	285 (8%) 18 21	12, 34, 66, 148	93 (2%)

The worst 5 of 285 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	Q	5	VAL	14.7
6	F	1	ALA	13.5
6	S	97	ALA	12.4
6	S	1	ALA	10.4
6	S	94	HIS	9.2

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	TPO	G	11	11/12	0.51	0.26	90,110,131,132	0
9	SAC	V	1	9/10	0.59	0.21	102,119,126,133	0
9	SAC	I	1	9/10	0.73	0.22	60,74,81,81	0
7	TPO	T	11	11/12	0.74	0.22	94,107,122,130	0
1	FME	A	1	10/11	0.91	0.15	38,45,74,86	0
1	FME	N	1	10/11	0.92	0.14	42,51,69,79	0
2	FME	O	1	10/11	0.94	0.11	38,39,48,53	0
2	FME	B	1	10/11	0.96	0.10	29,32,42,66	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column

labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
22	CHD	W	101	29/29	0.65	0.24	61,91,115,121	0
22	CHD	J	101	29/29	0.70	0.20	51,75,90,93	0
28	PEK	C	309	53/53	0.70	0.28	48,90,151,153	0
27	CDL	T	102	100/100	0.74	0.28	49,89,137,155	0
24	PSC	B	303	52/52	0.74	0.27	36,84,154,154	0
28	PEK	G	103	53/53	0.74	0.23	48,86,146,153	0
28	PEK	C	307	53/53	0.75	0.26	38,77,135,141	0
25	DMU	C	310	33/33	0.75	0.17	52,73,102,105	0
20	PGV	C	308	51/51	0.75	0.26	44,74,136,146	0
20	PGV	P	302	51/51	0.76	0.25	51,81,132,151	0
27	CDL	N	601	100/100	0.76	0.28	54,90,129,155	0
27	CDL	P	305	100/100	0.77	0.27	36,82,118,137	0
19	TGL	Y	101	63/63	0.78	0.27	45,74,111,138	0
25	DMU	P	309	33/33	0.78	0.17	45,72,92,97	0
19	TGL	Q	201	63/63	0.78	0.27	52,75,93,100	0
24	PSC	N	612	52/52	0.78	0.28	42,83,154,156	0
25	DMU	P	307	33/33	0.79	0.16	38,67,118,121	0
21	EDO	A	613	4/4	0.79	0.21	49,53,56,57	0
28	PEK	P	308	53/53	0.79	0.25	41,72,131,139	0
19	TGL	A	608	63/63	0.80	0.23	40,75,93,101	0
25	DMU	P	310	33/33	0.80	0.16	62,79,108,112	0
25	DMU	C	302	33/33	0.80	0.19	30,73,105,109	0
20	PGV	N	609	51/51	0.81	0.21	43,77,115,131	0
19	TGL	N	611	63/63	0.81	0.24	51,77,97,104	0
19	TGL	A	611	63/63	0.81	0.25	34,60,92,110	0
21	EDO	L	101	4/4	0.81	0.23	55,62,78,81	0
20	PGV	A	610	51/51	0.82	0.23	34,75,106,125	0
27	CDL	C	305	100/100	0.82	0.22	35,75,109,115	0
19	TGL	D	201	63/63	0.82	0.24	32,62,89,92	0
25	DMU	C	311	33/33	0.84	0.15	46,77,99,109	0
22	CHD	P	306	29/29	0.86	0.14	43,49,52,68	0
22	CHD	C	306	29/29	0.86	0.14	43,49,54,66	0
21	EDO	S	104	4/4	0.86	0.14	40,52,60,65	0
21	EDO	W	102	4/4	0.87	0.21	50,55,66,73	0
21	EDO	C	312	4/4	0.87	0.20	41,43,72,85	0
21	EDO	A	617	4/4	0.87	0.17	40,42,43,44	0
26	UNX	C	303	1/1	0.87	0.25	35,35,35,35	0
21	EDO	N	620	4/4	0.87	0.18	47,49,54,58	0
21	EDO	A	618	4/4	0.87	0.19	51,51,53,63	0
21	EDO	Y	102	4/4	0.88	0.17	61,65,66,67	0
21	EDO	N	616	4/4	0.88	0.14	52,54,55,56	0
25	DMU	Z	101	33/33	0.89	0.13	45,54,70,76	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
21	EDO	N	617	4/4	0.89	0.16	52,55,58,61	0
21	EDO	N	621	4/4	0.90	0.17	41,51,53,61	0
21	EDO	A	619	4/4	0.90	0.19	44,60,74,82	0
21	EDO	B	305	4/4	0.91	0.17	35,45,48,51	0
21	EDO	D	202	4/4	0.91	0.15	33,38,47,66	0
21	EDO	G	105	4/4	0.91	0.13	53,56,64,79	0
21	EDO	A	614	4/4	0.92	0.12	30,31,32,37	0
21	EDO	A	620	4/4	0.92	0.15	52,56,57,64	0
21	EDO	M	102	4/4	0.92	0.18	60,63,63,69	0
21	EDO	B	304	4/4	0.92	0.14	46,59,60,85	0
21	EDO	A	615	4/4	0.93	0.14	53,63,65,70	0
21	EDO	B	306	4/4	0.93	0.15	50,51,55,57	0
25	DMU	M	101	33/33	0.93	0.11	38,44,58,67	0
21	EDO	P	312	4/4	0.93	0.13	33,36,43,49	0
21	EDO	S	103	4/4	0.93	0.12	39,42,45,47	0
21	EDO	D	203	4/4	0.93	0.19	63,67,71,74	0
21	EDO	E	202	4/4	0.93	0.11	38,40,44,47	0
21	EDO	E	203	4/4	0.93	0.11	42,42,50,55	0
21	EDO	N	615	4/4	0.94	0.11	41,53,57,60	0
21	EDO	O	302	4/4	0.94	0.11	38,39,40,42	0
21	EDO	F	104	4/4	0.94	0.11	36,36,38,42	0
21	EDO	F	103	4/4	0.94	0.13	37,41,42,52	0
21	EDO	N	619	4/4	0.94	0.12	40,43,46,49	0
21	EDO	N	614	4/4	0.94	0.11	29,33,36,36	0
28	PEK	T	101	53/53	0.94	0.15	29,49,85,92	0
18	AZI	A	606[B]	3/3	0.95	0.11	22,22,23,24	3
21	EDO	P	311	4/4	0.95	0.10	39,42,43,46	0
21	EDO	A	616	4/4	0.95	0.10	22,27,28,42	0
21	EDO	N	618	4/4	0.95	0.10	37,40,43,45	0
21	EDO	N	613	4/4	0.95	0.10	33,34,37,41	0
28	PEK	G	101	53/53	0.95	0.14	29,46,79,95	0
21	EDO	T	103	4/4	0.95	0.11	36,37,42,44	0
18	AZI	N	607[B]	3/3	0.95	0.12	23,23,27,28	3
21	EDO	B	307	4/4	0.95	0.10	29,30,35,37	0
20	PGV	A	609	51/51	0.96	0.10	25,33,67,71	0
22	CHD	C	301	29/29	0.96	0.07	27,30,34,36	0
20	PGV	N	610	51/51	0.96	0.11	27,35,68,72	0
21	EDO	R	201	4/4	0.96	0.09	44,45,45,45	0
22	CHD	P	301	29/29	0.96	0.07	28,32,36,39	0
18	AZI	N	608[A]	3/3	0.96	0.13	29,29,38,40	3
21	EDO	A	612	4/4	0.96	0.10	37,41,41,53	0
20	PGV	C	304	51/51	0.96	0.12	27,34,86,96	0

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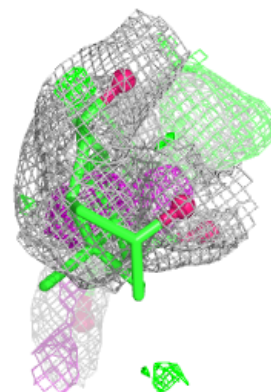
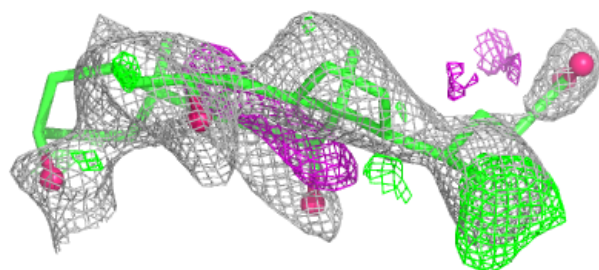
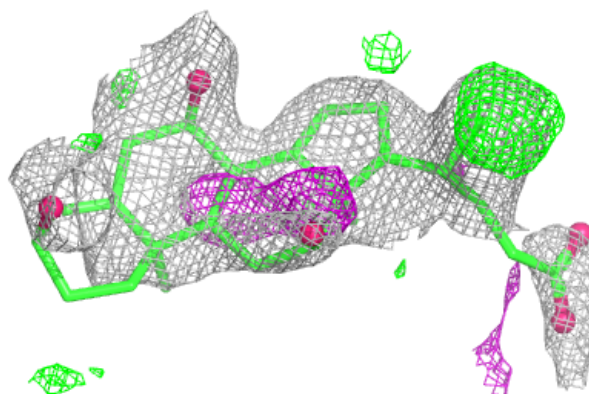
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
18	AZI	N	608[B]	3/3	0.96	0.13	24,24,26,26	3
26	UNX	P	303	1/1	0.96	0.09	32,32,32,32	0
18	AZI	A	607[A]	3/3	0.97	0.09	20,20,24,27	3
18	AZI	A	607[B]	3/3	0.97	0.09	23,23,24,29	3
22	CHD	B	301	29/29	0.97	0.07	25,30,33,41	0
16	MG	N	605	1/1	0.97	0.06	31,31,31,31	0
20	PGV	P	304	51/51	0.97	0.09	26,36,79,89	0
22	CHD	G	102	29/29	0.97	0.06	28,30,36,39	0
21	EDO	G	104	4/4	0.97	0.07	34,37,38,41	0
21	EDO	E	201	4/4	0.97	0.07	40,41,43,44	0
14	HEA	A	601	60/60	0.98	0.07	21,24,48,50	0
17	NA	N	606	1/1	0.98	0.07	35,35,35,35	0
14	HEA	A	602[A]	60/60	0.98	0.07	18,22,26,28	60
14	HEA	A	602[B]	60/60	0.98	0.07	21,25,34,37	60
21	EDO	S	102	4/4	0.98	0.07	29,29,30,30	0
14	HEA	N	602	60/60	0.98	0.08	26,30,49,52	0
14	HEA	N	603[A]	60/60	0.98	0.07	21,25,28,30	60
14	HEA	N	603[B]	60/60	0.98	0.07	24,29,42,46	60
21	EDO	F	102	4/4	0.99	0.06	27,27,28,30	0
23	CUA	O	301	2/2	0.99	0.02	32,32,32,32	0
16	MG	A	604	1/1	0.99	0.02	26,26,26,26	0
17	NA	A	605	1/1	0.99	0.04	30,30,30,30	0
29	ZN	S	101	1/1	0.99	0.03	33,33,33,33	0
15	CU	N	604	1/1	1.00	0.02	30,30,30,30	0
15	CU	A	603	1/1	1.00	0.01	26,26,26,26	0
29	ZN	F	101	1/1	1.00	0.02	31,31,31,31	0
23	CUA	B	302	2/2	1.00	0.02	27,27,27,28	0

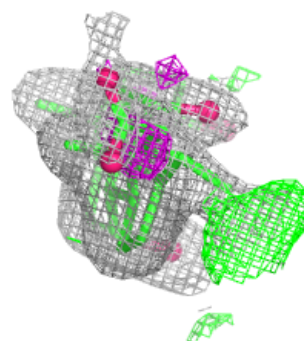
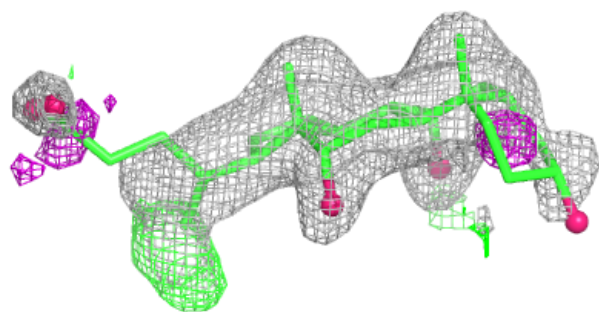
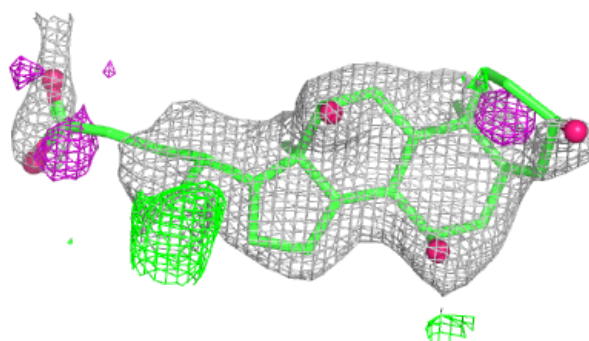
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around CHD W 101:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

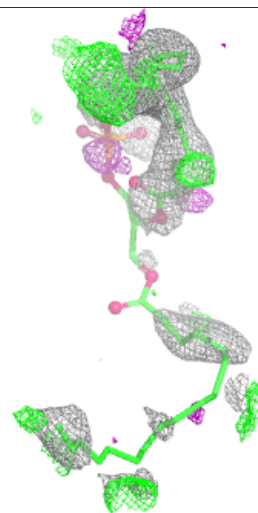
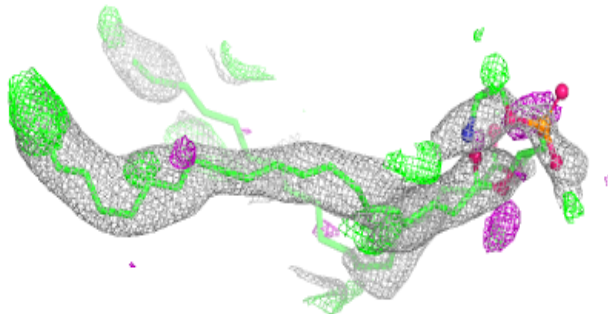
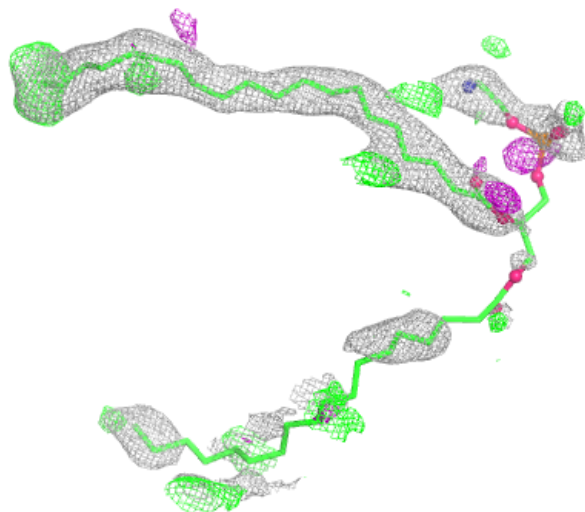
**Electron density around CHD J 101:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



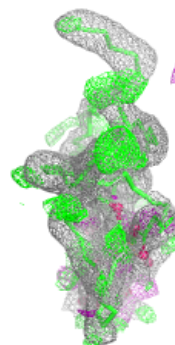
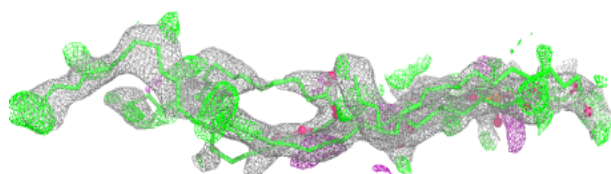
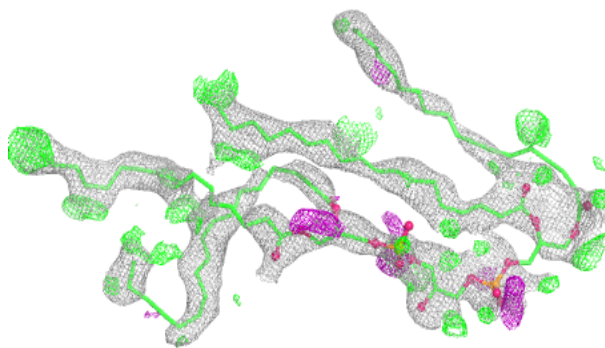
Electron density around PEK C 309:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

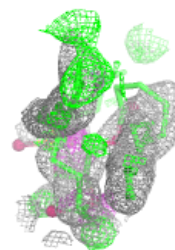
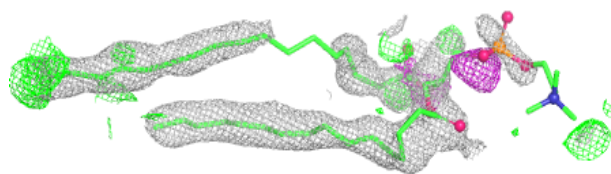
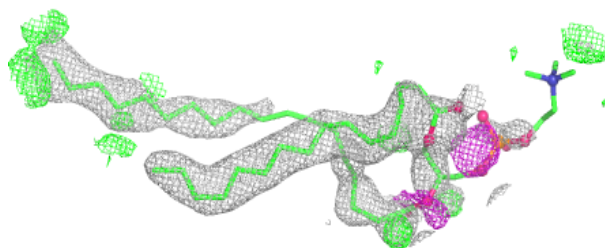


Electron density around CDL T 102:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

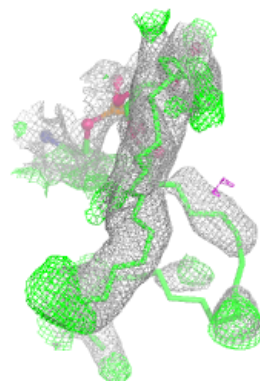
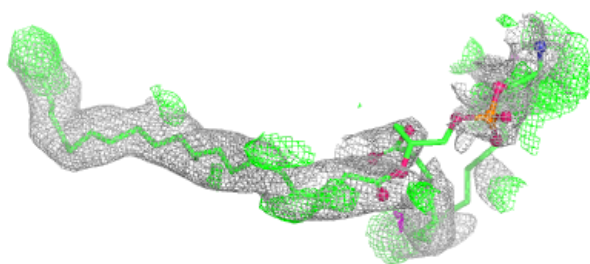
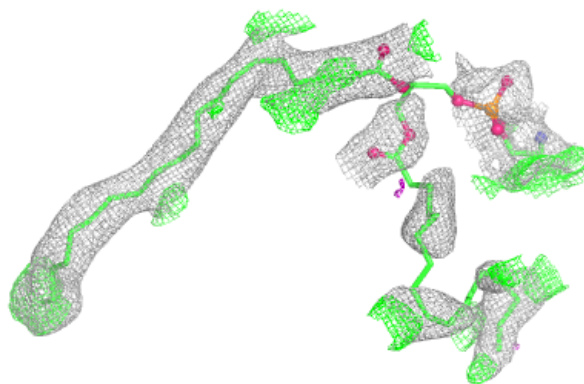
**Electron density around PSC B 303:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



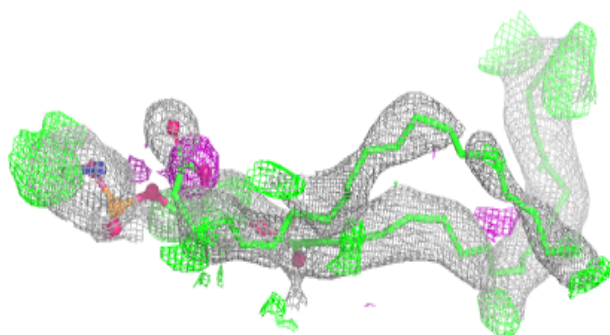
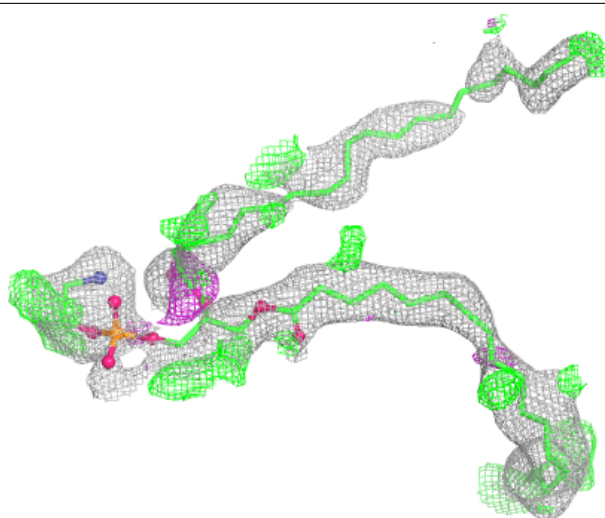
Electron density around PEK G 103:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



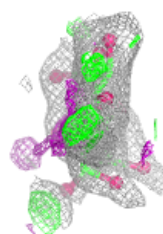
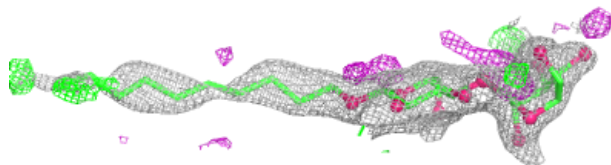
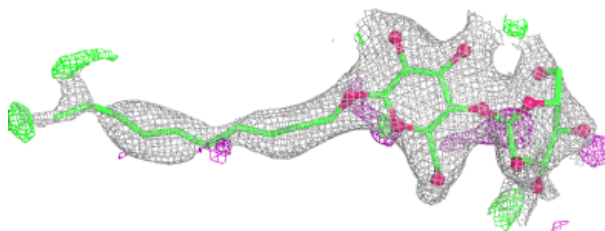
Electron density around PEK C 307:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

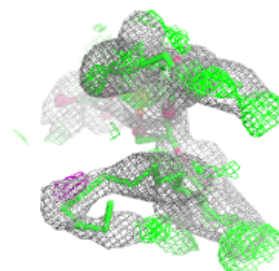
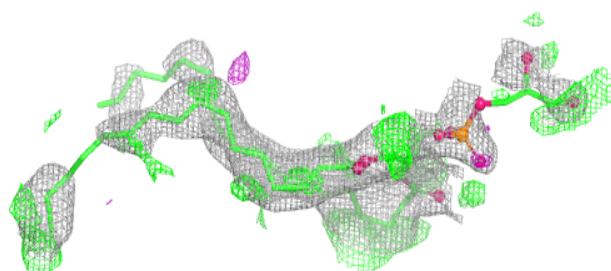
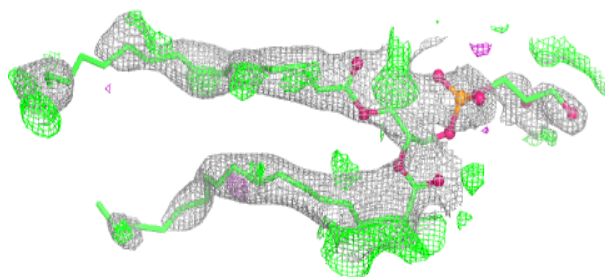


Electron density around DMU C 310:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

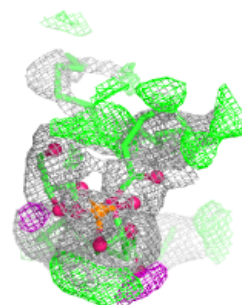
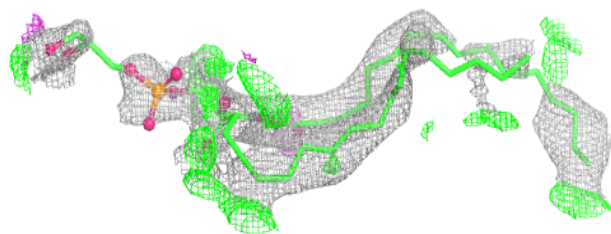
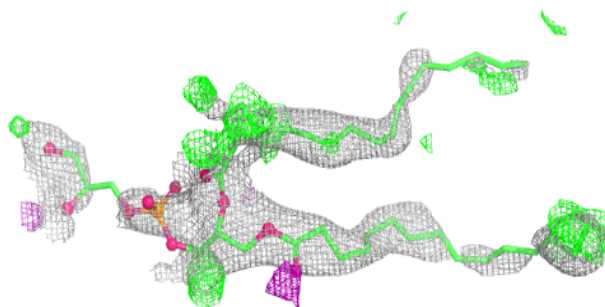
**Electron density around PGV C 308:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

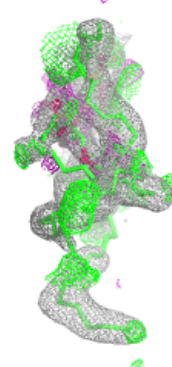
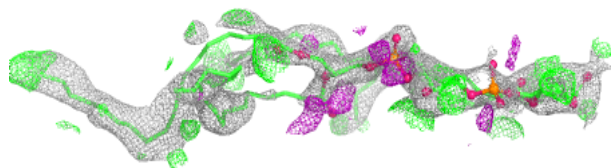
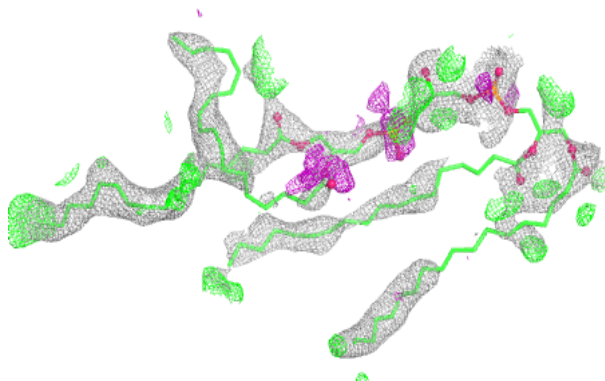


Electron density around PGV P 302:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

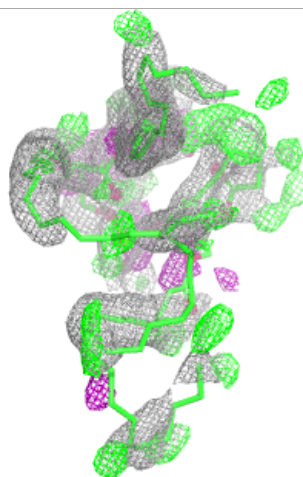
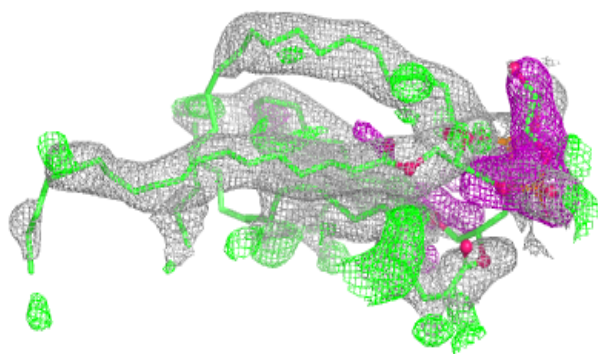
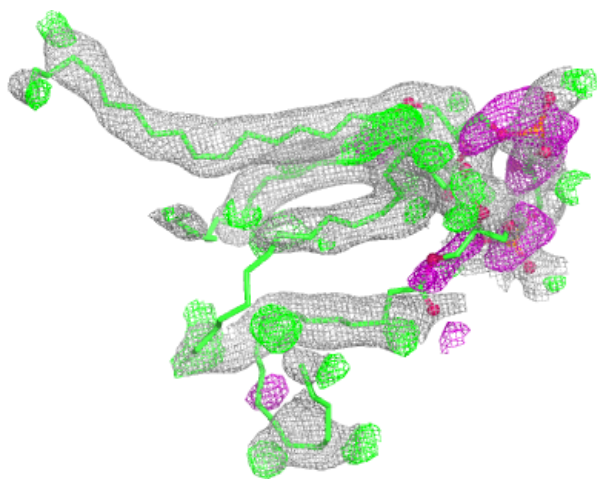
**Electron density around CDL N 601:**

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and green (positive)



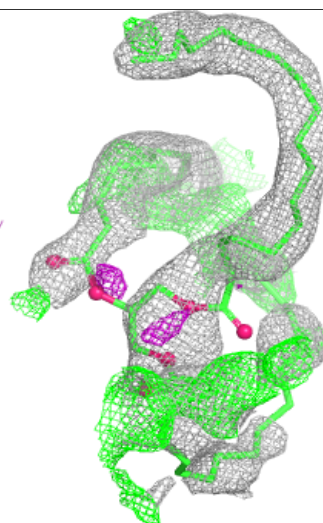
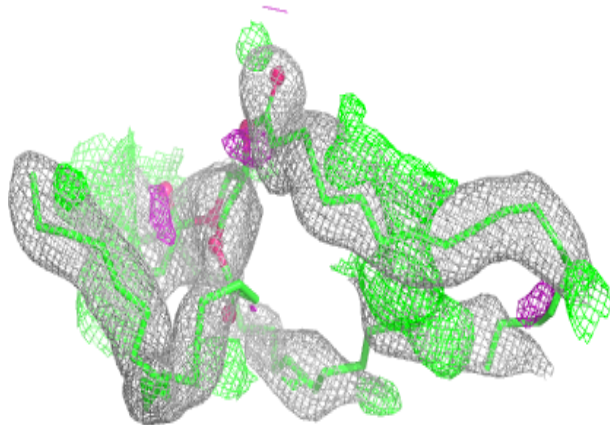
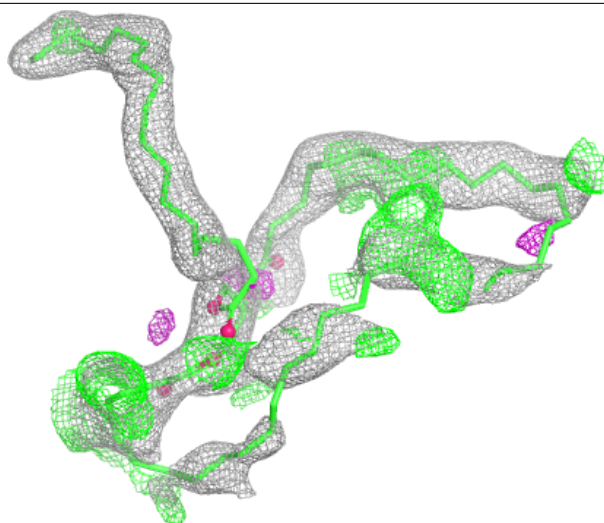
Electron density around CDL P 305:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



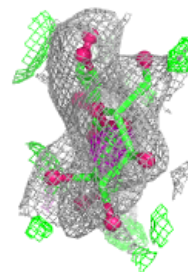
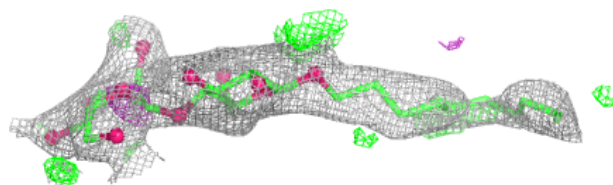
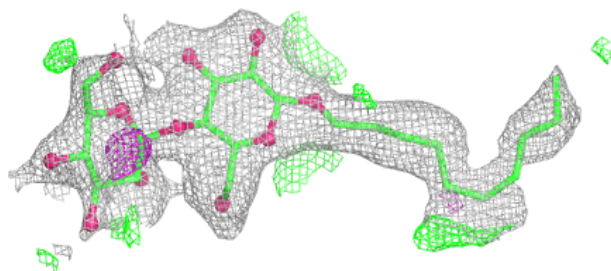
Electron density around TGL Y 101:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

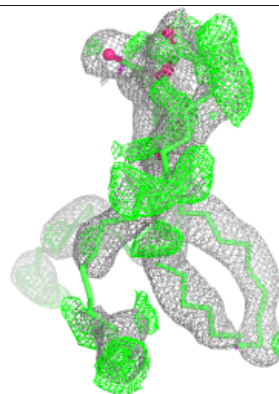
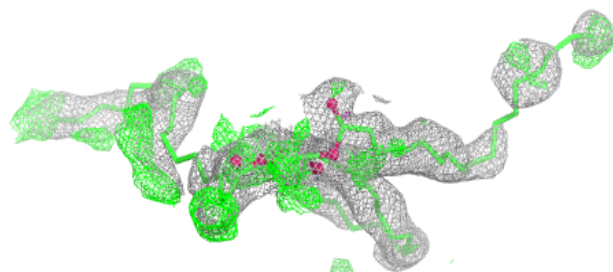
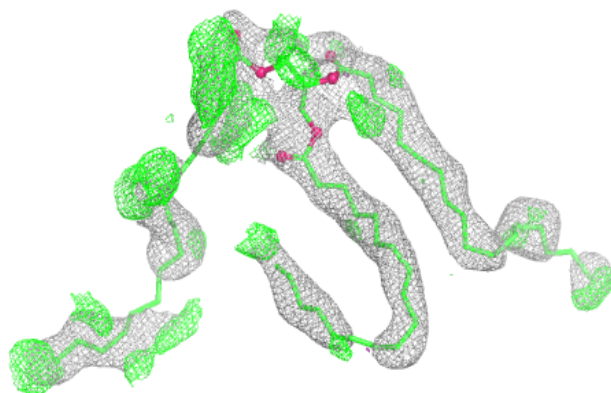


Electron density around DMU P 309:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

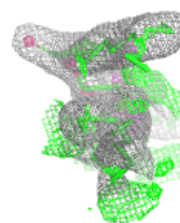
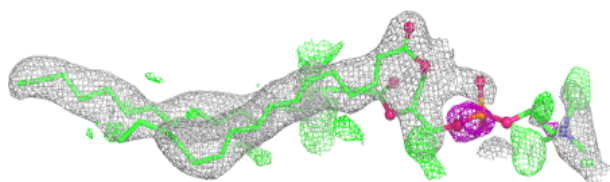
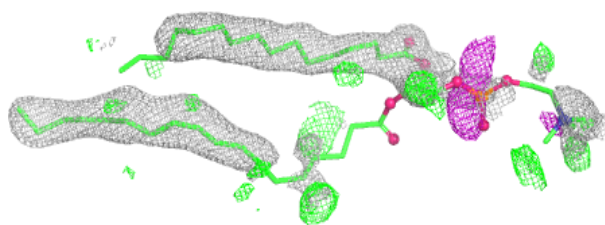
**Electron density around TGL Q 201:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

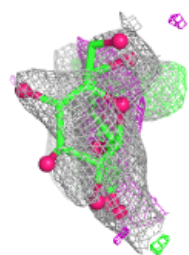
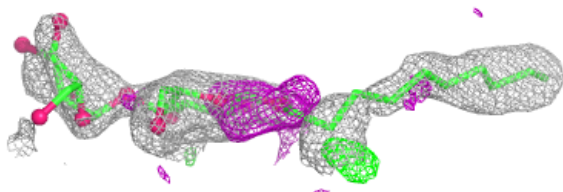
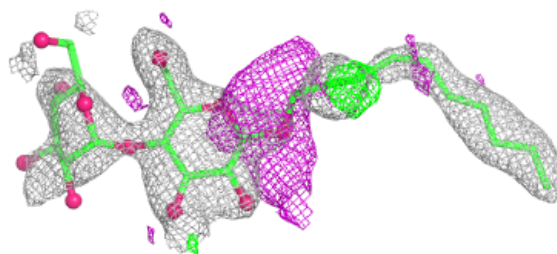


Electron density around PSC N 612:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

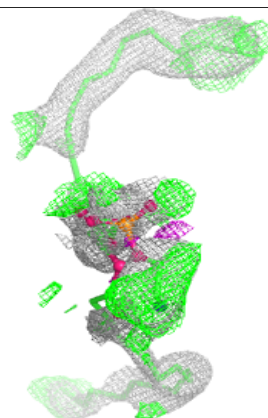
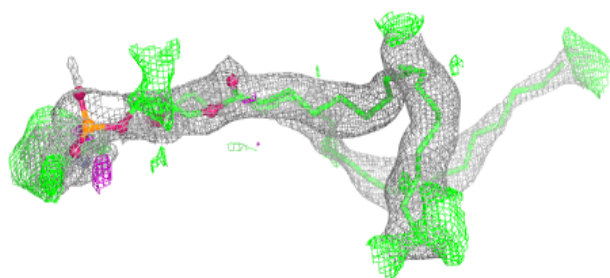
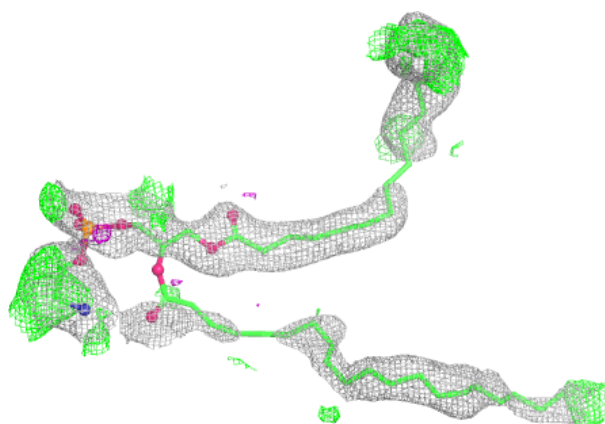
**Electron density around DMU P 307:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

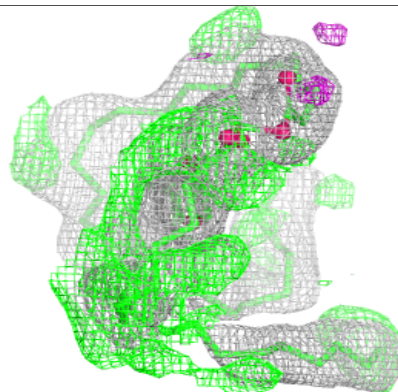
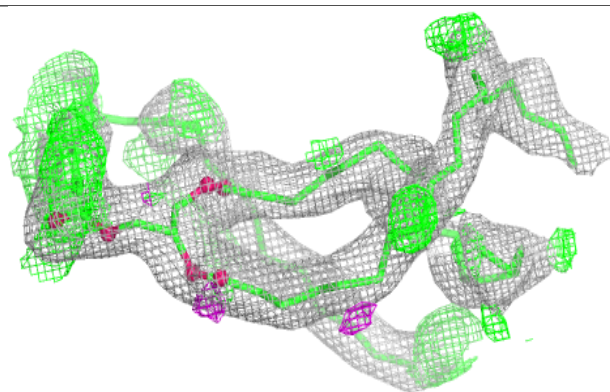
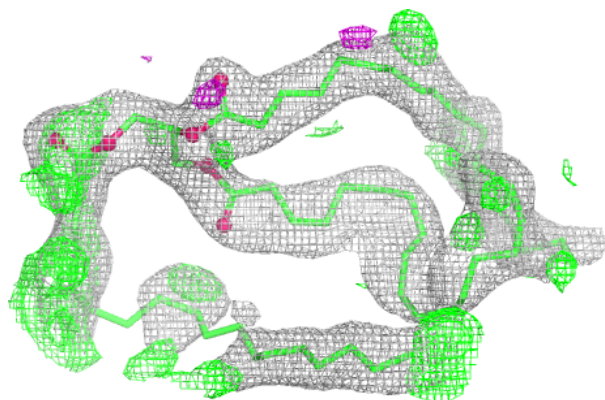


Electron density around PEK P 308:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

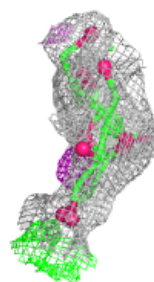
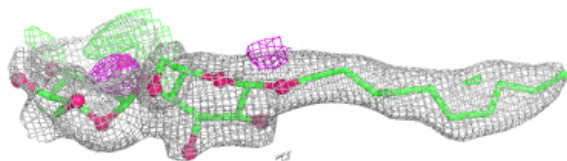
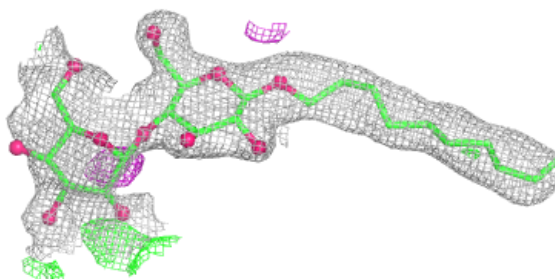
**Electron density around TGL A 608:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

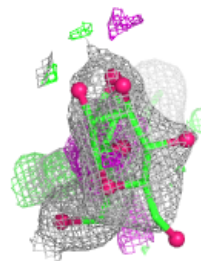
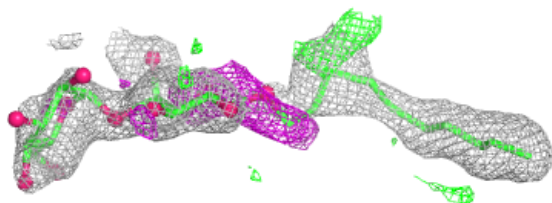
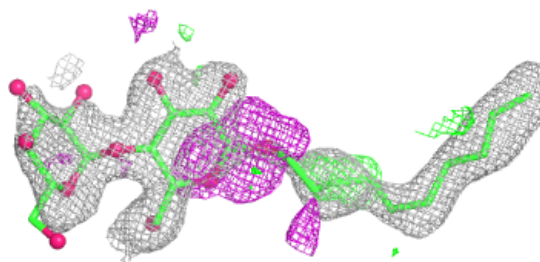


Electron density around DMU P 310:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

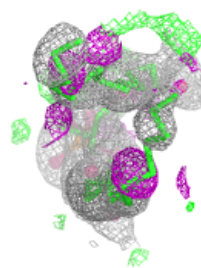
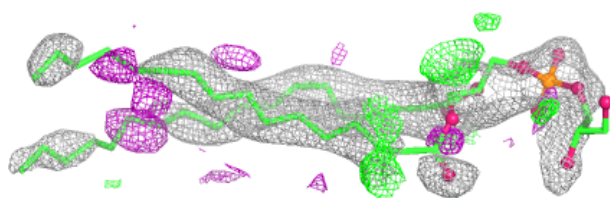
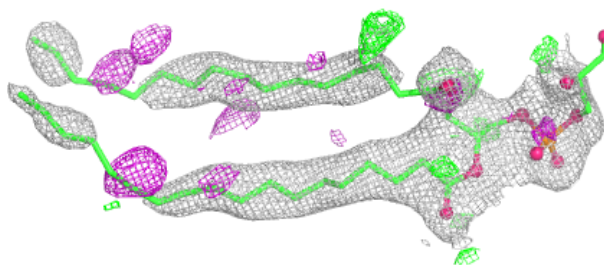
**Electron density around DMU C 302:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

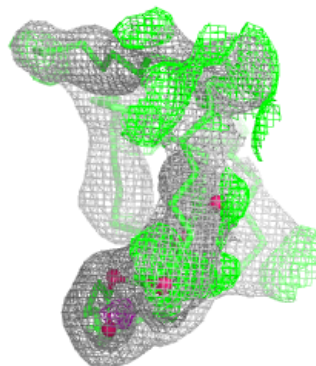
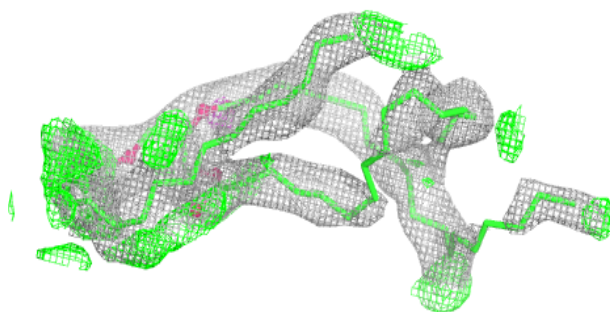
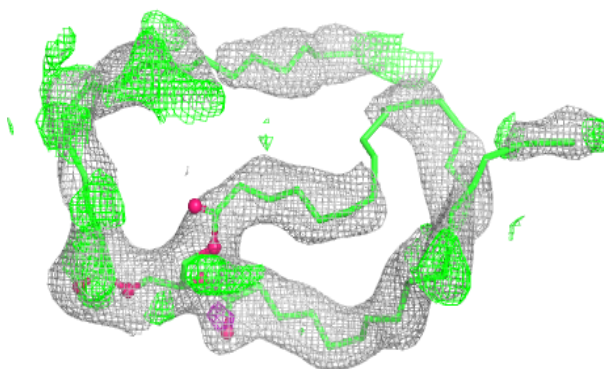


Electron density around PGV N 609:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

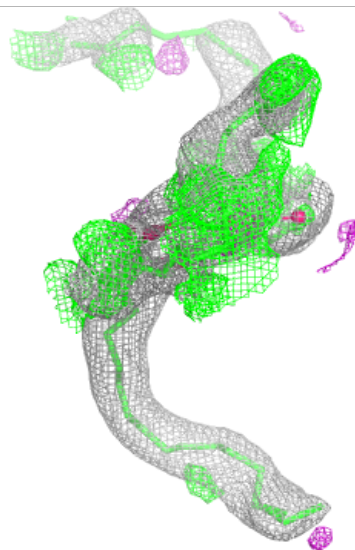
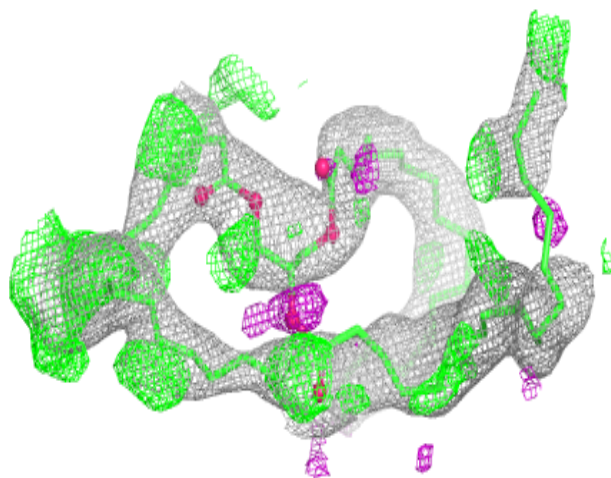
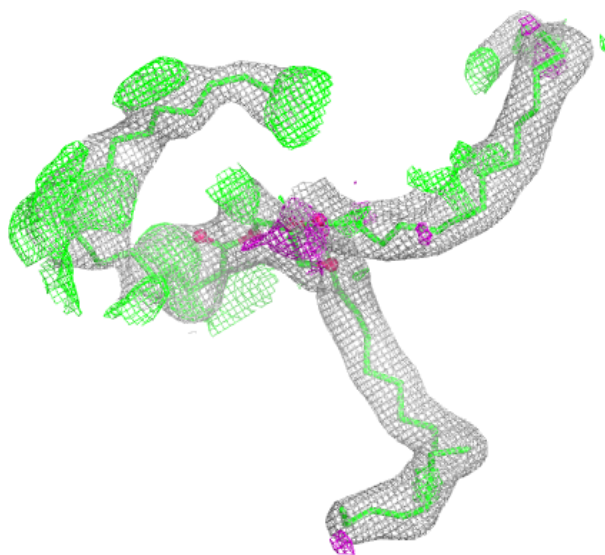
**Electron density around TGL N 611:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



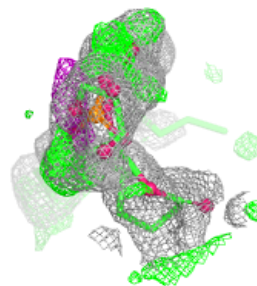
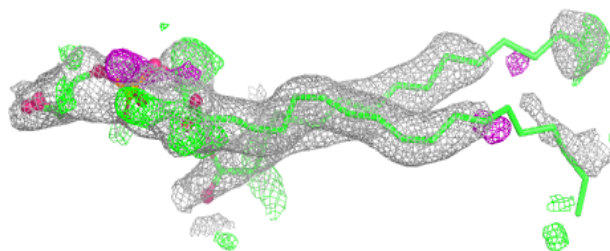
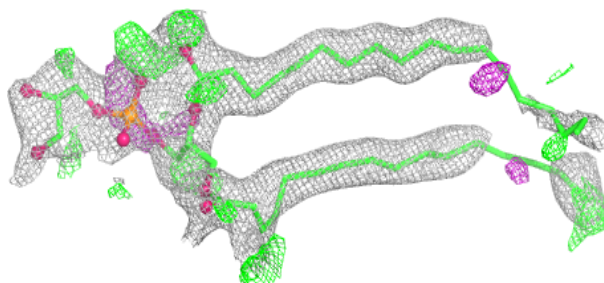
Electron density around TGL A 611:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



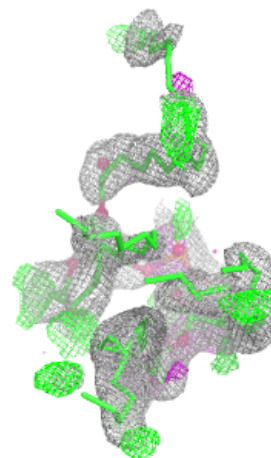
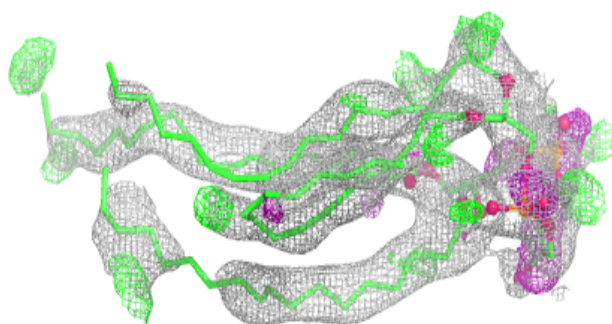
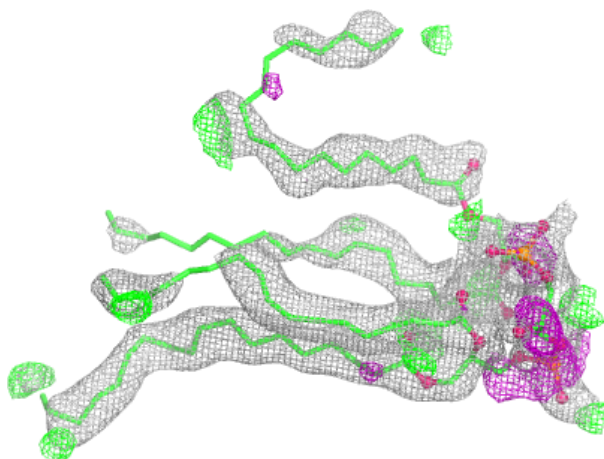
Electron density around PGV A 610:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



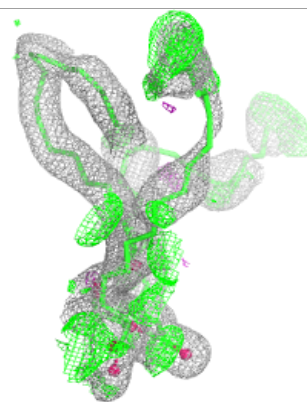
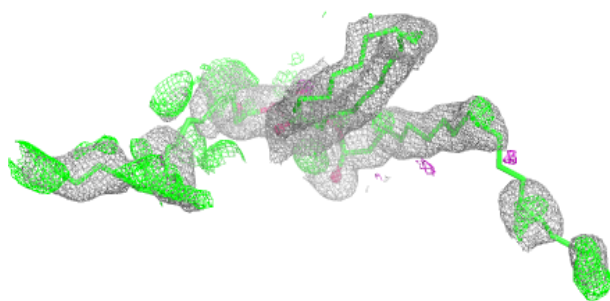
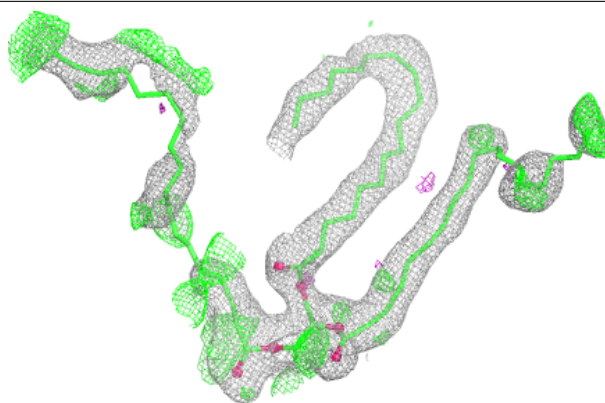
Electron density around CDL C 305:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

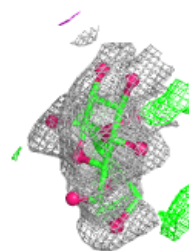
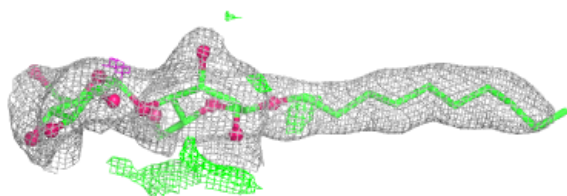
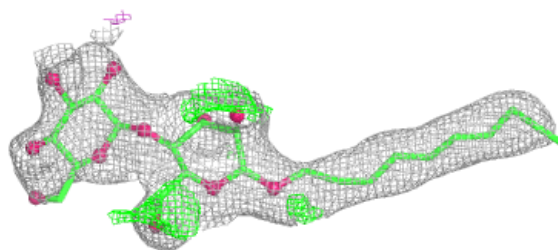


Electron density around TGL D 201:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

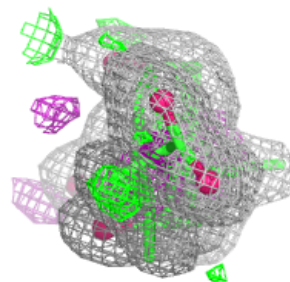
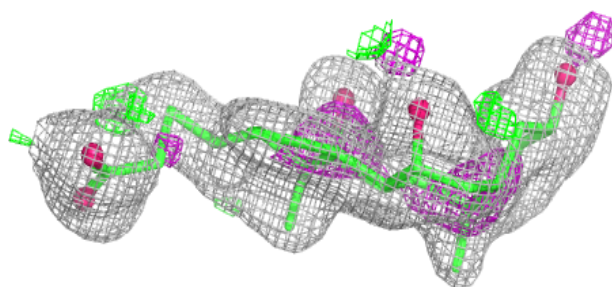
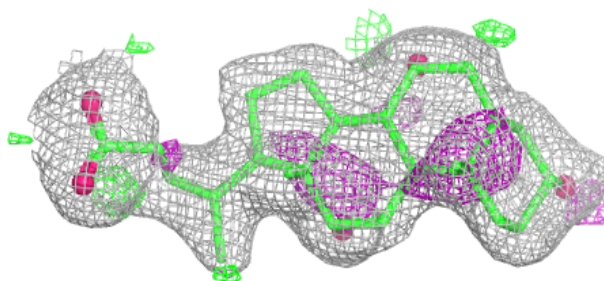
**Electron density around DMU C 311:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

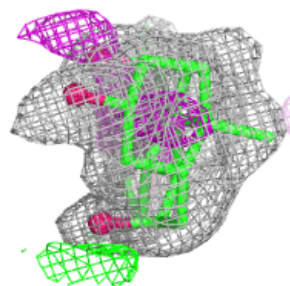
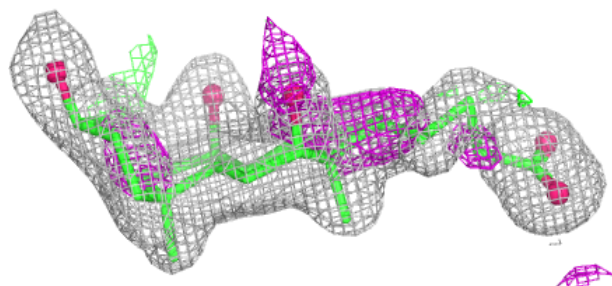
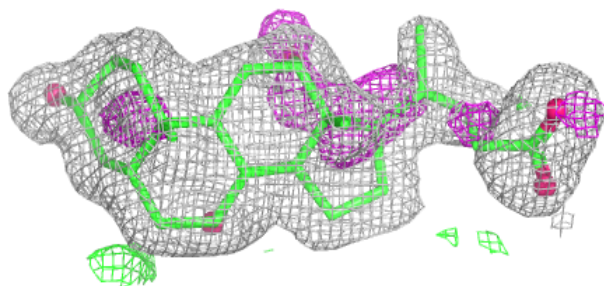


Electron density around CHD P 306:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

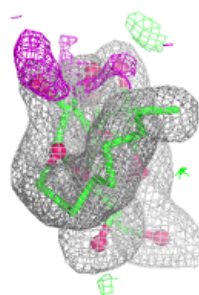
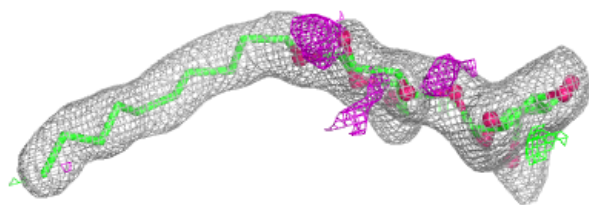
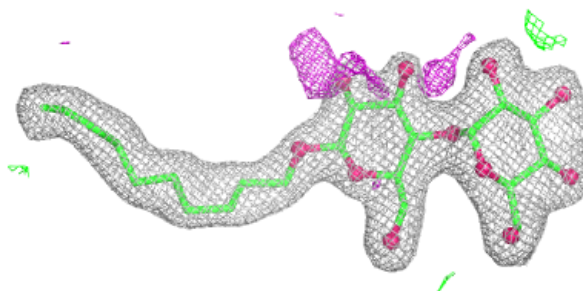
**Electron density around CHD C 306:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

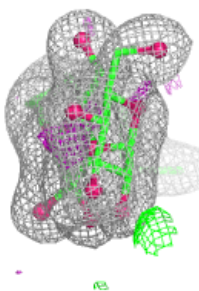
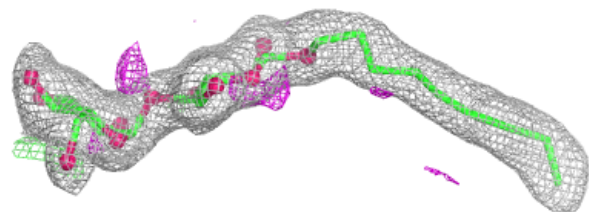
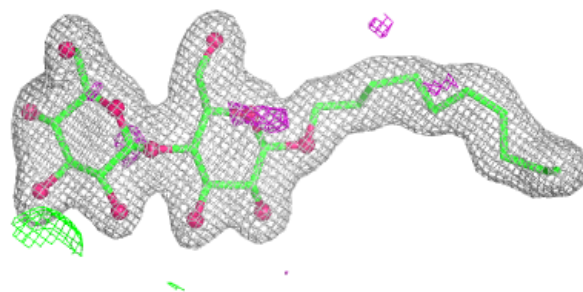


Electron density around DMU Z 101:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

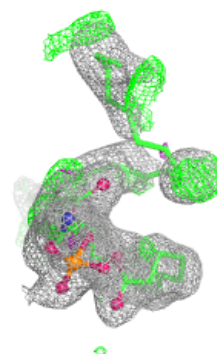
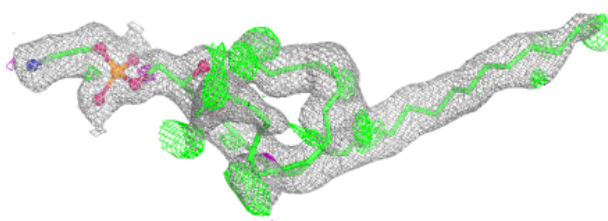
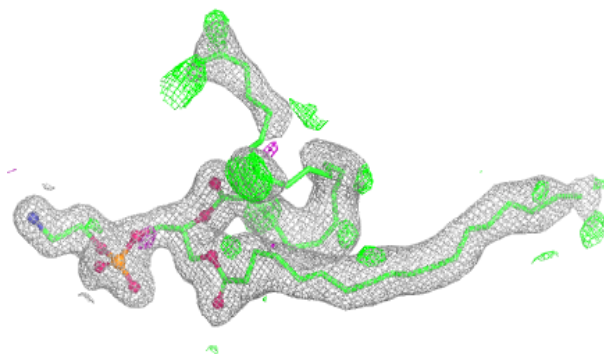
**Electron density around DMU M 101:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

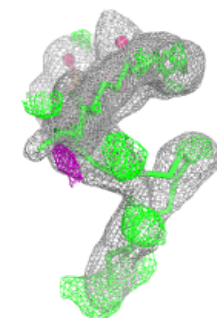
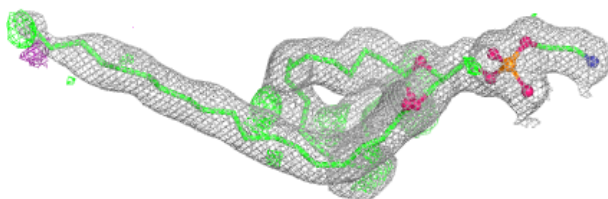
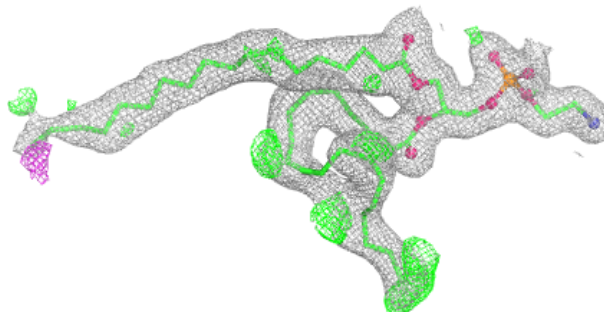


Electron density around PEK T 101:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

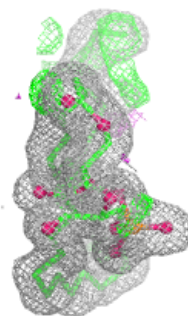
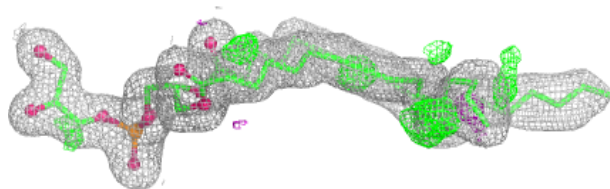
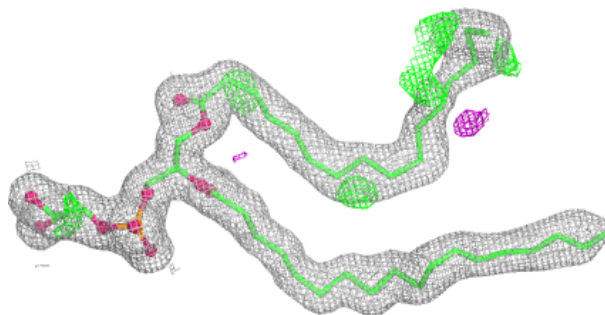
**Electron density around PEK G 101:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

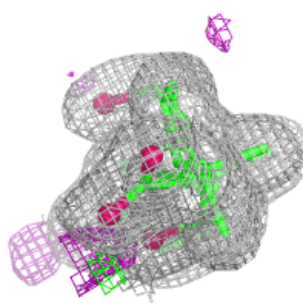
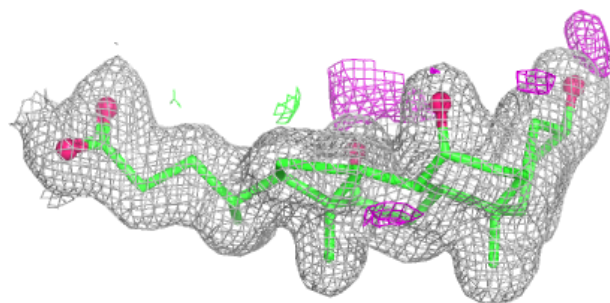
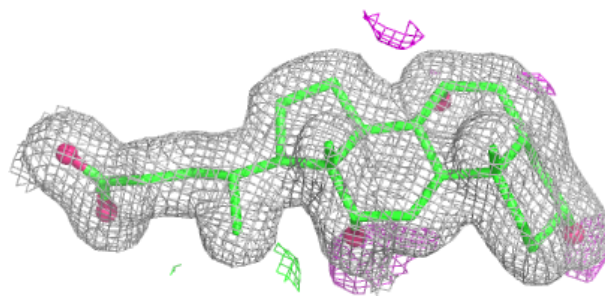


Electron density around PGV A 609:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

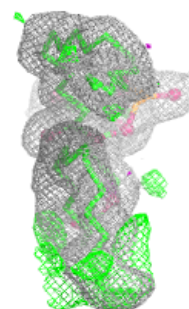
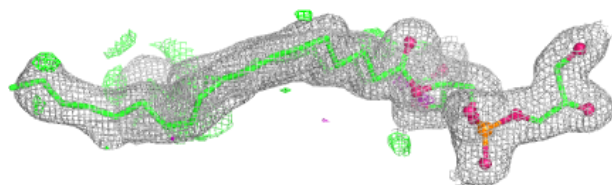
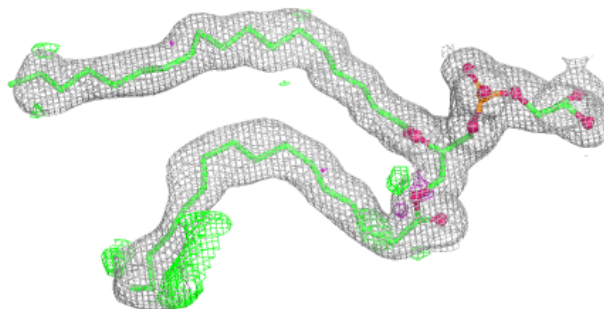
**Electron density around CHD C 301:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

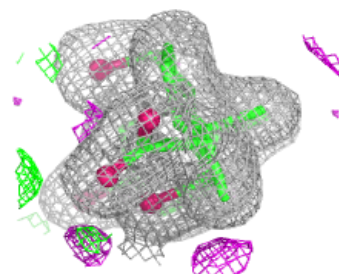
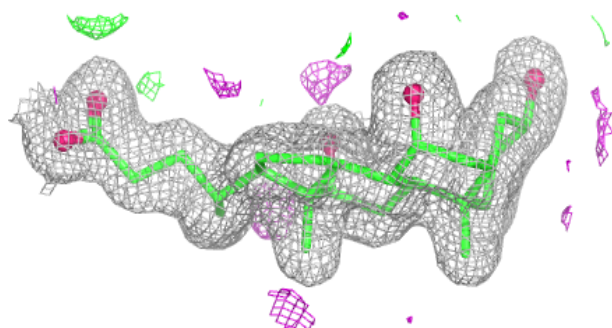
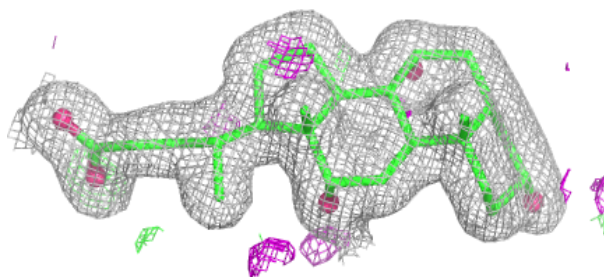


Electron density around PGV N 610:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

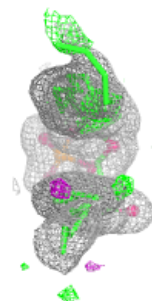
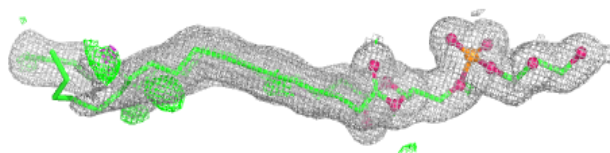
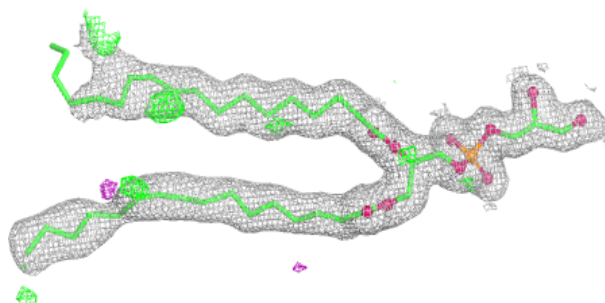
**Electron density around CHD P 301:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

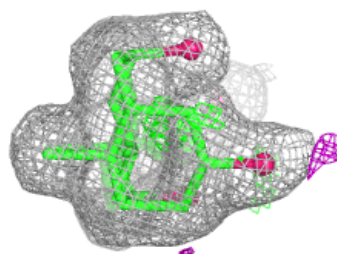
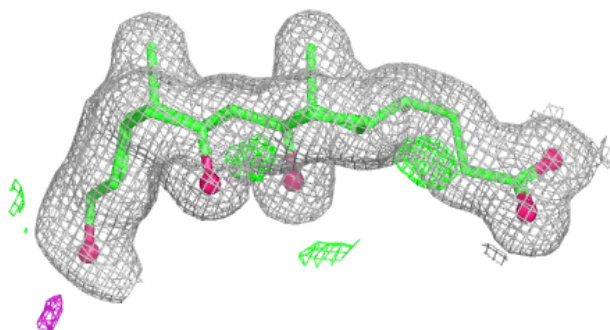
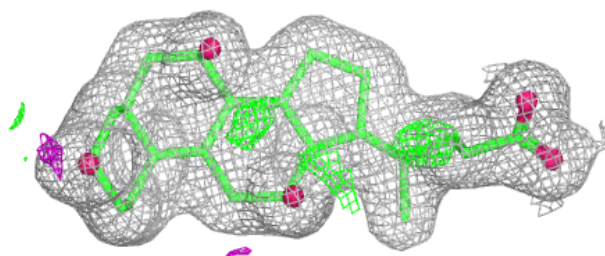


Electron density around PGV C 304:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

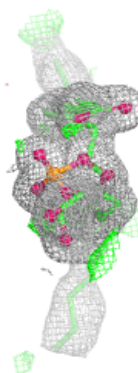
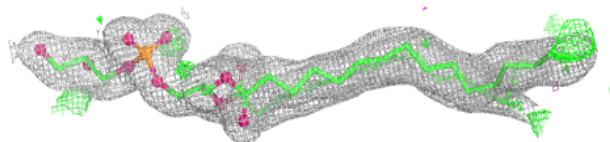
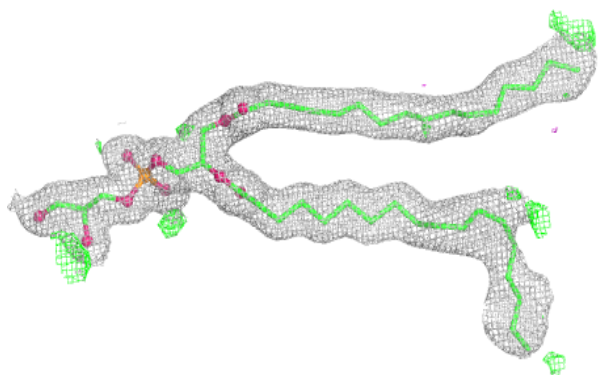
**Electron density around CHD B 301:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

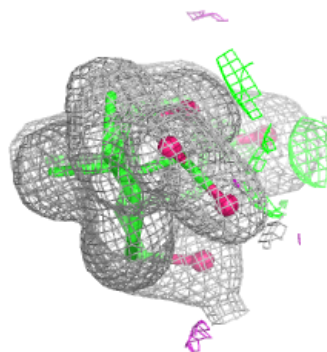
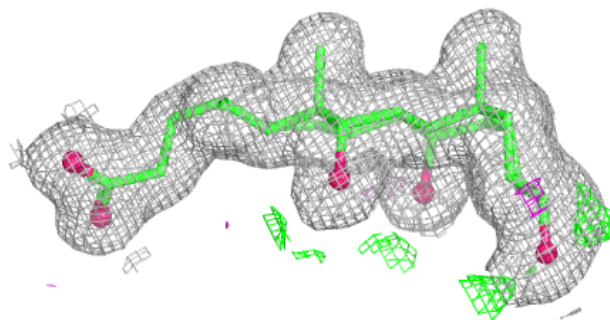
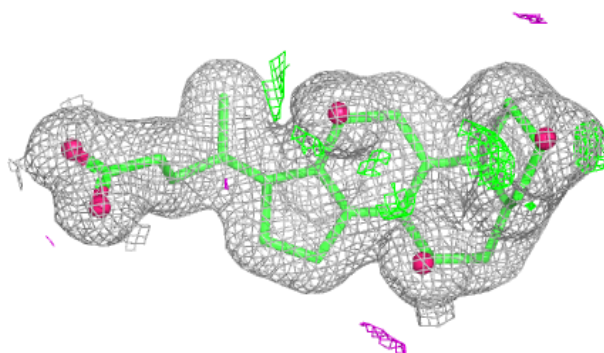


Electron density around PGV P 304:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

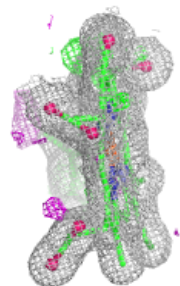
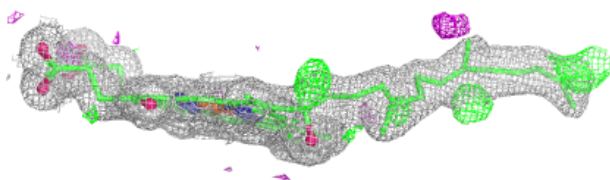
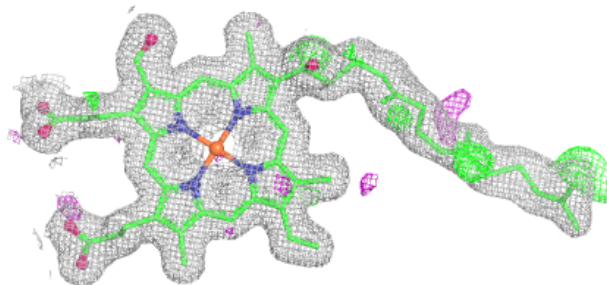
**Electron density around CHD G 102:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

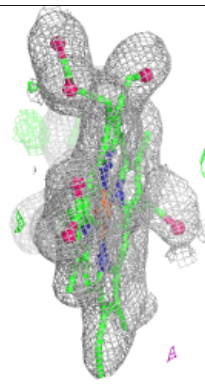
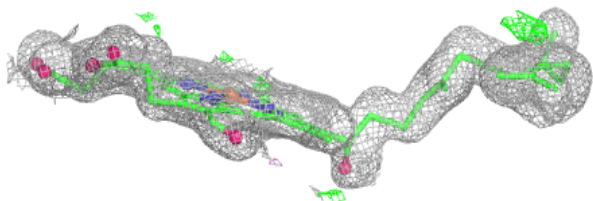
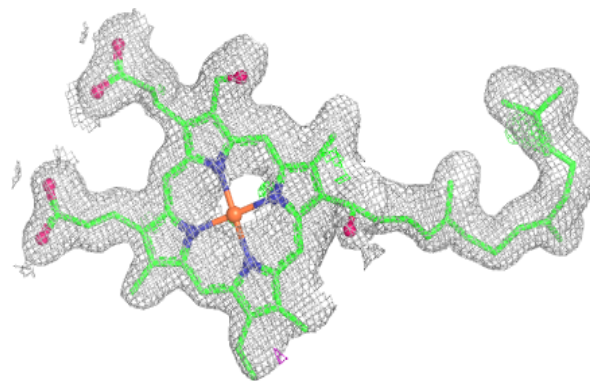


Electron density around HEA A 601:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

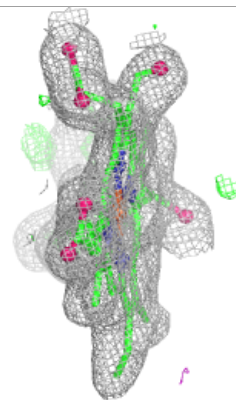
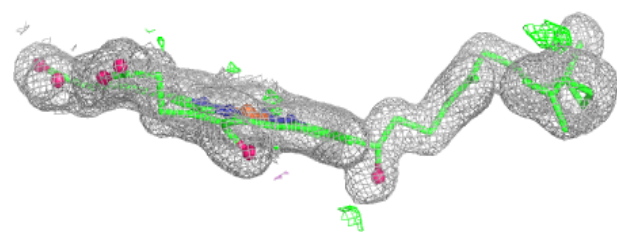
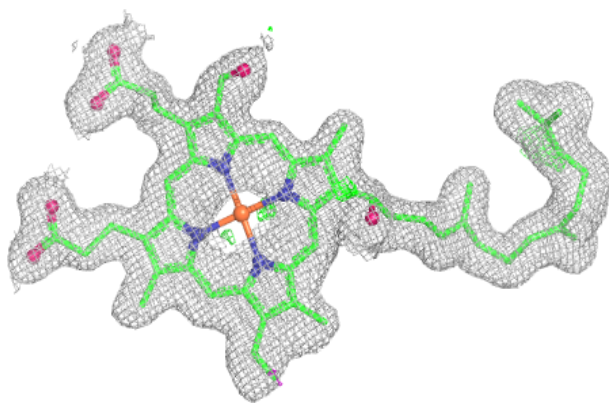
**Electron density around HEA A 602 (A):**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

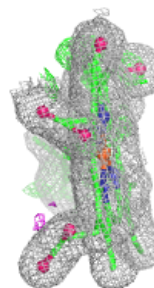
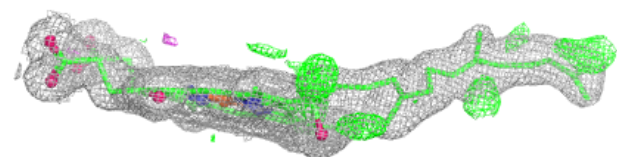
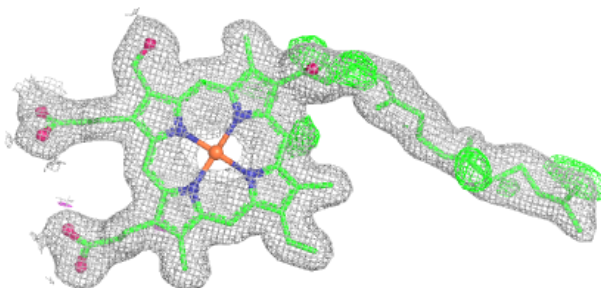


Electron density around HEA A 602 (B):

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

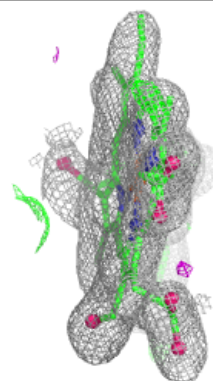
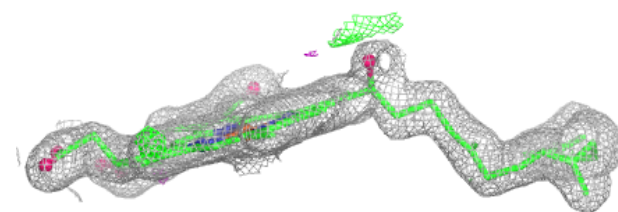
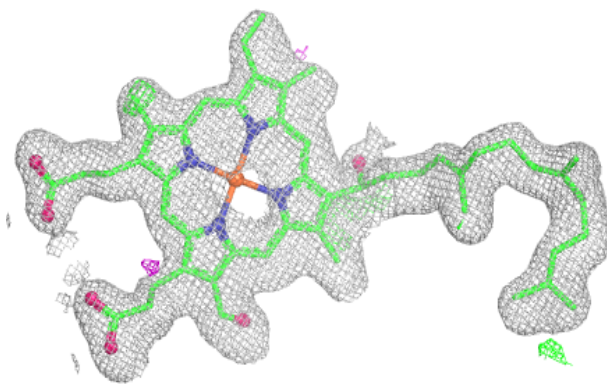
**Electron density around HEA N 602:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

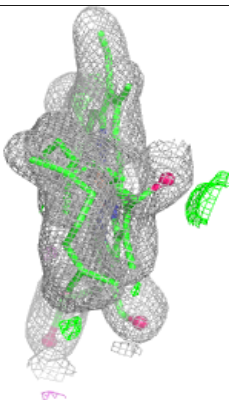
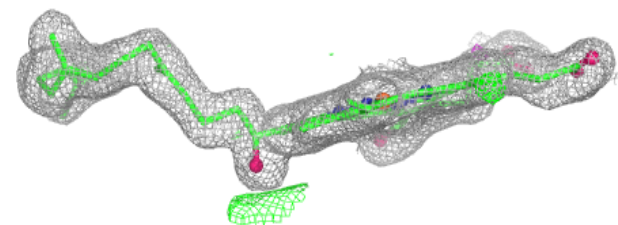
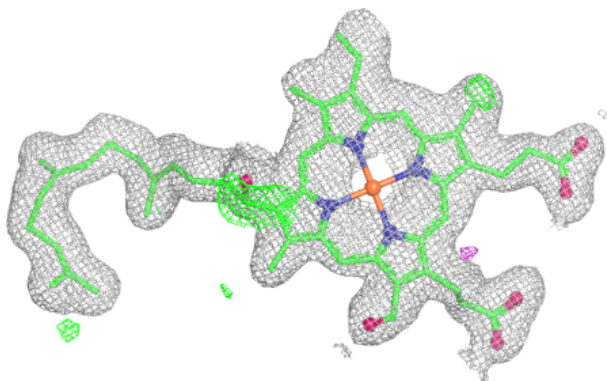


Electron density around HEA N 603 (A):

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around HEA N 603 (B):**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.