



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 10, 2026 – 05:11 PM UTC

PDB ID : 5Z85 / pdb_00005z85
Title : The structure of azide-bound cytochrome c oxidase determined using the another batch crystals exposed to 20 mM azide solution for 2 days
Authors : Shimada, A.; Hatano, K.; Tadehara, H.; Tsukihara, T.
Deposited on : 2018-01-31
Resolution : 1.85 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

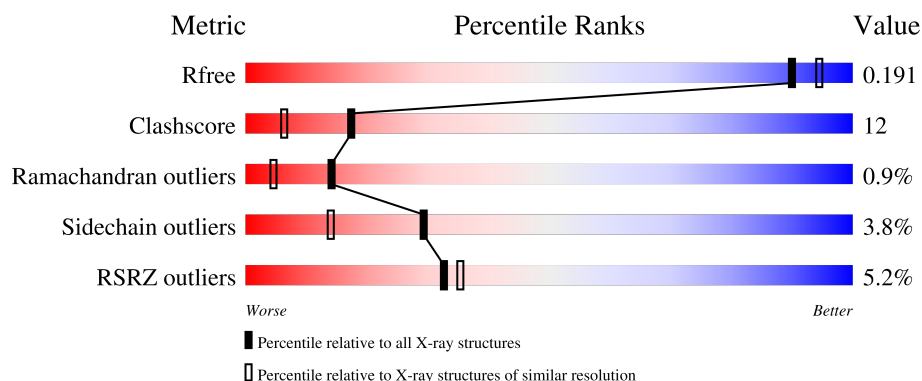
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 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	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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	6% 74% 23% ..
1	N	514	72% 24% .
2	B	227	6% 71% 24% 5%
2	O	227	5% 71% 24% .
3	C	261	2% 71% 26% ..

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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	606	-	X	-	-
18	AZI	A	607	-	X	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
18	AZI	N	606	-	X	-	-
18	AZI	N	607	-	X	X	-
21	EDO	A	612	-	-	X	-
21	EDO	A	614	-	X	X	-
21	EDO	A	618	-	-	X	-
21	EDO	D	202	-	-	X	-
21	EDO	D	203	-	-	X	-
21	EDO	N	618	-	-	X	-
21	EDO	N	621	-	-	X	-
21	EDO	Q	204	-	-	X	-
27	CDL	C	305	-	-	X	-
27	CDL	G	102	-	-	X	-
27	CDL	P	306	-	-	X	-
27	CDL	T	103	-	-	X	-

2 Entry composition

There are 30 unique types of molecules in this entry. The entry contains 33411 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	18	0
			4165	2777	645	703	40			
1	N	514	Total	C	N	O	S	0	18	0
			4160	2774	642	702	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	4	0
			1859	1211	285	344	19			
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	7	0
			2171	1449	347	360	15			
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	1	0
			1204	783	198	219	4			
4	Q	144	Total	C	N	O	S	0	2	0
			1213	791	198	220	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	0	0
			852	544	144	162	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	3	0
			771	477	138	150	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 675	C 431	N 129	O 113	P 1	S 1	0	0	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	0	0
			460	297	78	82	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			

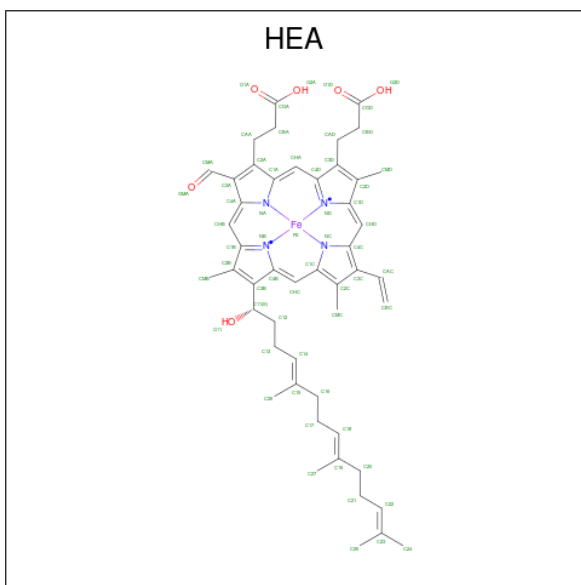
- Molecule 12 is a protein called Cytochrome c oxidase subunit 7C, mitochondrial.

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		

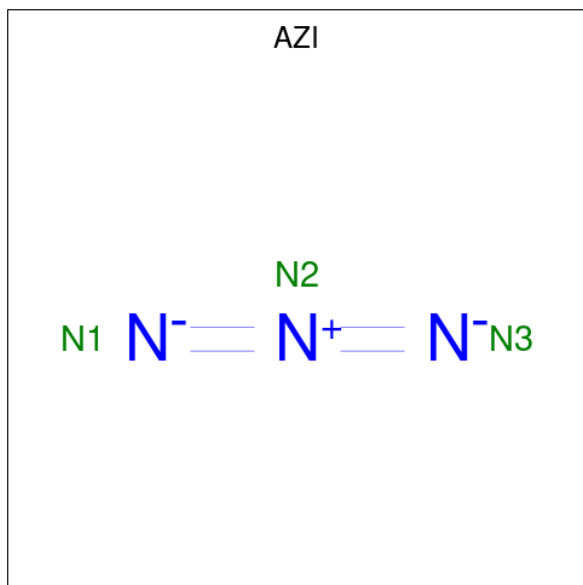
- Molecule 16 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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).

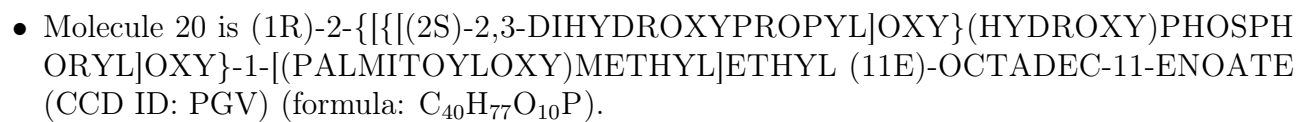
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
17	A	1	Total Na 1 1	0	0
17	N	1	Total Na 1 1	0	0

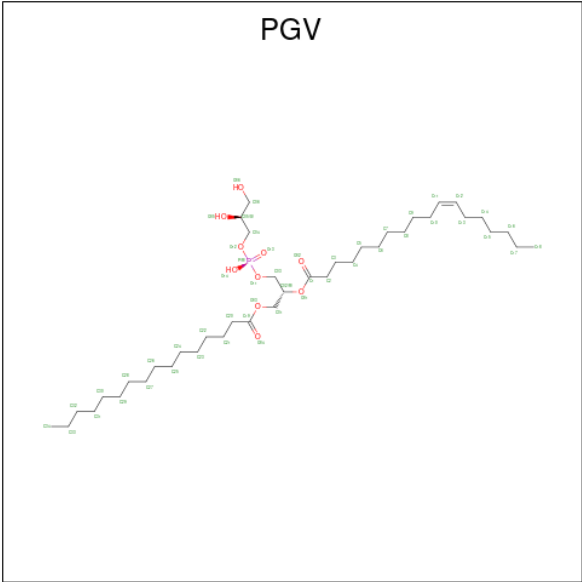
- Molecule 18 is AZIDE ION (CCD ID: AZI) (formula: N₃).



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

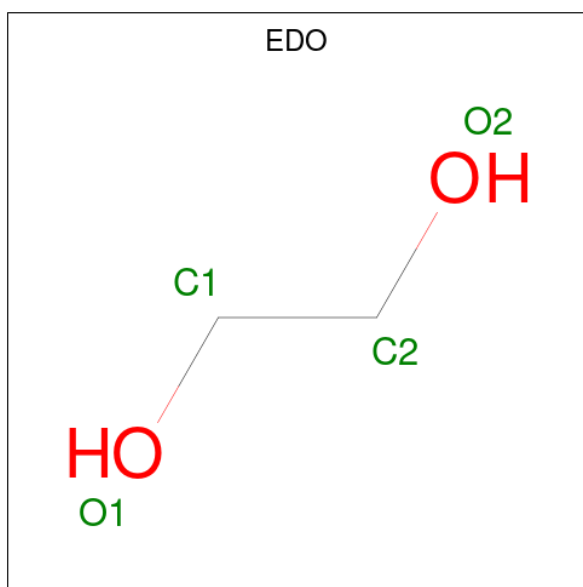
- Molecule 19 is TRISTEAROYLGLYCEROL (CCD ID: TGL) (formula: C₅₇H₁₁₀O₆).





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

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

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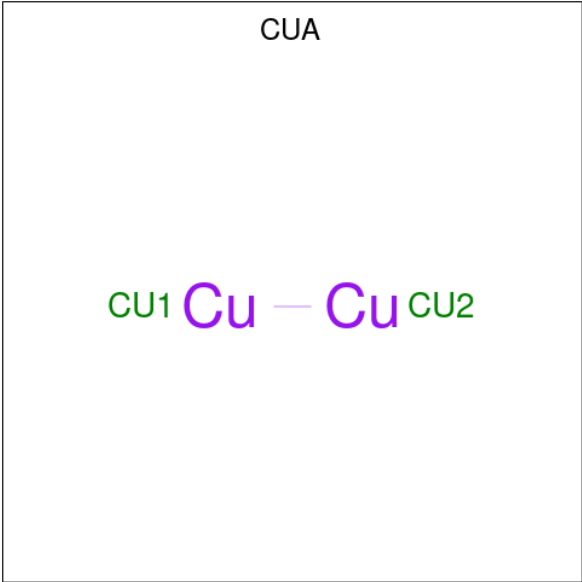
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
21	G	1	Total 4	C 2	O 2	0	0
21	H	1	Total 4	C 2	O 2	0	0
21	N	1	Total 4	C 2	O 2	0	0
21	N	1	Total 4	C 2	O 2	0	0
21	N	1	Total 4	C 2	O 2	0	0
21	N	1	Total 4	C 2	O 2	0	0
21	N	1	Total 4	C 2	O 2	0	0
21	N	1	Total 4	C 2	O 2	0	0
21	N	1	Total 4	C 2	O 2	0	0
21	N	1	Total 4	C 2	O 2	0	0
21	N	1	Total 4	C 2	O 2	0	0
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	O	1	Total 4	C 2	O 2	0	0
21	O	1	Total 4	C 2	O 2	0	0
21	O	1	Total 4	C 2	O 2	0	0
21	P	1	Total 4	C 2	O 2	0	0
21	P	1	Total 4	C 2	O 2	0	0
21	P	1	Total 4	C 2	O 2	0	0
21	P	1	Total 4	C 2	O 2	0	0
21	Q	1	Total 4	C 2	O 2	0	0

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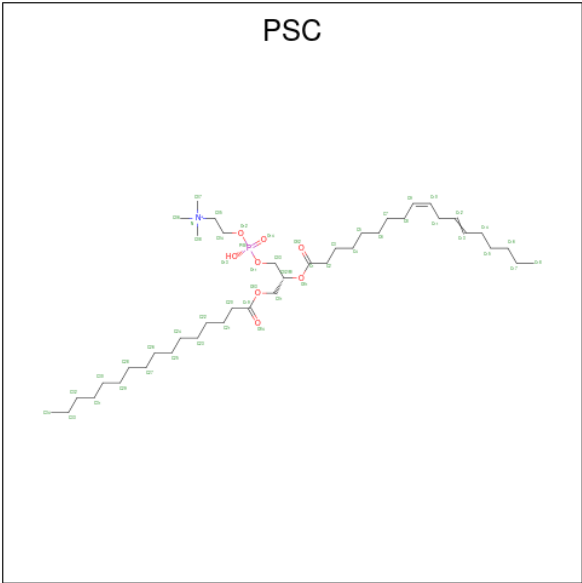
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
21	Q	1	Total	C	O	0	0
			4	2	2		
21	Q	1	Total	C	O	0	0
			4	2	2		
21	R	1	Total	C	O	0	0
			4	2	2		
21	R	1	Total	C	O	0	0
			4	2	2		
21	R	1	Total	C	O	0	0
			4	2	2		
21	R	1	Total	C	O	0	0
			4	2	2		
21	R	1	Total	C	O	0	0
			4	2	2		
21	S	1	Total	C	O	0	0
			4	2	2		
21	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	U	1	Total	C	O	0	0
			4	2	2		
21	V	1	Total	C	O	0	0
			4	2	2		

- Molecule 22 is DINUCLEAR COPPER ION (CCD ID: CUA) (formula: Cu₂).



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

- Molecule 23 is (7R,17E,20E)-4-HYDROXY-N,N,N-TRIMETHYL-9-OXO-7-[(PALMITO YLOXY)METHYL]-3,5,8-TRIOXA-4-PHOSPHAHEXACOSA-17,20-DIEN-1-AMINIUM 4-OXIDE (CCD ID: PSC) (formula: C₄₂H₈₁NO₈P).



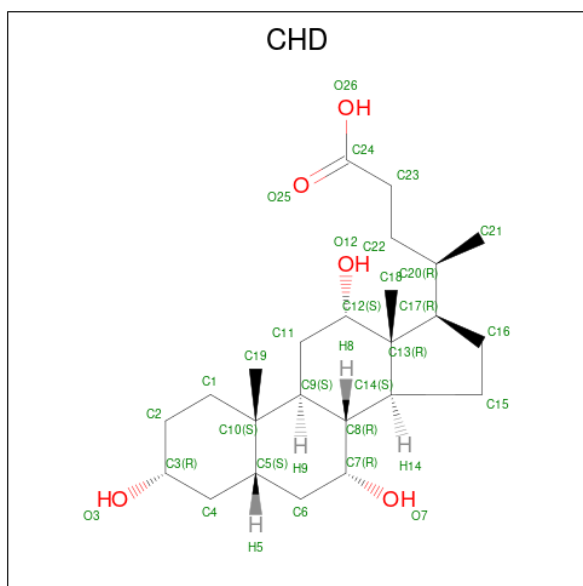
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
23	B	1	Total	C	N	O	P	0	0
			52	42	1	8	1		

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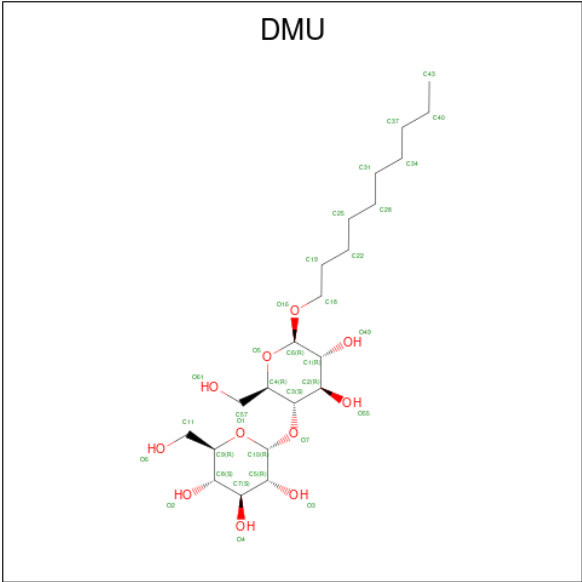
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
23	O	1	Total	C	N	O	P	0	0
			52	42	1	8	1		

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



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

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

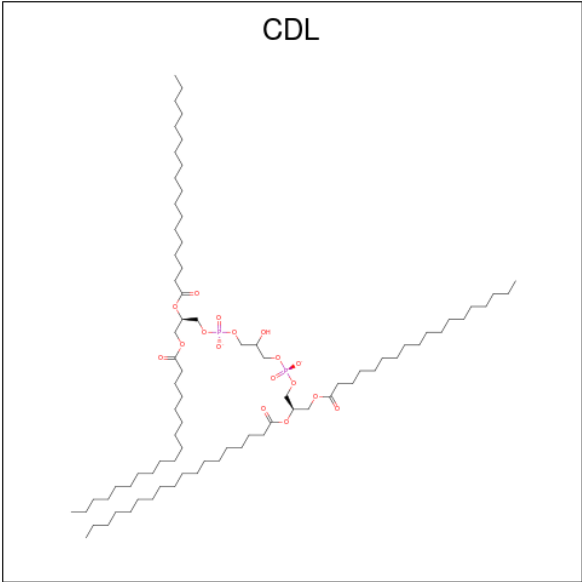


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
25	C	1	Total	C	O	0	0
			33	22	11		
25	C	1	Total	C	O	0	0
			33	22	11		
25	C	1	Total	C	O	0	0
			33	22	11		
25	M	1	Total	C	O	0	0
			33	22	11		
25	P	1	Total	C	O	0	0
			33	22	11		
25	P	1	Total	C	O	0	0
			33	22	11		
25	P	1	Total	C	O	0	0
			33	22	11		
25	Z	1	Total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

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

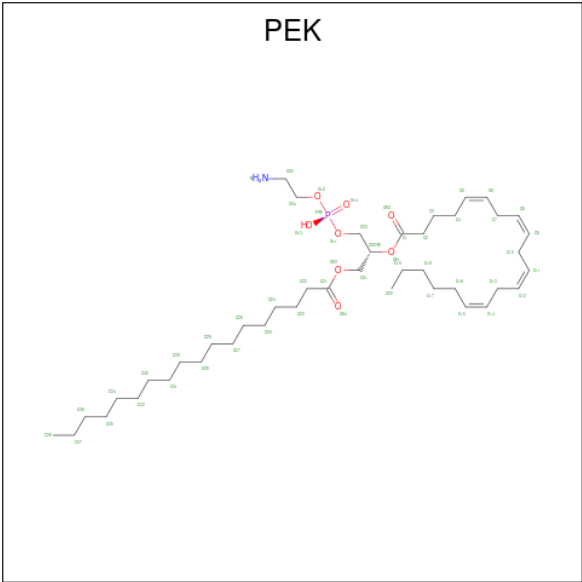
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
26	C	1	Total	X	0	0
			1	1		
26	P	1	Total	X	0	0
			1	1		

- Molecule 27 is CARDIOLIPIN (CCD ID: CDL) (formula: C₈₁H₁₅₆O₁₇P₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
27	C	1	Total	C	O	P	0	0
			100	81	17	2		
27	G	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-[(STEAROYLOXY)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	C	N	O	P	0	0
			53	43	1	8	1		
28	G	1	Total	C	N	O	P	0	0
			53	43	1	8	1		
28	G	1	Total	C	N	O	P	0	0
			53	43	1	8	1		
28	P	1	Total	C	N	O	P	0	0
			53	43	1	8	1		
28	P	1	Total	C	N	O	P	0	0
			53	43	1	8	1		
28	T	1	Total	C	N	O	P	0	0
			53	43	1	8	1		

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

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

- Molecule 30 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
30	A	219	Total	O	0	0
			219	219		
30	B	131	Total	O	0	1
			132	132		
30	C	96	Total	O	0	0
			96	96		
30	D	93	Total	O	0	0
			93	93		
30	E	85	Total	O	0	0
			85	85		
30	F	78	Total	O	0	0
			78	78		
30	G	37	Total	O	0	0
			37	37		
30	H	30	Total	O	0	0
			30	30		
30	I	20	Total	O	0	0
			20	20		
30	J	22	Total	O	0	0
			22	22		

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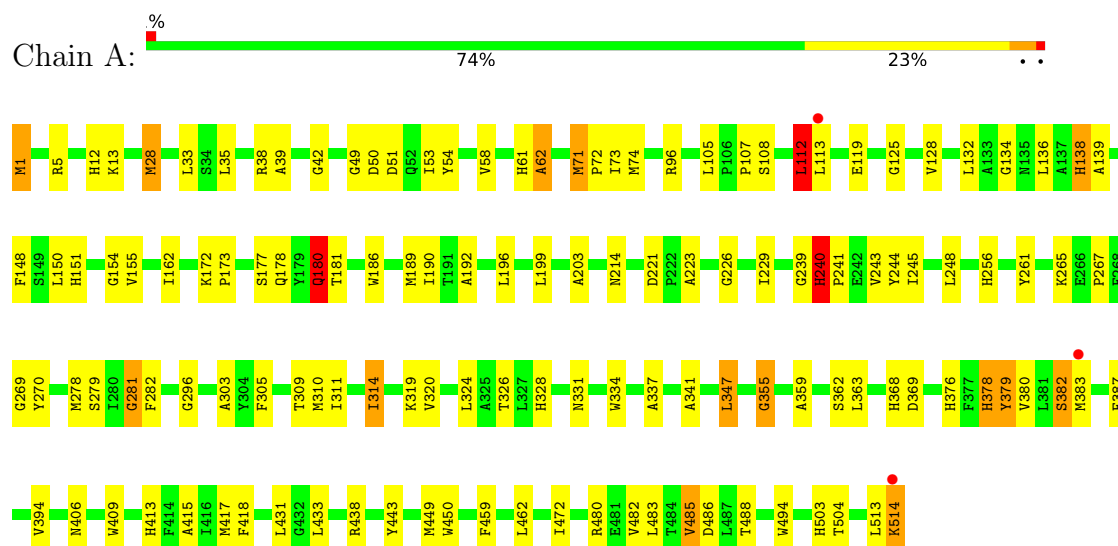
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
30	K	17	Total 17	O 17	0	0
30	L	21	Total 21	O 21	0	0
30	M	22	Total 22	O 22	0	0
30	N	217	Total 217	O 217	0	0
30	O	114	Total 115	O 115	0	1
30	P	95	Total 95	O 95	0	0
30	Q	57	Total 57	O 57	0	0
30	R	56	Total 56	O 56	0	0
30	S	65	Total 65	O 65	0	0
30	T	33	Total 33	O 33	0	0
30	U	38	Total 38	O 38	0	0
30	V	16	Total 16	O 16	0	0
30	W	13	Total 13	O 13	0	0
30	X	15	Total 15	O 15	0	0
30	Y	14	Total 14	O 14	0	0
30	Z	12	Total 12	O 12	0	0

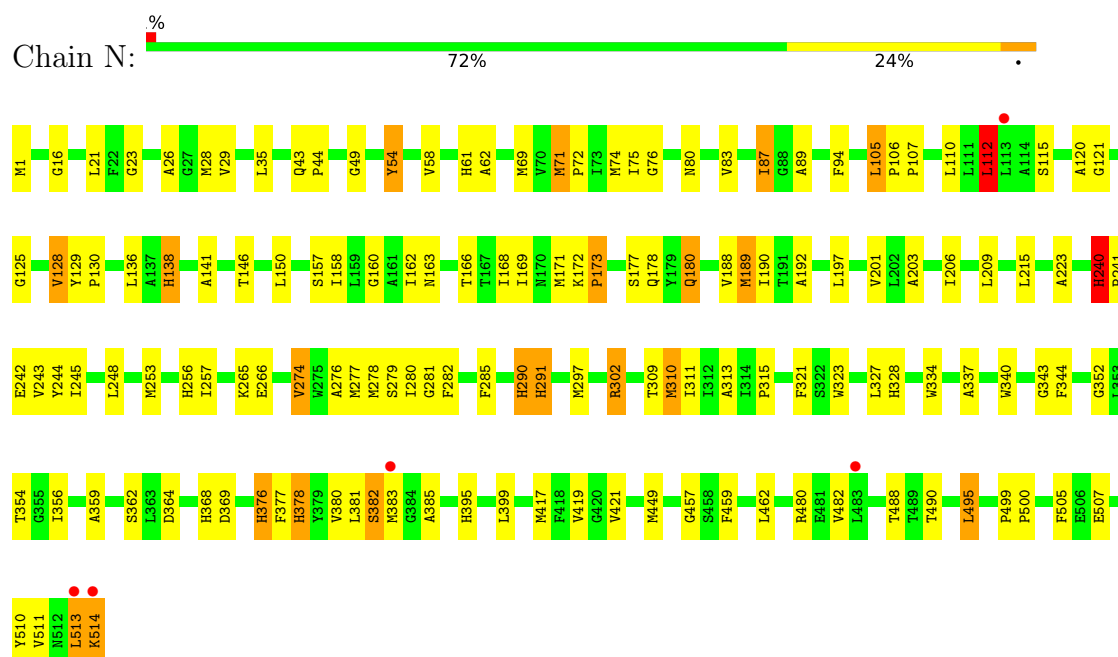
3 Residue-property plots

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.

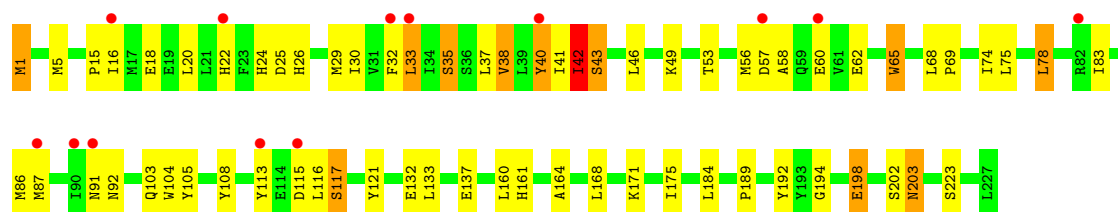
• Molecule 1: Cytochrome c oxidase subunit 1



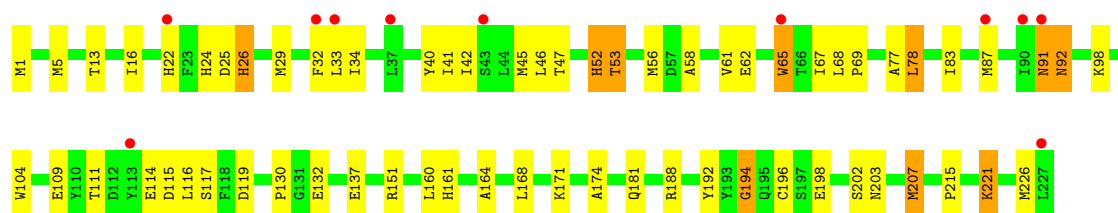
• Molecule 1: Cytochrome c oxidase subunit 1



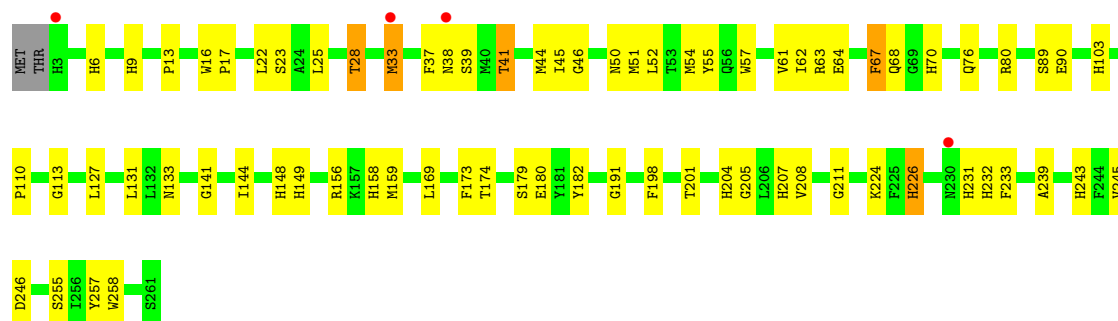
- Molecule 2: Cytochrome c oxidase subunit 2



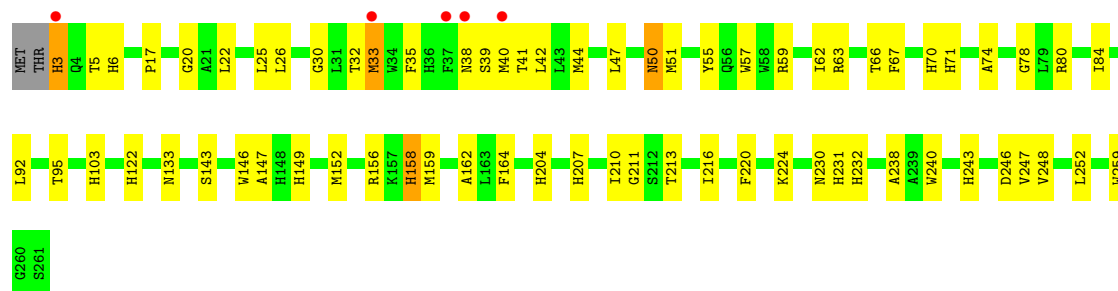
- Molecule 2: Cytochrome c oxidase subunit 2



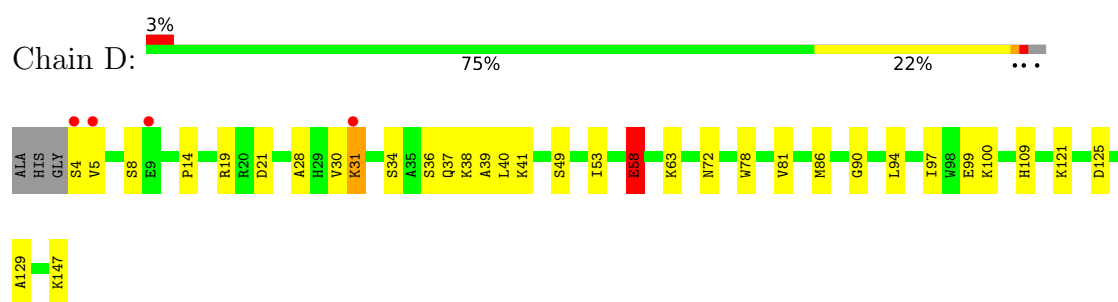
- Molecule 3: Cytochrome c oxidase subunit 3



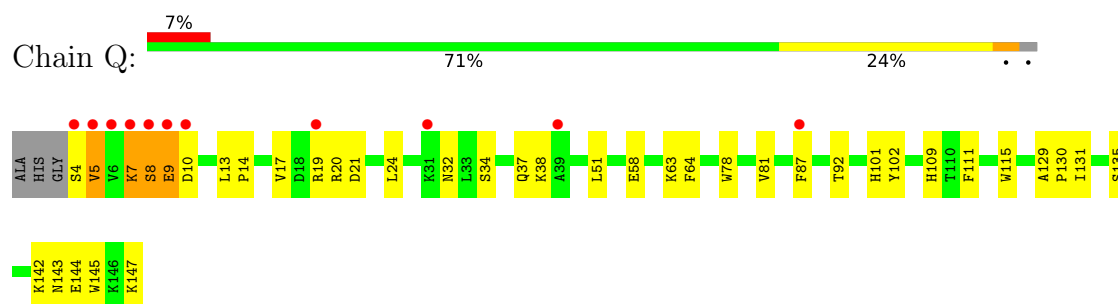
- Molecule 3: Cytochrome c oxidase subunit 3



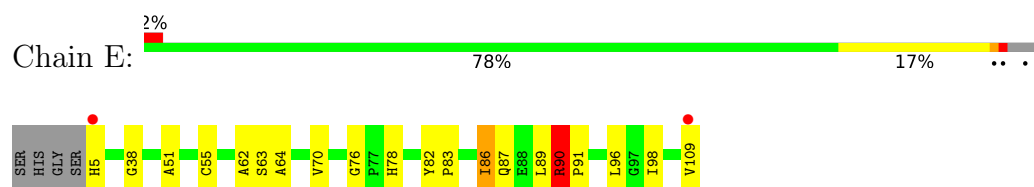
- Molecule 4: Cytochrome c oxidase subunit 4 isoform 1, mitochondrial



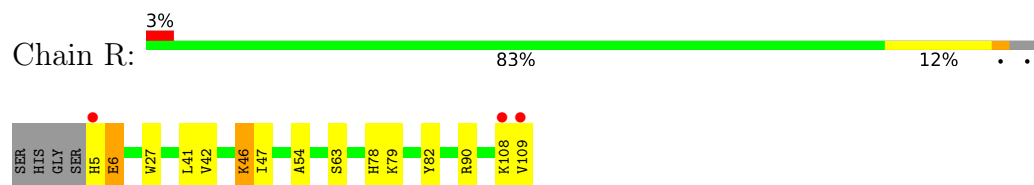
- Molecule 4: Cytochrome c oxidase subunit 4 isoform 1, mitochondrial



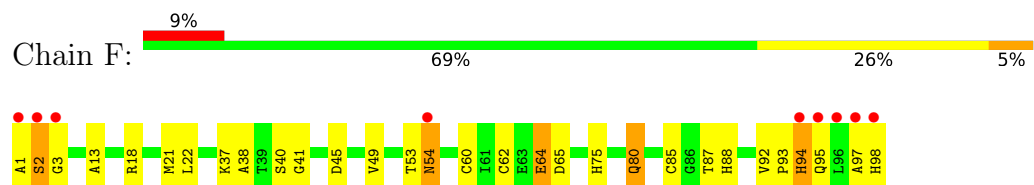
- Molecule 5: Cytochrome c oxidase subunit 5A, mitochondrial



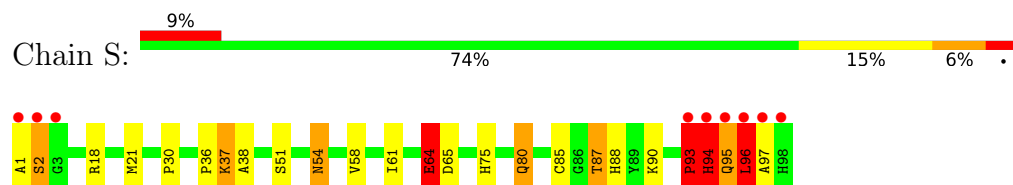
- Molecule 5: Cytochrome c oxidase subunit 5A, mitochondrial



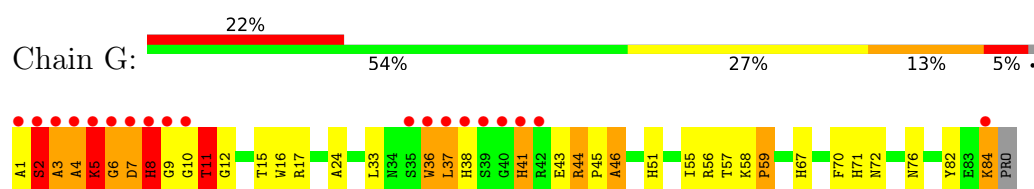
- Molecule 6: Cytochrome c oxidase subunit 5B, mitochondrial



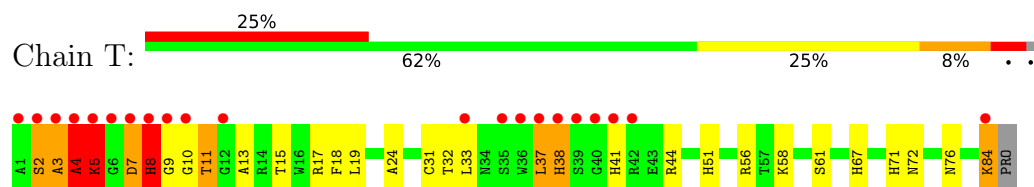
- Molecule 6: Cytochrome c oxidase subunit 5B, mitochondrial



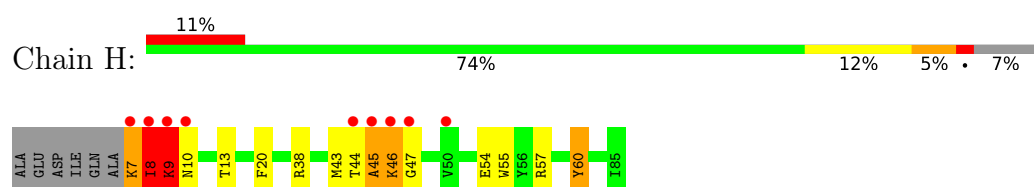
- Molecule 7: Cytochrome c oxidase subunit 6A2, mitochondrial



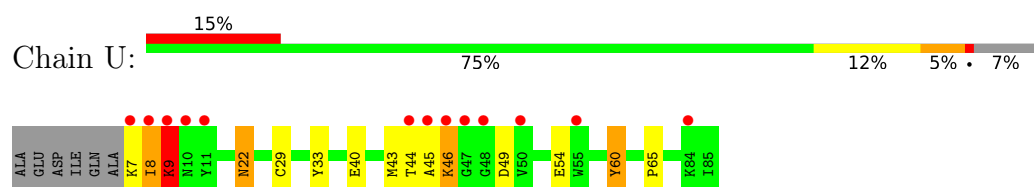
- Molecule 7: Cytochrome c oxidase subunit 6A2, mitochondrial



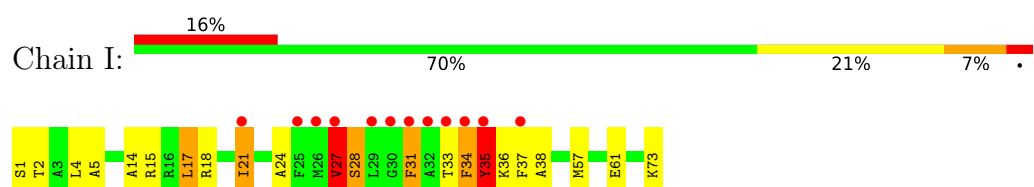
- Molecule 8: Cytochrome c oxidase subunit 6B1



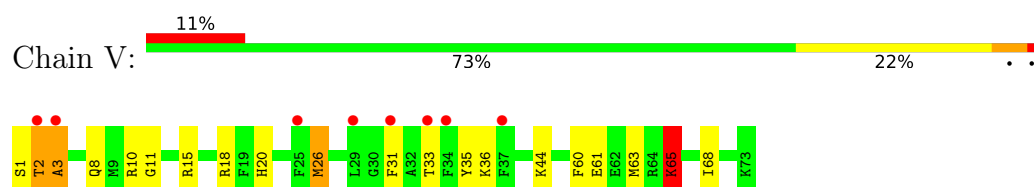
- Molecule 8: Cytochrome c oxidase subunit 6B1



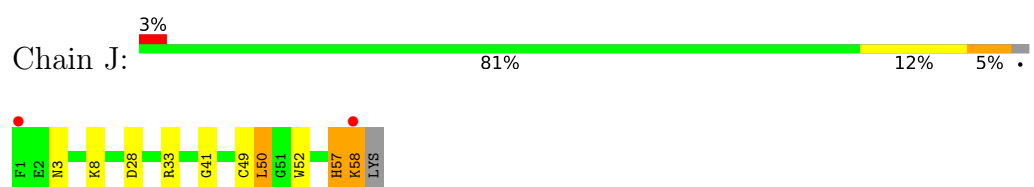
- Molecule 9: Cytochrome c oxidase subunit 6C



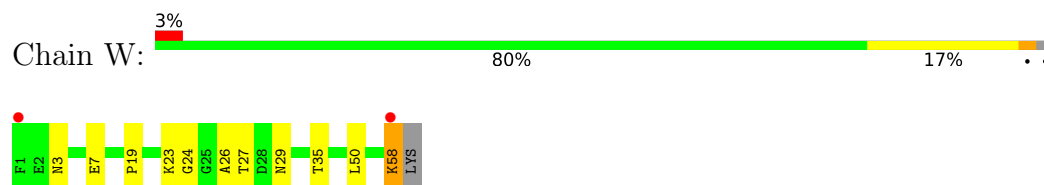
- Molecule 9: Cytochrome c oxidase subunit 6C



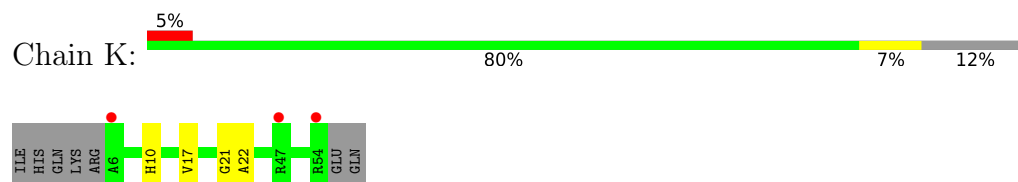
- Molecule 10: Cytochrome c oxidase subunit 7A1, mitochondrial



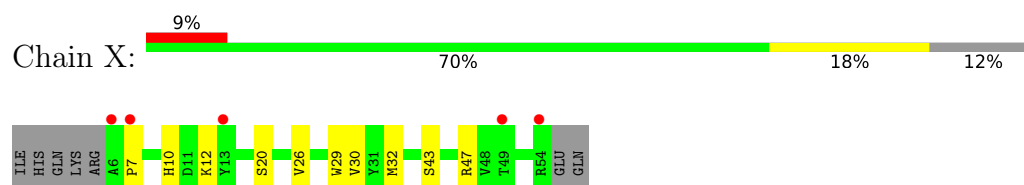
- Molecule 10: Cytochrome c oxidase subunit 7A1, mitochondrial



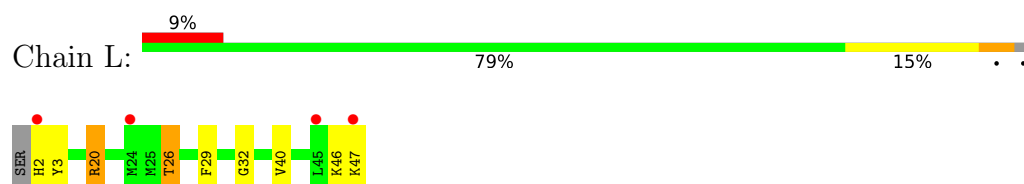
- Molecule 11: Cytochrome c oxidase subunit 7B, mitochondrial



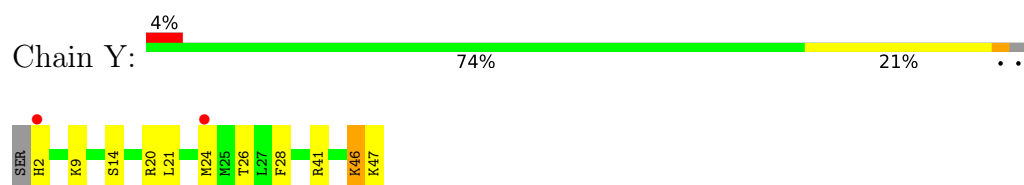
- Molecule 11: Cytochrome c oxidase subunit 7B, mitochondrial



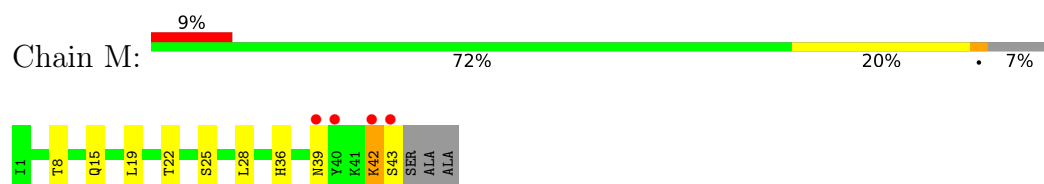
- Molecule 12: Cytochrome c oxidase subunit 7C, mitochondrial



- 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	183.13Å 205.89Å 177.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 1.85 40.00 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.3 (40.00-1.85) 99.3 (40.00-1.85)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.13 (at 1.86Å)	Xtriage
Refinement program	REFMAC 5.8.0048	Depositor
R, R_{free}	0.163 , 0.189 0.165 , 0.191	Depositor DCC
R_{free} test set	28242 reflections (2.23%)	wwPDB-VP
Wilson B-factor (Å ²)	27.6	Xtriage
Anisotropy	0.677	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 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	33411	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% 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: EDO, PEK, UNX, CHD, PSC, CUA, NA, PGV, CU, DMU, CDL, TPO, SAC, MG, AZI, ZN, FME, TGL, HEA

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	1.98	76/4294 (1.8%)	1.53	37/5861 (0.6%)
1	N	1.93	79/4289 (1.8%)	1.46	32/5853 (0.5%)
2	B	1.82	24/1896 (1.3%)	1.44	10/2582 (0.4%)
2	O	1.57	19/1908 (1.0%)	1.33	8/2597 (0.3%)
3	C	1.89	38/2258 (1.7%)	1.35	12/3084 (0.4%)
3	P	1.90	36/2272 (1.6%)	1.34	4/3102 (0.1%)
4	D	1.74	14/1238 (1.1%)	1.41	4/1669 (0.2%)
4	Q	1.38	4/1248 (0.3%)	1.18	4/1684 (0.2%)
5	E	1.93	16/871 (1.8%)	1.43	7/1182 (0.6%)
5	R	1.60	4/871 (0.5%)	1.26	4/1182 (0.3%)
6	F	1.76	7/788 (0.9%)	1.49	7/1069 (0.7%)
6	S	1.68	6/780 (0.8%)	1.44	8/1058 (0.8%)
7	G	1.93	18/702 (2.6%)	1.40	5/953 (0.5%)
7	T	1.90	17/690 (2.5%)	1.47	9/937 (1.0%)
8	H	1.57	3/682 (0.4%)	1.32	2/921 (0.2%)
8	U	1.36	1/682 (0.1%)	1.21	2/921 (0.2%)
9	I	1.59	5/605 (0.8%)	1.43	5/802 (0.6%)
9	V	1.32	2/605 (0.3%)	1.36	2/802 (0.2%)
10	J	1.68	3/471 (0.6%)	1.38	1/636 (0.2%)
10	W	1.58	4/471 (0.8%)	1.26	2/636 (0.3%)
11	K	1.69	1/398 (0.3%)	1.33	2/546 (0.4%)
11	X	1.34	2/405 (0.5%)	1.13	0/556
12	L	1.74	2/393 (0.5%)	1.46	3/526 (0.6%)
12	Y	1.47	1/401 (0.2%)	1.25	0/536
13	M	1.73	4/345 (1.2%)	1.27	0/470
13	Z	1.52	4/345 (1.2%)	1.21	0/470
All	All	1.78	390/29908 (1.3%)	1.40	170/40635 (0.4%)

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	2
1	N	0	1
4	Q	0	1
5	R	0	1
6	F	0	1
6	S	0	1
7	G	0	1
8	H	0	2
8	U	0	1
9	V	0	1
12	Y	0	1
All	All	0	13

The worst 5 of 390 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	O	65	TRP	NE1-CE2	-12.61	1.23	1.37
5	E	78	HIS	N-CA	10.08	1.59	1.46
1	A	310	MET	SD-CE	-9.30	1.56	1.79
1	N	240	HIS	N-CA	9.28	1.53	1.46
1	A	49	GLY	C-O	9.19	1.36	1.23

The worst 5 of 170 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	O	65	TRP	CA-CB-CG	12.87	138.06	113.60
1	A	278	MET	CG-SD-CE	-12.25	73.96	100.90
1	A	71	MET	CG-SD-CE	-11.81	74.93	100.90
1	N	71	MET	CG-SD-CE	-11.36	75.92	100.90
1	A	112	LEU	CD1-CG-CD2	-10.26	88.22	110.80

There are no chirality outliers.

5 of 13 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	240	HIS	Sidechain
1	A	296	GLY	Mainchain
6	F	93	PRO	Peptide
7	G	11	TPO	Peptide
8	H	8	ILE	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	4165	0	4137	83	0
1	N	4160	0	4132	92	0
2	B	1859	0	1869	46	0
2	O	1870	0	1868	41	0
3	C	2171	0	2085	48	0
3	P	2185	0	2097	55	0
4	D	1204	0	1195	24	0
4	Q	1213	0	1199	34	0
5	E	852	0	845	4	0
5	R	852	0	845	4	0
6	F	771	0	747	25	0
6	S	763	0	742	17	0
7	G	686	0	652	53	0
7	T	675	0	643	37	0
8	H	662	0	623	12	0
8	U	662	0	623	11	0
9	I	601	0	613	18	0
9	V	601	0	613	18	0
10	J	460	0	459	7	0
10	W	460	0	459	7	0
11	K	384	0	366	1	0
11	X	391	0	374	13	0
12	L	380	0	380	11	0
12	Y	388	0	388	12	0
13	M	335	0	352	5	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	0	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	6	0	0	6	0
18	N	6	0	0	6	0
19	A	63	0	110	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
19	D	63	0	110	13	0
19	L	63	0	110	17	0
19	N	63	0	110	4	0
19	Q	63	0	110	12	0
19	Y	63	0	110	16	0
20	A	51	0	76	4	0
20	C	102	0	152	7	0
20	M	51	0	76	3	0
20	N	102	0	152	14	0
20	P	102	0	152	12	0
21	A	44	0	66	20	0
21	B	4	0	6	0	0
21	C	36	0	54	3	0
21	D	20	0	30	12	0
21	E	20	0	30	0	0
21	F	12	0	18	1	0
21	G	8	0	12	1	0
21	H	4	0	6	3	0
21	N	44	0	66	15	0
21	O	12	0	18	0	0
21	P	16	0	24	1	0
21	Q	12	0	18	6	0
21	R	20	0	30	3	0
21	S	12	0	18	3	0
21	T	4	0	6	0	0
21	U	4	0	6	0	0
21	V	4	0	6	0	0
22	B	2	0	0	0	0
22	O	2	0	0	0	0
23	B	52	0	80	13	0
23	O	52	0	80	16	0
24	C	58	0	78	5	0
24	G	29	0	39	1	0
24	P	58	0	77	6	0
24	T	29	0	39	0	0
25	C	99	0	126	19	0
25	M	33	0	42	0	0
25	P	99	0	126	14	0
25	Z	33	0	42	0	0
26	C	1	0	0	0	0
26	P	1	0	0	0	0
27	C	100	0	156	25	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
27	G	100	0	156	24	0
27	P	100	0	156	22	0
27	T	100	0	156	25	0
28	C	53	0	77	10	0
28	G	106	0	154	23	0
28	P	106	0	154	5	0
28	T	53	0	77	10	0
29	F	1	0	0	0	0
29	S	1	0	0	0	0
30	A	219	0	0	13	0
30	B	132	0	0	15	1
30	C	96	0	0	5	0
30	D	93	0	0	8	1
30	E	85	0	0	1	0
30	F	78	0	0	3	0
30	G	37	0	0	4	0
30	H	30	0	0	0	0
30	I	20	0	0	2	0
30	J	22	0	0	0	0
30	K	17	0	0	0	0
30	L	21	0	0	0	0
30	M	22	0	0	3	0
30	N	217	0	0	13	0
30	O	115	0	0	4	0
30	P	95	0	0	3	0
30	Q	57	0	0	10	0
30	R	56	0	0	4	0
30	S	65	0	0	1	0
30	T	33	0	0	0	0
30	U	38	0	0	2	0
30	V	16	0	0	3	0
30	W	13	0	0	1	0
30	X	15	0	0	2	0
30	Y	14	0	0	0	0
30	Z	12	0	0	2	0
All	All	33411	0	32479	803	1

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 803 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)
21:A:614:EDO:C1	21:A:614:EDO:C2	1.83	1.54
20:N:608:PGV:C2	20:N:608:PGV:H011	1.28	1.46
19:L:101:TGL:HC62	19:L:101:TGL:CC2	1.34	1.38
21:A:614:EDO:C2	21:A:614:EDO:O1	1.74	1.29
1:N:297[B]:MET:HB2	30:N:782:HOH:O	1.36	1.25

All (1) 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:468:HOH:O	30:D:353:HOH:O[2_584]	1.99	0.21

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	530/514 (103%)	514 (97%)	16 (3%)	0	100	100
1	N	530/514 (103%)	516 (97%)	14 (3%)	0	100	100
2	B	229/227 (101%)	223 (97%)	6 (3%)	0	100	100
2	O	230/227 (101%)	224 (97%)	5 (2%)	1 (0%)	30	17
3	C	264/261 (101%)	260 (98%)	4 (2%)	0	100	100
3	P	266/261 (102%)	261 (98%)	5 (2%)	0	100	100
4	D	143/147 (97%)	139 (97%)	4 (3%)	0	100	100
4	Q	144/147 (98%)	138 (96%)	4 (3%)	2 (1%)	9	2
5	E	103/109 (94%)	103 (100%)	0	0	100	100
5	R	103/109 (94%)	102 (99%)	0	1 (1%)	12	4
6	F	99/98 (101%)	93 (94%)	2 (2%)	4 (4%)	2	0
6	S	98/98 (100%)	89 (91%)	4 (4%)	5 (5%)	1	0
7	G	82/85 (96%)	69 (84%)	6 (7%)	7 (8%)	0	0

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
7	T	81/85 (95%)	70 (86%)	7 (9%)	4 (5%)	1	0
8	H	77/85 (91%)	71 (92%)	3 (4%)	3 (4%)	2	0
8	U	77/85 (91%)	68 (88%)	6 (8%)	3 (4%)	2	0
9	I	71/73 (97%)	68 (96%)	1 (1%)	2 (3%)	4	0
9	V	71/73 (97%)	68 (96%)	3 (4%)	0	100	100
10	J	56/59 (95%)	56 (100%)	0	0	100	100
10	W	56/59 (95%)	56 (100%)	0	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%)	42 (96%)	2 (4%)	0	100	100
12	Y	45/47 (96%)	43 (96%)	2 (4%)	0	100	100
13	M	41/46 (89%)	39 (95%)	1 (2%)	1 (2%)	4	1
13	Z	41/46 (89%)	41 (100%)	0	0	100	100
All	All	3576/3614 (99%)	3444 (96%)	99 (3%)	33 (1%)	14	4

5 of 33 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
6	F	2	SER
6	F	94	HIS
6	F	95	GLN
7	G	2	SER
7	G	4	ALA

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	443/426 (104%)	433 (98%)	10 (2%)	44	30
1	N	443/426 (104%)	433 (98%)	10 (2%)	44	30
2	B	214/210 (102%)	201 (94%)	13 (6%)	17	5

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	O	215/210 (102%)	209 (97%)	6 (3%)	38	23
3	C	231/226 (102%)	227 (98%)	4 (2%)	53	42
3	P	233/226 (103%)	227 (97%)	6 (3%)	40	25
4	D	129/129 (100%)	124 (96%)	5 (4%)	28	13
4	Q	130/129 (101%)	124 (95%)	6 (5%)	24	9
5	E	92/95 (97%)	88 (96%)	4 (4%)	26	11
5	R	92/95 (97%)	88 (96%)	4 (4%)	26	11
6	F	84/81 (104%)	79 (94%)	5 (6%)	17	5
6	S	83/81 (102%)	71 (86%)	12 (14%)	3	0
7	G	68/68 (100%)	61 (90%)	7 (10%)	7	1
7	T	67/68 (98%)	63 (94%)	4 (6%)	17	5
8	H	71/75 (95%)	66 (93%)	5 (7%)	14	3
8	U	71/75 (95%)	67 (94%)	4 (6%)	19	6
9	I	57/57 (100%)	54 (95%)	3 (5%)	20	7
9	V	57/57 (100%)	54 (95%)	3 (5%)	20	7
10	J	49/50 (98%)	48 (98%)	1 (2%)	48	36
10	W	49/50 (98%)	46 (94%)	3 (6%)	17	5
11	K	39/46 (85%)	39 (100%)	0	100	100
11	X	40/46 (87%)	40 (100%)	0	100	100
12	L	39/40 (98%)	38 (97%)	1 (3%)	40	25
12	Y	40/40 (100%)	36 (90%)	4 (10%)	7	1
13	M	37/38 (97%)	36 (97%)	1 (3%)	39	24
13	Z	37/38 (97%)	36 (97%)	1 (3%)	39	24
All	All	3110/3082 (101%)	2988 (96%)	122 (4%)	29	13

5 of 122 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
12	L	47	LYS
8	U	65	PRO
2	O	91	ASN
8	U	60	TYR
12	Y	9	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 47 such sidechains are listed below:

Mol	Chain	Res	Type
4	Q	76	ASN
7	T	8	HIS
4	Q	101	HIS
6	S	54	ASN
8	U	22	ASN

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	1.15	0	8,9,11	1.81	4 (50%)
7	TPO	G	11	7	8,10,11	2.22	3 (37%)	10,14,16	1.73	2 (20%)
2	FME	O	1	2	8,9,10	1.39	1 (12%)	8,9,11	1.55	2 (25%)
9	SAC	I	1	9	7,8,9	1.30	1 (14%)	7,9,11	2.12	2 (28%)
1	FME	N	1	1	8,9,10	1.10	1 (12%)	8,9,11	1.27	0
2	FME	B	1	2	8,9,10	1.78	3 (37%)	8,9,11	1.67	2 (25%)
7	TPO	T	11	7	8,10,11	1.77	1 (12%)	10,14,16	1.02	1 (10%)
9	SAC	V	1	9	7,8,9	1.61	1 (14%)	7,9,11	1.46	2 (28%)

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	-	3/7/9/11	-
7	TPO	G	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	-	2/7/9/11	-
2	FME	B	1	2	-	0/7/9/11	-
7	TPO	T	11	7	-	3/9/11/13	-
9	SAC	V	1	9	-	5/7/8/10	-

The worst 5 of 11 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	V	1	SAC	CA-N	3.99	1.52	1.46
7	T	11	TPO	P-O1P	3.67	1.61	1.50
7	G	11	TPO	P-O1P	3.33	1.60	1.50
7	G	11	TPO	P-OG1	3.32	1.65	1.59
9	I	1	SAC	CA-N	2.94	1.50	1.46

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	I	1	SAC	OG-CB-CA	-3.78	100.70	111.13
9	I	1	SAC	CB-CA-C	-3.15	104.43	111.35
2	B	1	FME	C-CA-N	-3.09	103.53	109.50
2	B	1	FME	CG-CB-CA	-3.01	103.67	112.87
7	G	11	TPO	O2P-P-OG1	2.99	117.50	105.85

There are no chirality outliers.

5 of 23 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-CG2
7	G	11	TPO	N-CA-CB-OG1
7	G	11	TPO	CA-CB-OG1-P
7	G	11	TPO	CG2-CB-OG1-P

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	1	FME	1	0
7	G	11	TPO	3	0
2	B	1	FME	2	0
7	T	11	TPO	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 131 ligands modelled in this entry, 8 are monoatomic and 2 are unknown - leaving 121 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	E	201	-	3,3,3	0.59	0	2,2,2	0.45	0
18	AZI	A	606	15	2,2,2	4.66	2 (100%)	0,1,1	-	-
20	PGV	P	302	-	50,50,50	1.18	2 (4%)	53,56,56	1.43	8 (15%)
21	EDO	F	103	-	3,3,3	0.78	0	2,2,2	0.69	0
23	PSC	O	302	-	51,51,51	1.29	3 (5%)	57,59,59	1.30	5 (8%)
25	DMU	C	309	-	34,34,34	0.89	2 (5%)	45,45,45	2.65	11 (24%)
22	CUA	B	301	2	0,1,1	-	-	-	-	-
21	EDO	G	105	-	3,3,3	0.56	0	2,2,2	0.57	0
14	HEA	A	602[A]	1,18	67,67,67	1.96	27 (40%)	81,103,103	2.13	25 (30%)
21	EDO	C	311	-	3,3,3	1.01	0	2,2,2	0.36	0
21	EDO	P	312	-	3,3,3	0.74	0	2,2,2	1.78	1 (50%)
21	EDO	E	203	-	3,3,3	0.34	0	2,2,2	1.52	0
21	EDO	A	619	-	3,3,3	0.56	0	2,2,2	0.93	0
21	EDO	A	610	-	3,3,3	0.31	0	2,2,2	1.39	0
18	AZI	N	606	15	2,2,2	4.26	2 (100%)	0,1,1	-	-
21	EDO	D	202	-	3,3,3	0.42	0	2,2,2	0.49	0
14	HEA	N	601	1	67,67,67	2.02	23 (34%)	81,103,103	2.51	40 (49%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
21	EDO	A	615	-	3,3,3	1.25	0	2,2,2	0.29	0
21	EDO	A	617	-	3,3,3	0.26	0	2,2,2	1.68	1 (50%)
21	EDO	C	318	-	3,3,3	0.45	0	2,2,2	0.39	0
19	TGL	N	610	-	62,62,62	1.23	4 (6%)	65,65,65	1.74	11 (16%)
21	EDO	O	305	-	3,3,3	0.42	0	2,2,2	0.40	0
24	CHD	C	306	-	32,32,32	1.27	4 (12%)	51,51,51	3.05	18 (35%)
28	PEK	T	102	-	52,52,52	1.28	2 (3%)	55,57,57	1.51	7 (12%)
21	EDO	R	203	-	3,3,3	0.82	0	2,2,2	0.51	0
21	EDO	O	304	-	3,3,3	0.35	0	2,2,2	1.00	0
20	PGV	C	308	-	50,50,50	1.26	2 (4%)	53,56,56	1.37	5 (9%)
21	EDO	N	616	-	3,3,3	0.52	0	2,2,2	1.25	0
18	AZI	N	607	14	2,2,2	4.67	2 (100%)	0,1,1	-	-
25	DMU	C	310	-	34,34,34	0.98	1 (2%)	45,45,45	2.24	12 (26%)
24	CHD	G	103	-	32,32,32	1.61	8 (25%)	51,51,51	1.76	17 (33%)
21	EDO	N	620	-	3,3,3	0.45	0	2,2,2	1.42	0
20	PGV	M	101	-	50,50,50	1.42	4 (8%)	53,56,56	1.62	9 (16%)
25	DMU	P	311	-	34,34,34	0.86	1 (2%)	45,45,45	1.61	8 (17%)
18	AZI	A	607	14	2,2,2	5.49	2 (100%)	0,1,1	-	-
27	CDL	C	305	-	99,99,99	1.64	18 (18%)	105,111,111	1.59	18 (17%)
21	EDO	R	201	-	3,3,3	0.36	0	2,2,2	0.93	0
21	EDO	C	317	-	3,3,3	0.70	0	2,2,2	0.48	0
21	EDO	B	303	-	3,3,3	0.74	0	2,2,2	0.37	0
19	TGL	A	608	-	62,62,62	1.34	5 (8%)	65,65,65	2.28	11 (16%)
21	EDO	S	103	-	3,3,3	0.78	0	2,2,2	0.43	0
19	TGL	D	201	-	62,62,62	1.86	6 (9%)	65,65,65	2.32	12 (18%)
21	EDO	C	313	-	3,3,3	0.54	0	2,2,2	0.57	0
21	EDO	Q	204	-	3,3,3	0.69	0	2,2,2	0.17	0
21	EDO	Q	202	-	3,3,3	0.27	0	2,2,2	1.13	0
22	CUA	O	301	2	0,1,1	-	-	-	-	-
21	EDO	N	612	-	3,3,3	1.20	0	2,2,2	1.70	0
25	DMU	P	310	-	34,34,34	0.79	1 (2%)	45,45,45	2.40	12 (26%)
20	PGV	C	304	-	50,50,50	0.98	3 (6%)	53,56,56	1.55	10 (18%)
28	PEK	G	104	-	52,52,52	1.33	4 (7%)	55,57,57	1.54	6 (10%)
27	CDL	T	103	-	99,99,99	1.52	12 (12%)	105,111,111	1.49	19 (18%)
24	CHD	T	101	-	32,32,32	2.21	14 (43%)	51,51,51	1.96	18 (35%)
25	DMU	P	308	-	34,34,34	1.12	1 (2%)	45,45,45	1.56	4 (8%)
21	EDO	F	102	-	3,3,3	1.11	0	2,2,2	0.29	0
21	EDO	D	203	-	3,3,3	0.28	0	2,2,2	0.71	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
25	DMU	M	102	-	34,34,34	0.60	0	45,45,45	1.43	5 (11%)
24	CHD	P	301	-	32,32,32	1.69	8 (25%)	51,51,51	2.07	18 (35%)
21	EDO	A	613	-	3,3,3	0.30	0	2,2,2	1.36	0
19	TGL	Q	201	-	62,62,62	1.55	4 (6%)	65,65,65	1.46	9 (13%)
14	HEA	A	601	1	67,67,67	1.92	24 (35%)	81,103,103	2.72	42 (51%)
21	EDO	C	312	-	3,3,3	0.33	0	2,2,2	0.63	0
21	EDO	A	620	-	3,3,3	0.30	0	2,2,2	1.05	0
21	EDO	P	314	-	3,3,3	0.50	0	2,2,2	0.67	0
21	EDO	O	303	-	3,3,3	0.75	0	2,2,2	0.94	0
27	CDL	P	306	-	99,99,99	1.73	19 (19%)	105,111,111	1.57	17 (16%)
24	CHD	P	307	-	32,32,32	1.43	4 (12%)	51,51,51	2.95	25 (49%)
20	PGV	A	609	-	50,50,50	1.13	2 (4%)	53,56,56	1.19	5 (9%)
21	EDO	N	619	-	3,3,3	0.50	0	2,2,2	0.22	0
19	TGL	L	101	-	62,62,62	1.19	3 (4%)	65,65,65	1.77	15 (23%)
21	EDO	D	204	-	3,3,3	0.72	0	2,2,2	0.73	0
14	HEA	N	602[B]	1,18	67,67,67	2.00	21 (31%)	81,103,103	1.78	20 (24%)
21	EDO	N	618	-	3,3,3	0.72	0	2,2,2	0.75	0
25	DMU	C	302	-	34,34,34	1.18	2 (5%)	45,45,45	1.83	7 (15%)
21	EDO	N	613	-	3,3,3	0.65	0	2,2,2	0.51	0
21	EDO	A	616	-	3,3,3	0.57	0	2,2,2	1.10	0
21	EDO	U	101	-	3,3,3	0.43	0	2,2,2	0.53	0
20	PGV	P	305	-	50,50,50	1.07	4 (8%)	53,56,56	1.40	8 (15%)
21	EDO	V	101	-	3,3,3	0.40	0	2,2,2	0.47	0
21	EDO	N	617	-	3,3,3	0.97	0	2,2,2	0.21	0
21	EDO	E	204	-	3,3,3	0.63	0	2,2,2	0.35	0
20	PGV	N	609	-	50,50,50	1.10	5 (10%)	53,56,56	1.32	4 (7%)
21	EDO	C	315	-	3,3,3	1.04	0	2,2,2	1.05	0
21	EDO	S	104	-	3,3,3	0.70	0	2,2,2	1.55	1 (50%)
21	EDO	T	104	-	3,3,3	1.06	0	2,2,2	0.56	0
28	PEK	G	101	-	52,52,52	1.07	5 (9%)	55,57,57	1.30	10 (18%)
21	EDO	H	101	-	3,3,3	0.28	0	2,2,2	2.19	2 (100%)
23	PSC	B	302	-	51,51,51	1.45	4 (7%)	57,59,59	1.52	5 (8%)
21	EDO	Q	203	-	3,3,3	1.19	0	2,2,2	1.02	0
21	EDO	N	611	-	3,3,3	0.46	0	2,2,2	1.41	0
21	EDO	S	102	-	3,3,3	0.95	0	2,2,2	0.31	0
21	EDO	F	104	-	3,3,3	0.34	0	2,2,2	0.37	0
14	HEA	N	602[A]	1,18	67,67,67	1.93	20 (29%)	81,103,103	1.96	27 (33%)
21	EDO	D	206	-	3,3,3	0.29	0	2,2,2	1.19	0
24	CHD	C	301	-	32,32,32	1.84	10 (31%)	51,51,51	2.41	16 (31%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
20	PGV	N	608	-	50,50,50	1.21	2 (4%)	53,56,56	1.36	6 (11%)
21	EDO	N	614	-	3,3,3	0.48	0	2,2,2	0.94	0
19	TGL	Y	101	-	62,62,62	1.48	4 (6%)	65,65,65	1.57	11 (16%)
21	EDO	A	612	-	3,3,3	0.96	0	2,2,2	1.31	0
21	EDO	N	615	-	3,3,3	1.73	1 (33%)	2,2,2	0.29	0
28	PEK	C	307	-	52,52,52	1.22	3 (5%)	55,57,57	1.35	7 (12%)
21	EDO	R	205	-	3,3,3	1.09	0	2,2,2	0.38	0
28	PEK	P	309	-	52,52,52	1.25	2 (3%)	55,57,57	1.44	6 (10%)
21	EDO	R	204	-	3,3,3	0.24	0	2,2,2	0.49	0
21	EDO	C	314	-	3,3,3	0.66	0	2,2,2	0.62	0
21	EDO	A	614	-	3,3,3	2.72	1 (33%)	2,2,2	4.55	1 (50%)
21	EDO	P	313	-	3,3,3	0.92	0	2,2,2	0.33	0
21	EDO	R	202	-	3,3,3	0.48	0	2,2,2	1.01	0
21	EDO	P	315	-	3,3,3	0.42	0	2,2,2	0.70	0
21	EDO	E	202	-	3,3,3	0.49	0	2,2,2	0.89	0
27	CDL	G	102	-	99,99,99	1.59	15 (15%)	105,111,111	1.52	19 (18%)
21	EDO	D	205	-	3,3,3	0.44	0	2,2,2	1.13	0
28	PEK	P	304	-	52,52,52	1.02	4 (7%)	55,57,57	1.21	5 (9%)
21	EDO	G	106	-	3,3,3	0.96	0	2,2,2	0.59	0
25	DMU	Z	101	-	34,34,34	0.83	1 (2%)	45,45,45	1.25	5 (11%)
14	HEA	A	602[B]	1,18	67,67,67	1.97	21 (31%)	81,103,103	1.98	25 (30%)
21	EDO	A	618	-	3,3,3	0.39	0	2,2,2	1.35	0
21	EDO	C	316	-	3,3,3	0.55	0	2,2,2	0.38	0
21	EDO	A	611	-	3,3,3	0.54	0	2,2,2	0.99	0
21	EDO	E	205	-	3,3,3	0.51	0	2,2,2	0.79	0
21	EDO	N	621	-	3,3,3	0.77	0	2,2,2	0.19	0
21	EDO	C	319	-	3,3,3	0.85	0	2,2,2	1.42	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	E	201	-	-	0/1/1/1	-
20	PGV	P	302	-	-	23/55/55/55	-
21	EDO	F	103	-	-	0/1/1/1	-
23	PSC	O	302	-	-	30/55/55/55	-
25	DMU	C	309	-	-	6/19/59/59	0/2/2/2
21	EDO	G	105	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
14	HEA	A	602[A]	1,18	-	8/36/76/76	-
21	EDO	C	311	-	-	0/1/1/1	-
21	EDO	P	312	-	-	1/1/1/1	-
21	EDO	E	203	-	-	1/1/1/1	-
21	EDO	A	619	-	-	1/1/1/1	-
21	EDO	A	610	-	-	1/1/1/1	-
21	EDO	D	202	-	-	0/1/1/1	-
14	HEA	N	601	1	-	6/36/76/76	-
21	EDO	A	615	-	-	0/1/1/1	-
21	EDO	A	617	-	-	1/1/1/1	-
21	EDO	C	318	-	-	1/1/1/1	-
19	TGL	N	610	-	-	37/65/65/65	-
21	EDO	O	305	-	-	1/1/1/1	-
24	CHD	C	306	-	-	6/9/74/74	0/4/4/4
28	PEK	T	102	-	-	21/56/56/56	-
21	EDO	R	203	-	-	0/1/1/1	-
21	EDO	O	304	-	-	1/1/1/1	-
20	PGV	C	308	-	-	26/55/55/55	-
21	EDO	N	616	-	-	0/1/1/1	-
25	DMU	C	310	-	-	10/19/59/59	0/2/2/2
24	CHD	G	103	-	-	2/9/74/74	0/4/4/4
21	EDO	N	620	-	-	0/1/1/1	-
20	PGV	M	101	-	-	22/55/55/55	-
25	DMU	P	311	-	-	13/19/59/59	0/2/2/2
27	CDL	C	305	-	-	66/110/110/110	-
21	EDO	R	201	-	-	1/1/1/1	-
21	EDO	C	317	-	-	1/1/1/1	-
21	EDO	B	303	-	-	0/1/1/1	-
19	TGL	A	608	-	-	38/65/65/65	-
21	EDO	S	103	-	-	1/1/1/1	-
19	TGL	D	201	-	-	42/65/65/65	-
21	EDO	C	313	-	-	0/1/1/1	-
21	EDO	Q	204	-	-	0/1/1/1	-
21	EDO	Q	202	-	-	0/1/1/1	-
21	EDO	N	612	-	-	1/1/1/1	-
25	DMU	P	310	-	-	7/19/59/59	0/2/2/2
20	PGV	C	304	-	-	18/55/55/55	-
28	PEK	G	104	-	-	35/56/56/56	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
27	CDL	T	103	-	-	52/110/110/110	-
24	CHD	T	101	-	-	2/9/74/74	0/4/4/4
25	DMU	P	308	-	-	6/19/59/59	0/2/2/2
21	EDO	F	102	-	-	0/1/1/1	-
21	EDO	D	203	-	-	1/1/1/1	-
25	DMU	M	102	-	-	5/19/59/59	0/2/2/2
24	CHD	P	301	-	-	2/9/74/74	0/4/4/4
21	EDO	A	613	-	-	1/1/1/1	-
19	TGL	Q	201	-	-	39/65/65/65	-
14	HEA	A	601	1	-	5/36/76/76	-
21	EDO	C	312	-	-	1/1/1/1	-
21	EDO	P	314	-	-	0/1/1/1	-
21	EDO	A	620	-	-	0/1/1/1	-
21	EDO	O	303	-	-	0/1/1/1	-
27	CDL	P	306	-	-	61/110/110/110	-
24	CHD	P	307	-	-	7/9/74/74	0/4/4/4
20	PGV	A	609	-	-	8/55/55/55	-
21	EDO	N	619	-	-	1/1/1/1	-
19	TGL	L	101	-	-	43/65/65/65	-
21	EDO	D	204	-	-	1/1/1/1	-
14	HEA	N	602[B]	1,18	-	8/36/76/76	-
21	EDO	N	618	-	-	0/1/1/1	-
25	DMU	C	302	-	-	7/19/59/59	0/2/2/2
21	EDO	N	613	-	-	1/1/1/1	-
21	EDO	A	616	-	-	1/1/1/1	-
21	EDO	U	101	-	-	1/1/1/1	-
20	PGV	P	305	-	-	9/55/55/55	-
21	EDO	V	101	-	-	1/1/1/1	-
21	EDO	N	617	-	-	0/1/1/1	-
21	EDO	E	204	-	-	1/1/1/1	-
20	PGV	N	609	-	-	9/55/55/55	-
21	EDO	C	315	-	-	0/1/1/1	-
21	EDO	S	104	-	-	0/1/1/1	-
21	EDO	T	104	-	-	0/1/1/1	-
28	PEK	G	101	-	-	18/56/56/56	-
21	EDO	H	101	-	-	0/1/1/1	-
23	PSC	B	302	-	-	29/55/55/55	-
21	EDO	Q	203	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
21	EDO	N	611	-	-	1/1/1/1	-
21	EDO	S	102	-	-	0/1/1/1	-
21	EDO	F	104	-	-	0/1/1/1	-
14	HEA	N	602[A]	1,18	-	6/36/76/76	-
21	EDO	D	206	-	-	1/1/1/1	-
24	CHD	C	301	-	-	2/9/74/74	0/4/4/4
20	PGV	N	608	-	-	25/55/55/55	-
21	EDO	N	614	-	-	1/1/1/1	-
19	TGL	Y	101	-	-	34/65/65/65	-
21	EDO	A	612	-	-	1/1/1/1	-
21	EDO	N	615	-	-	0/1/1/1	-
28	PEK	C	307	-	-	30/56/56/56	-
21	EDO	R	205	-	-	1/1/1/1	-
28	PEK	P	309	-	-	27/56/56/56	-
21	EDO	R	204	-	-	1/1/1/1	-
21	EDO	C	314	-	-	1/1/1/1	-
21	EDO	A	614	-	-	1/1/1/1	-
21	EDO	P	313	-	-	0/1/1/1	-
21	EDO	R	202	-	-	1/1/1/1	-
21	EDO	P	315	-	-	0/1/1/1	-
21	EDO	E	202	-	-	0/1/1/1	-
27	CDL	G	102	-	-	53/110/110/110	-
21	EDO	D	205	-	-	0/1/1/1	-
28	PEK	P	304	-	-	16/56/56/56	-
21	EDO	G	106	-	-	0/1/1/1	-
25	DMU	Z	101	-	-	3/19/59/59	0/2/2/2
14	HEA	A	602[B]	1,18	-	7/36/76/76	-
21	EDO	A	618	-	-	0/1/1/1	-
21	EDO	C	316	-	-	1/1/1/1	-
21	EDO	A	611	-	-	1/1/1/1	-
21	EDO	E	205	-	-	1/1/1/1	-
21	EDO	N	621	-	-	0/1/1/1	-
21	EDO	C	319	-	-	1/1/1/1	-

The worst 5 of 344 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
19	D	201	TGL	OB1-CB1	8.46	1.47	1.22
19	Q	201	TGL	OB1-CB1	7.15	1.43	1.22
19	Y	101	TGL	OG2-CB1	6.54	1.52	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
19	D	201	TGL	OG1-CA1	6.49	1.52	1.33
19	D	201	TGL	OG2-CB1	6.34	1.52	1.34

The worst 5 of 609 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
24	C	306	CHD	C23-C22-C20	-11.76	92.49	114.46
19	D	201	TGL	OG2-CB1-CB2	-10.33	89.13	111.48
25	C	310	DMU	O16-C6-C1	9.86	123.24	108.27
25	C	309	DMU	C10-O1-C9	-9.21	95.73	113.72
19	D	201	TGL	OG2-CB1-OB1	8.91	144.53	123.70

There are no chirality outliers.

5 of 965 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
14	A	602[A]	HEA	C2A-C3A-CMA-OMA
14	A	602[B]	HEA	C2A-C3A-CMA-OMA
14	A	602[B]	HEA	C4D-C3D-CAD-CBD
14	N	602[A]	HEA	C2A-C3A-CMA-OMA
19	D	201	TGL	CG2-CG1-OG1-CA1

There are no ring outliers.

70 monomers are involved in 425 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
18	A	606	AZI	1	0
20	P	302	PGV	10	0
23	O	302	PSC	16	0
25	C	309	DMU	2	0
21	G	105	EDO	1	0
14	A	602[A]	HEA	5	0
21	P	312	EDO	1	0
21	A	610	EDO	1	0
18	N	606	AZI	1	0
21	D	202	EDO	6	0
14	N	601	HEA	9	0
21	A	617	EDO	2	0
19	N	610	TGL	4	0
24	C	306	CHD	5	0
28	T	102	PEK	10	0

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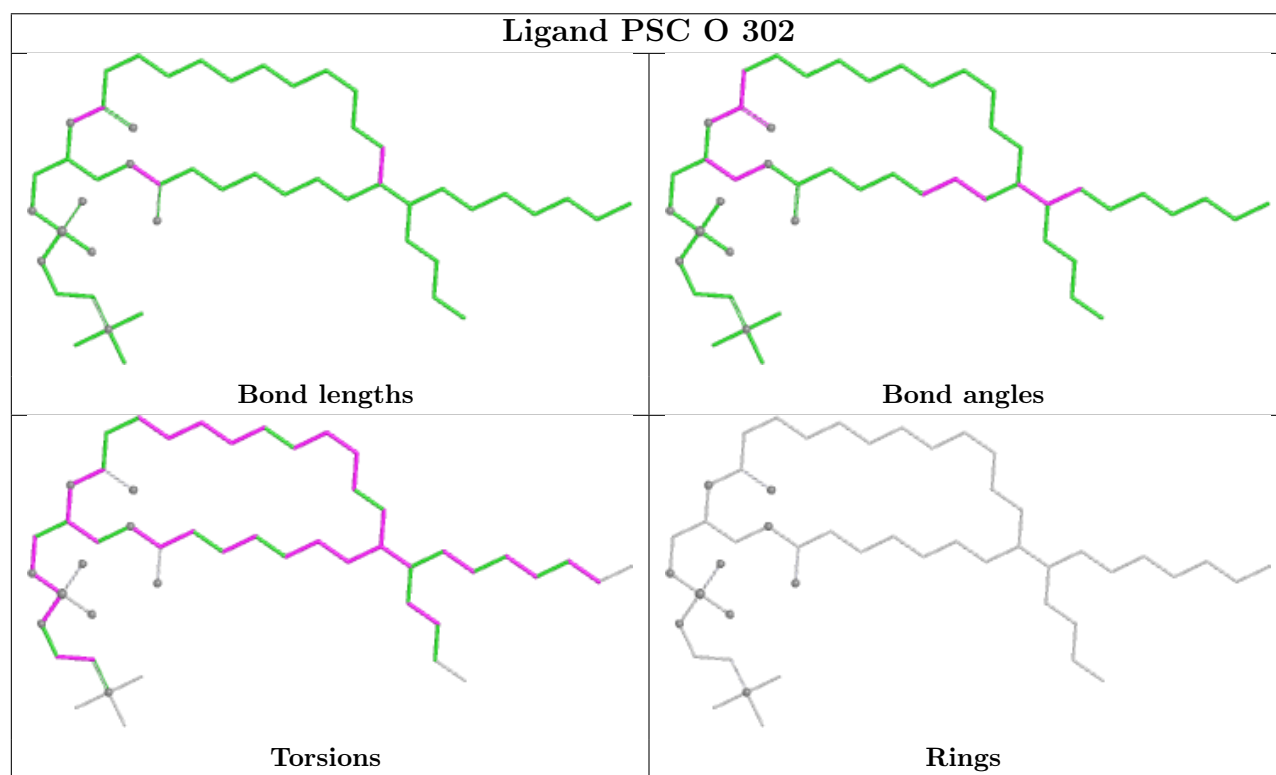
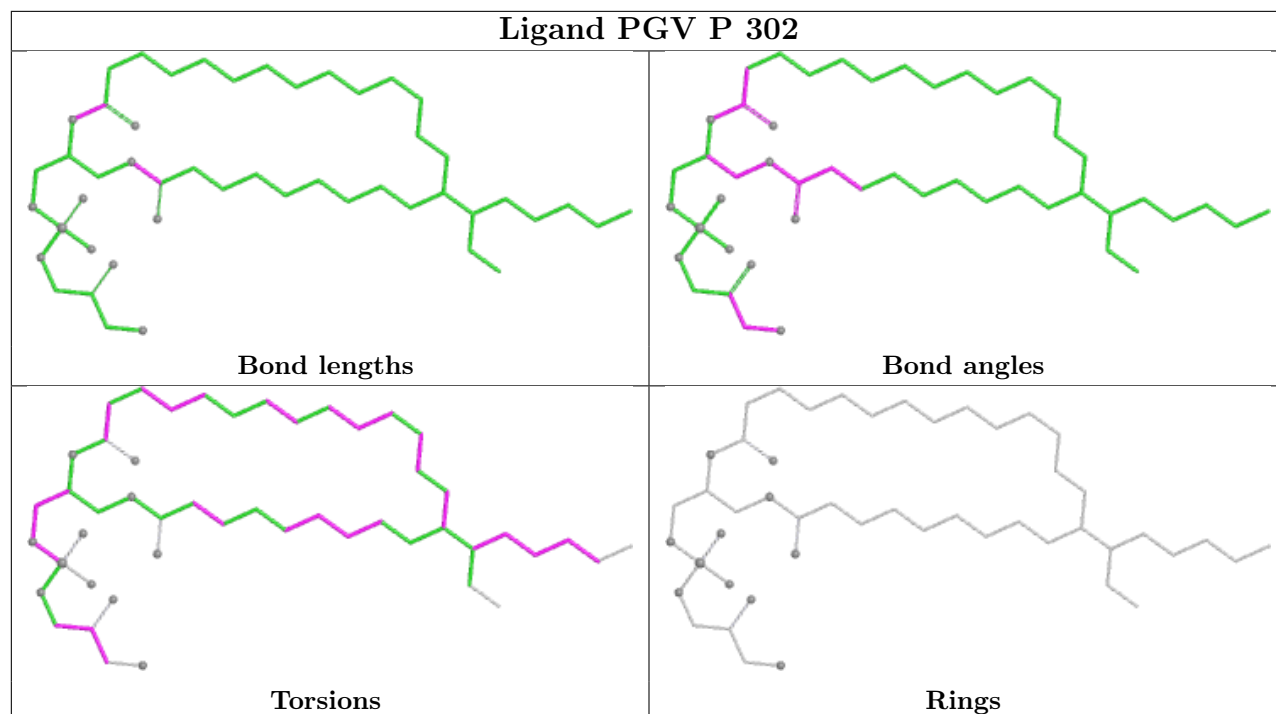
Mol	Chain	Res	Type	Clashes	Symm-Clashes
20	C	308	PGV	2	0
18	N	607	AZI	6	0
25	C	310	DMU	7	0
24	G	103	CHD	1	0
20	M	101	PGV	3	0
18	A	607	AZI	6	0
27	C	305	CDL	25	0
21	C	317	EDO	1	0
19	A	608	TGL	4	0
21	S	103	EDO	3	0
19	D	201	TGL	13	0
21	C	313	EDO	1	0
21	Q	204	EDO	5	0
21	N	612	EDO	2	0
25	P	310	DMU	2	0
20	C	304	PGV	5	0
28	G	104	PEK	14	0
27	T	103	CDL	25	0
25	P	308	DMU	12	0
21	D	203	EDO	4	0
24	P	301	CHD	2	0
19	Q	201	TGL	12	0
14	A	601	HEA	7	0
21	C	312	EDO	1	0
21	A	620	EDO	3	0
27	P	306	CDL	22	0
24	P	307	CHD	4	0
20	A	609	PGV	4	0
19	L	101	TGL	17	0
14	N	602[B]	HEA	7	0
21	N	618	EDO	5	0
25	C	302	DMU	13	0
21	A	616	EDO	1	0
20	P	305	PGV	2	0
20	N	609	PGV	2	0
28	G	101	PEK	9	0
21	H	101	EDO	3	0
23	B	302	PSC	13	0
21	Q	203	EDO	1	0
21	F	104	EDO	1	0
14	N	602[A]	HEA	6	0
21	D	206	EDO	2	0

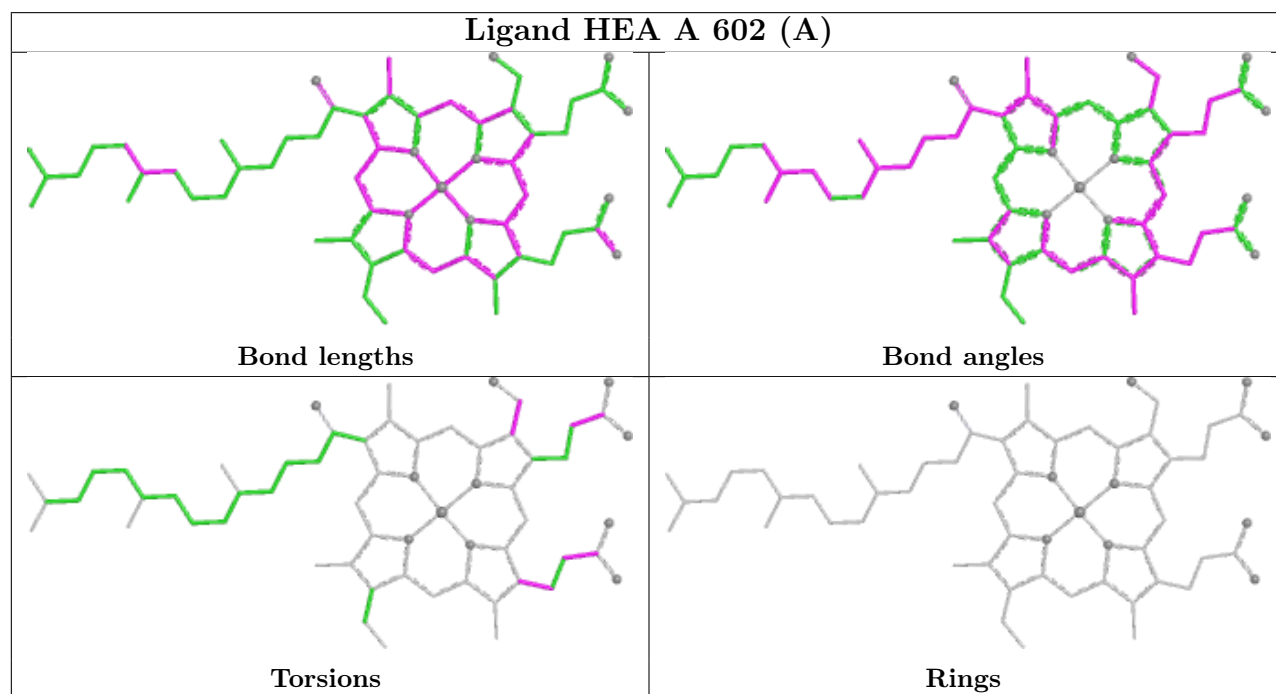
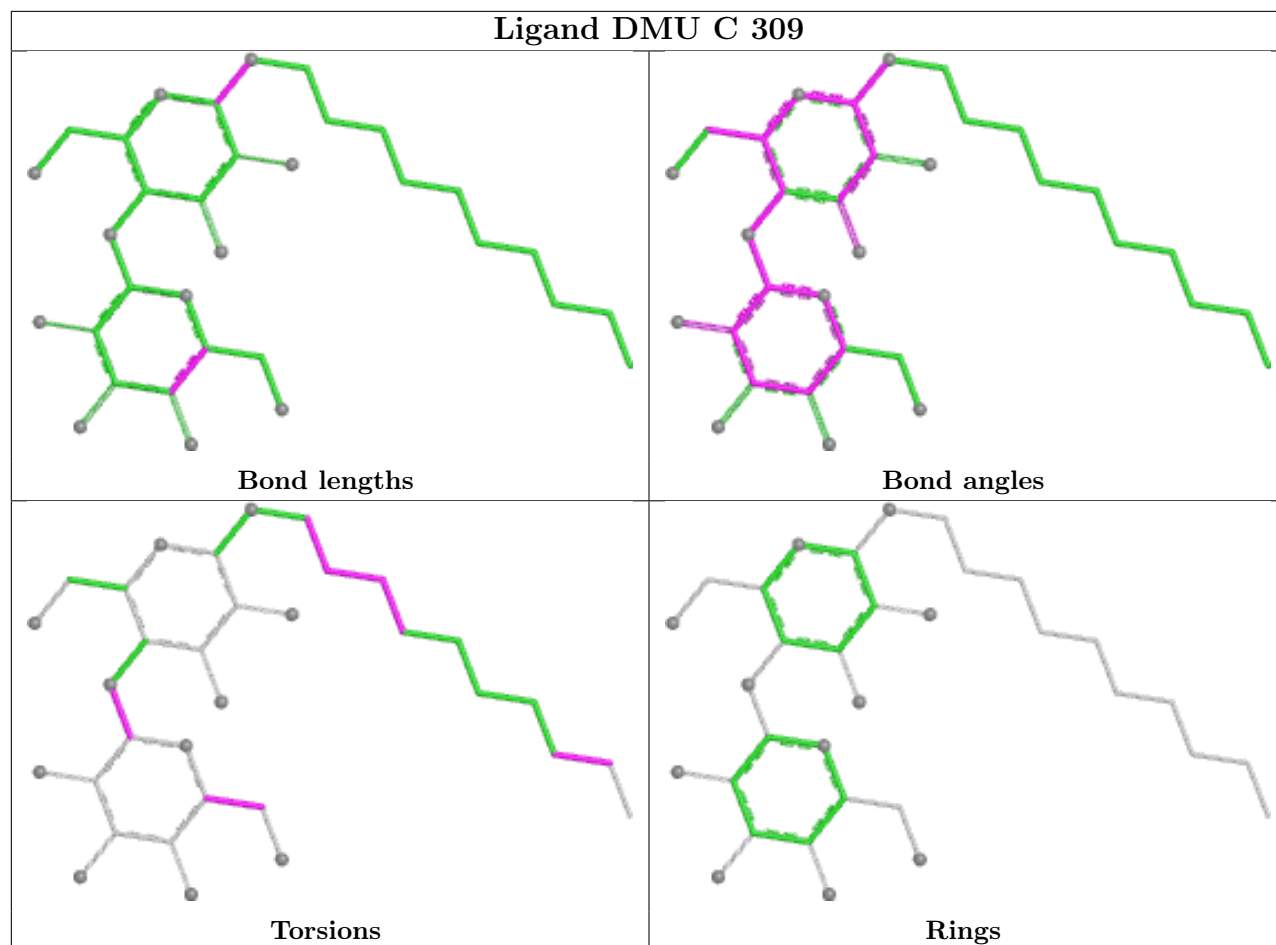
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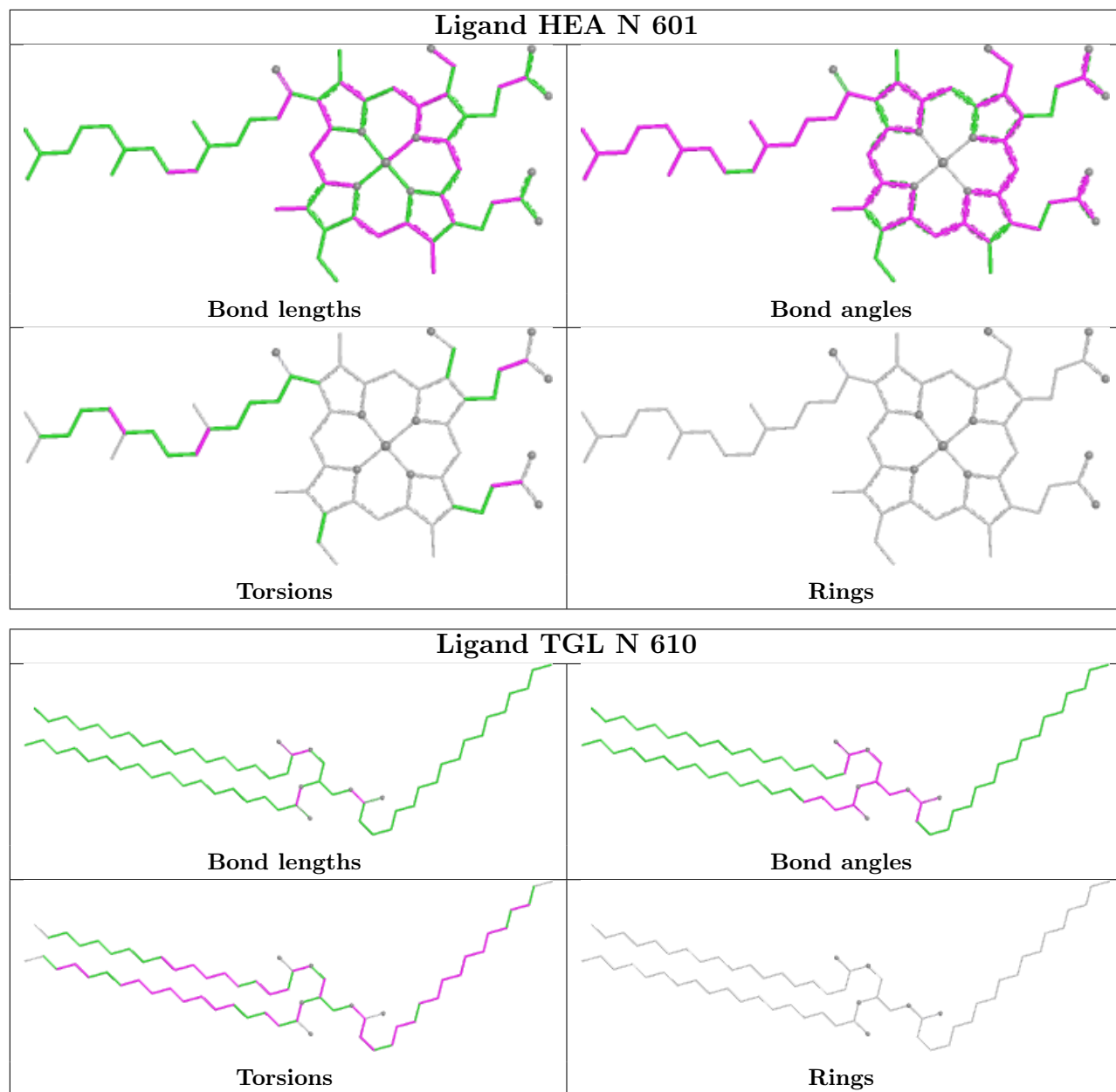
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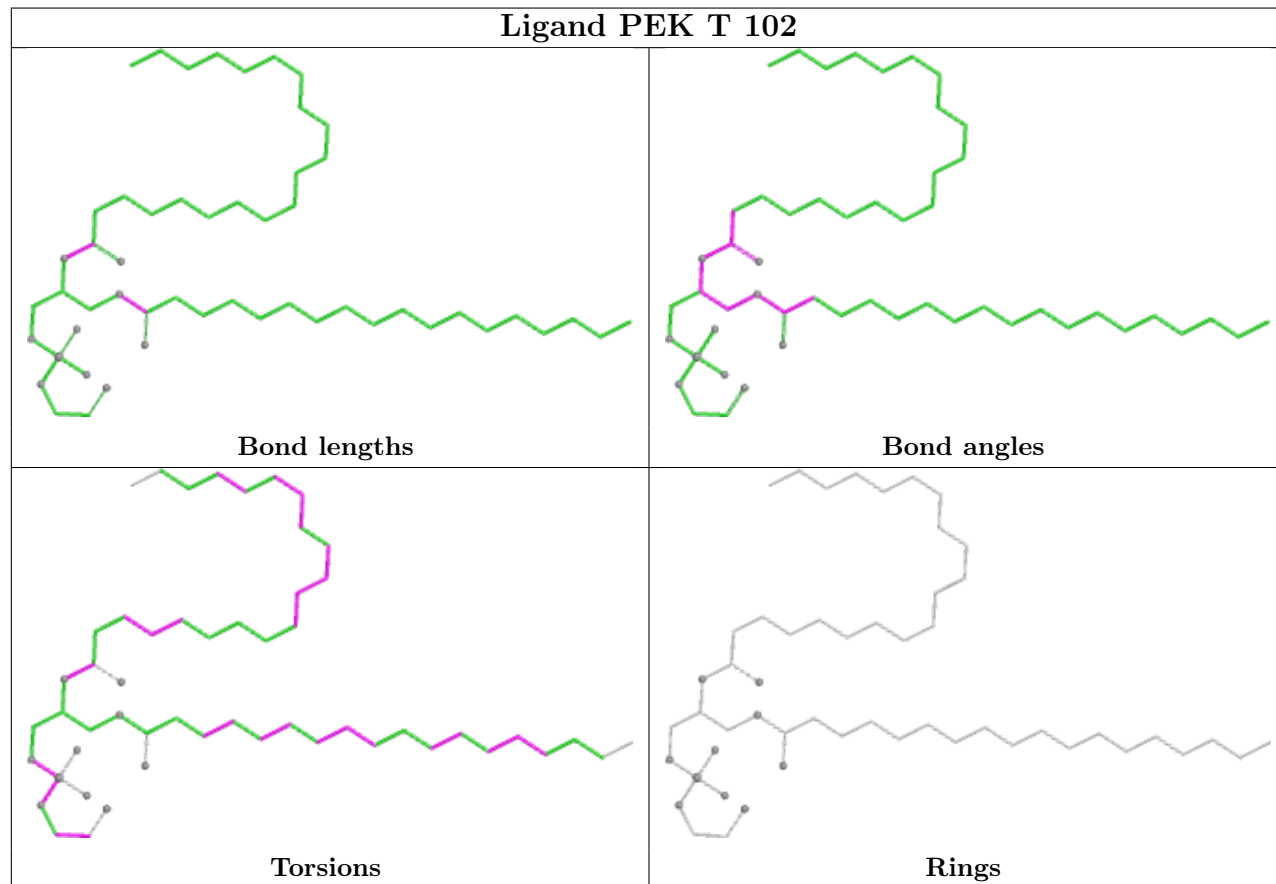
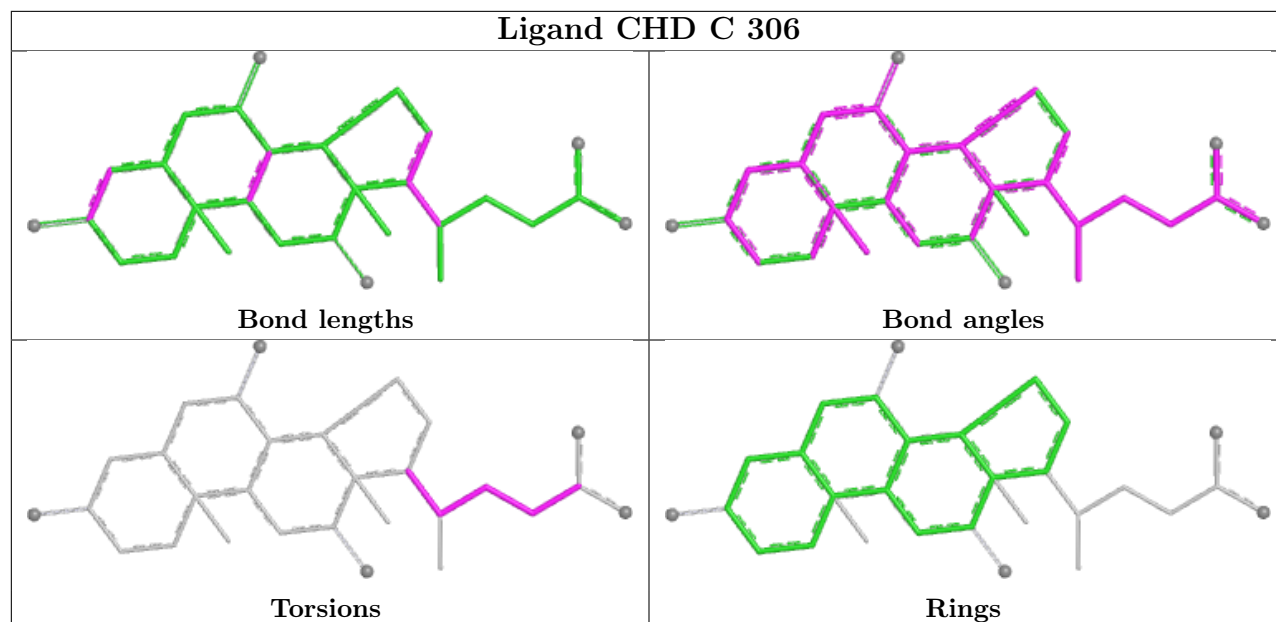
Mol	Chain	Res	Type	Clashes	Symm-Clashes
20	N	608	PGV	12	0
19	Y	101	TGL	16	0
21	A	612	EDO	4	0
28	C	307	PEK	10	0
28	P	309	PEK	2	0
21	R	204	EDO	3	0
21	C	314	EDO	1	0
21	A	614	EDO	5	0
27	G	102	CDL	24	0
28	P	304	PEK	3	0
14	A	602[B]	HEA	6	0
21	A	618	EDO	4	0
21	N	621	EDO	8	0

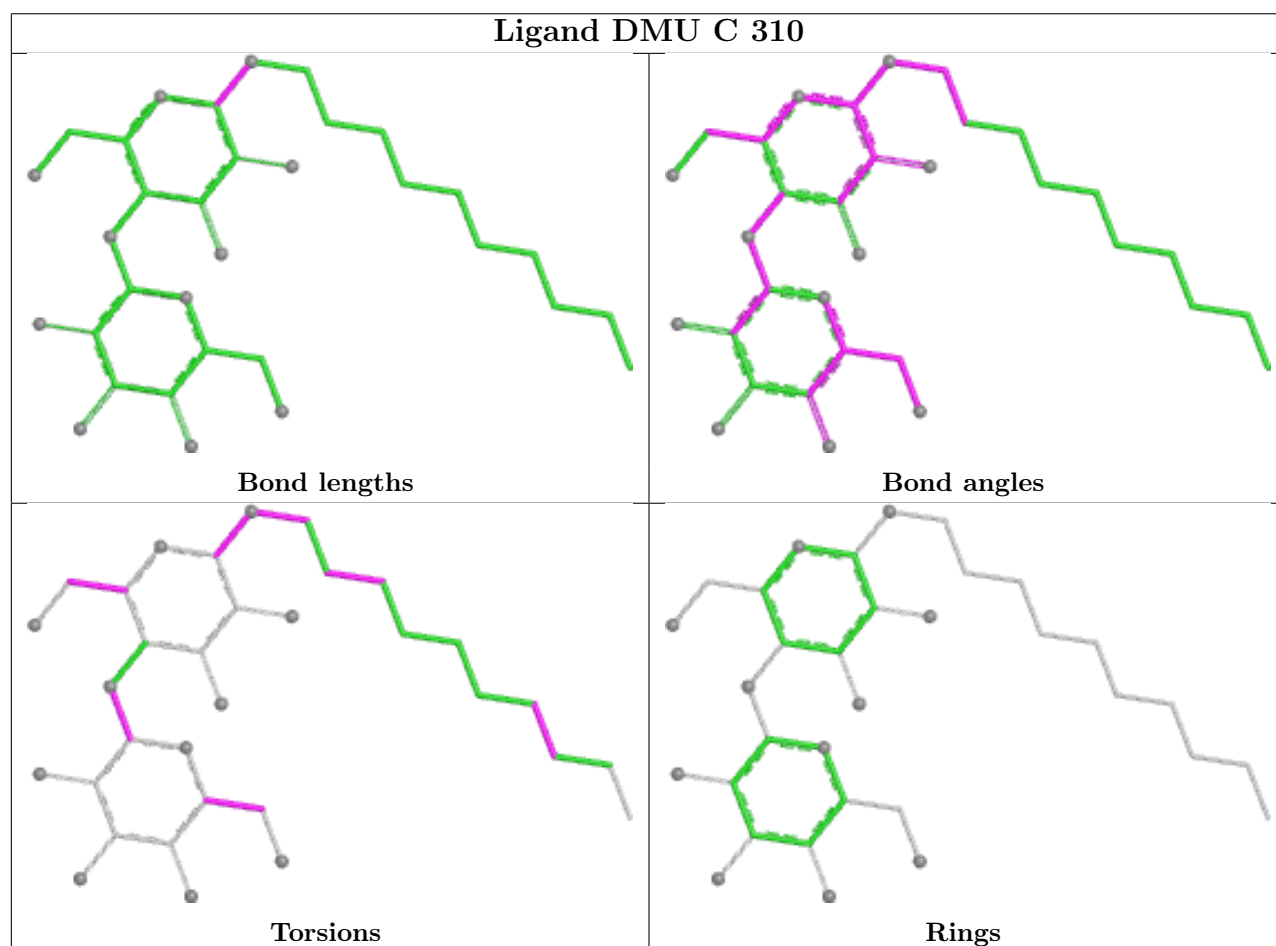
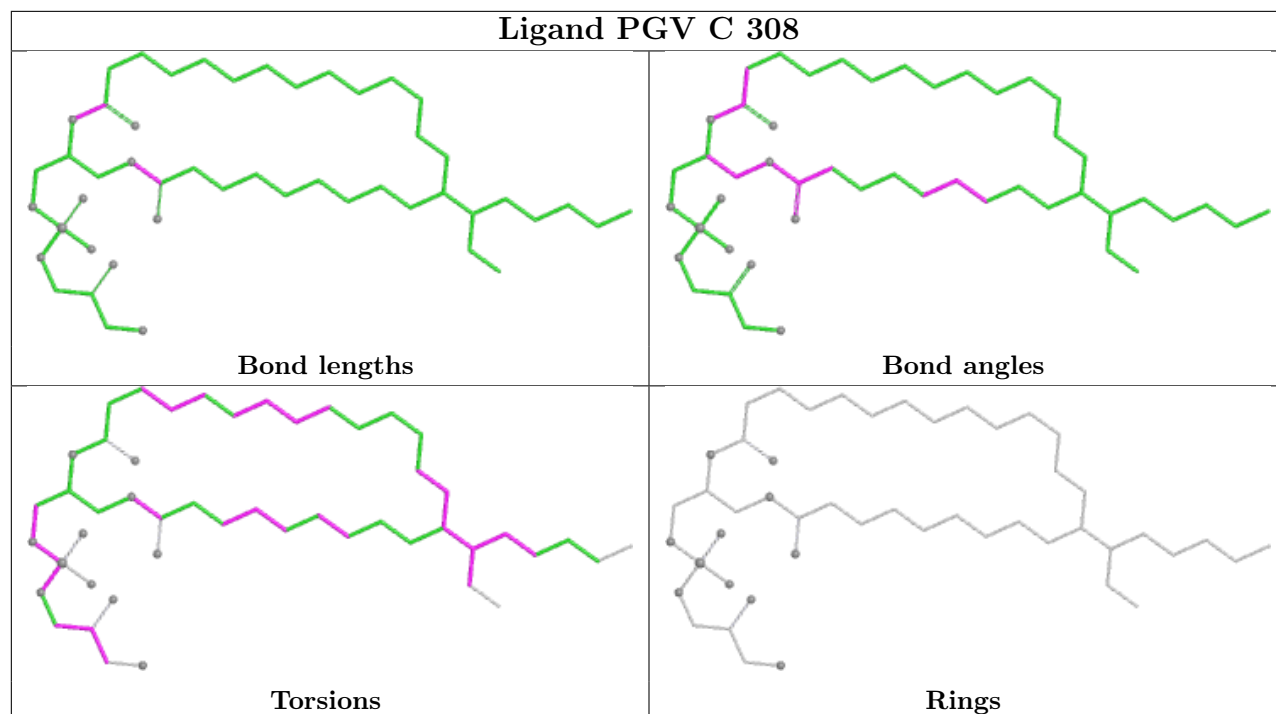
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

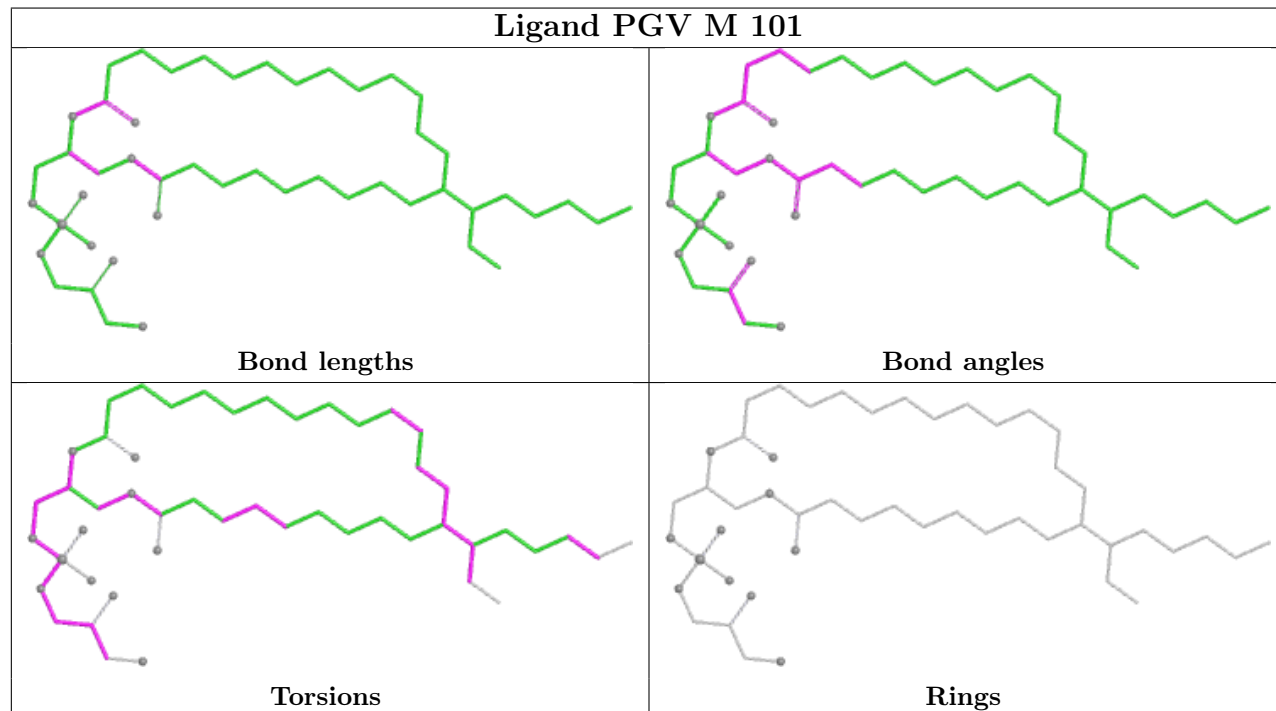
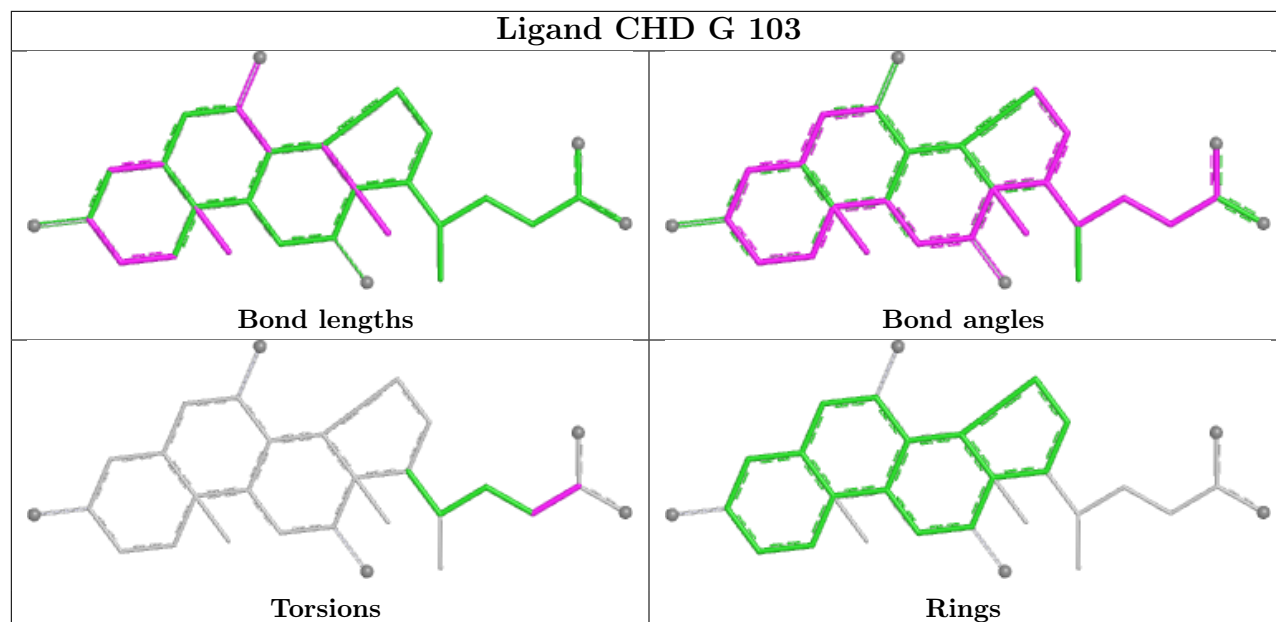


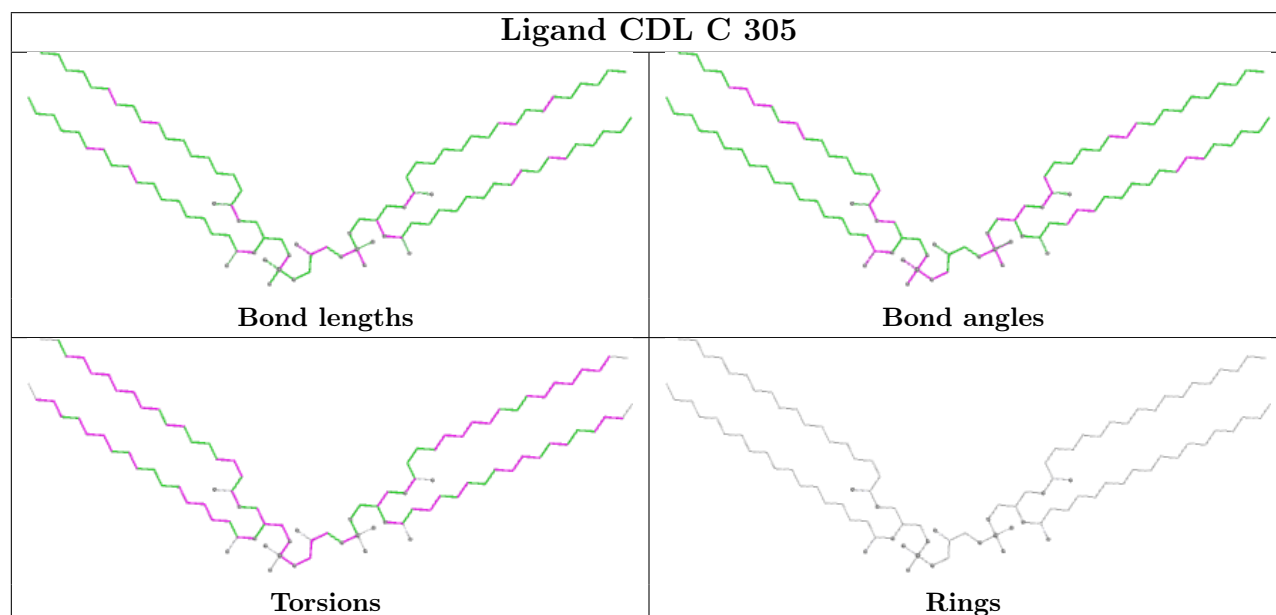
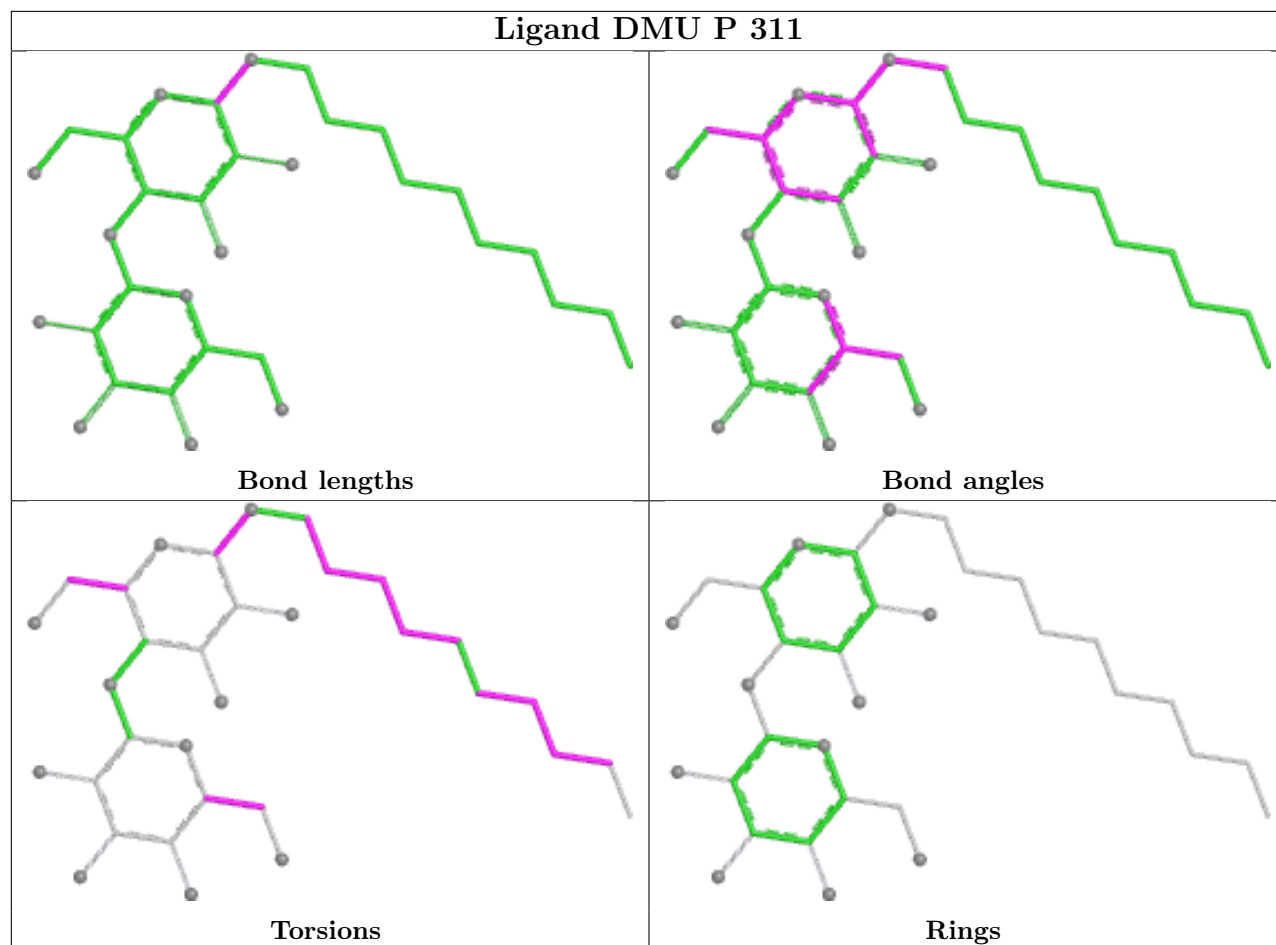


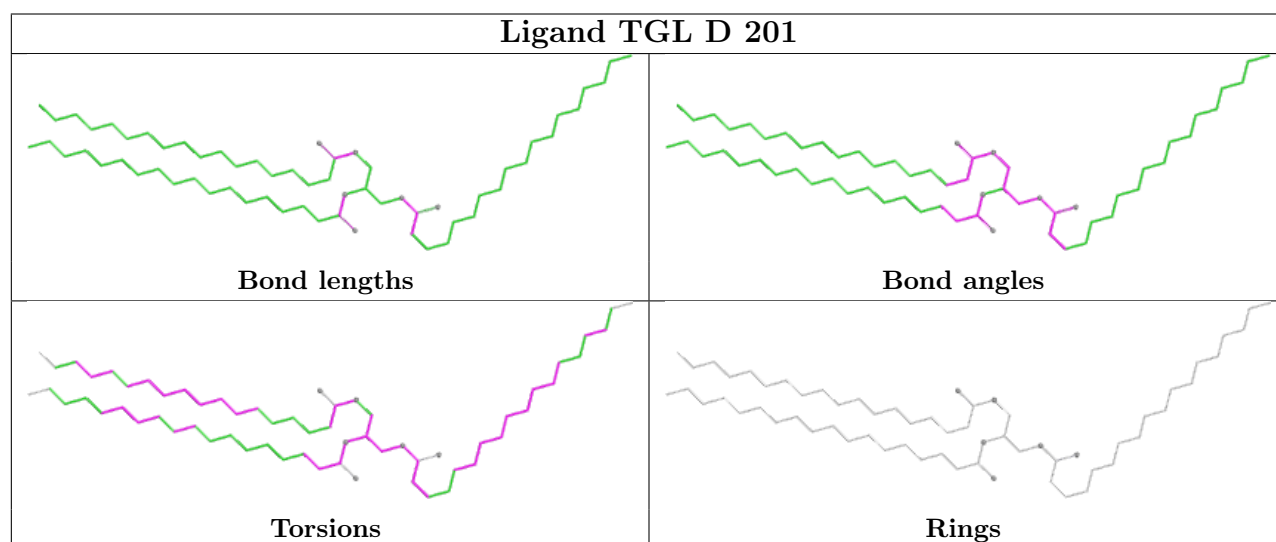
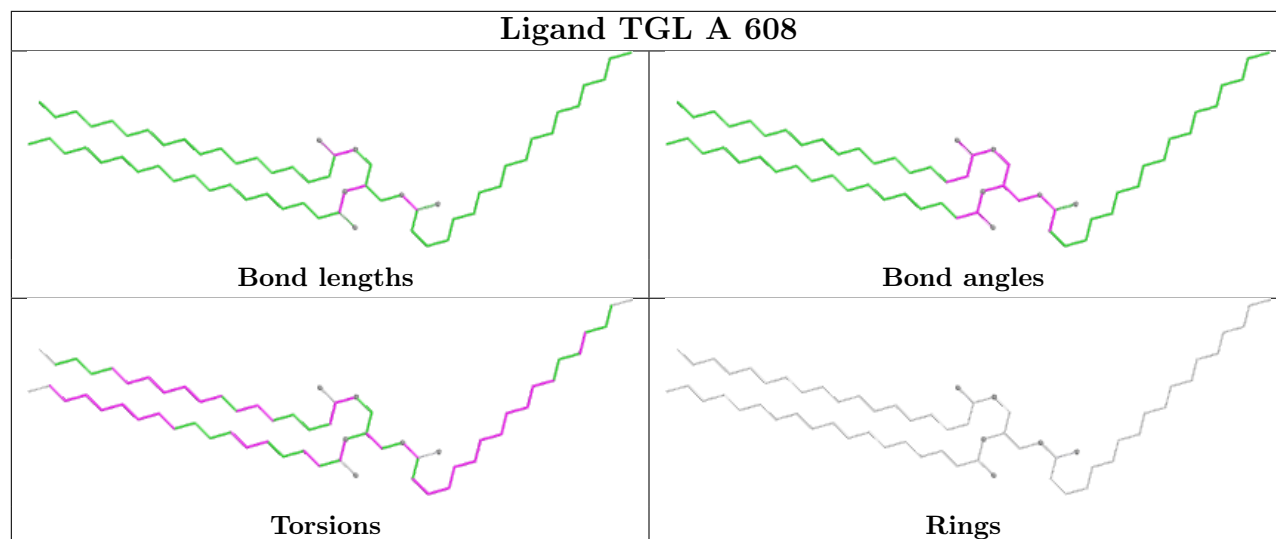


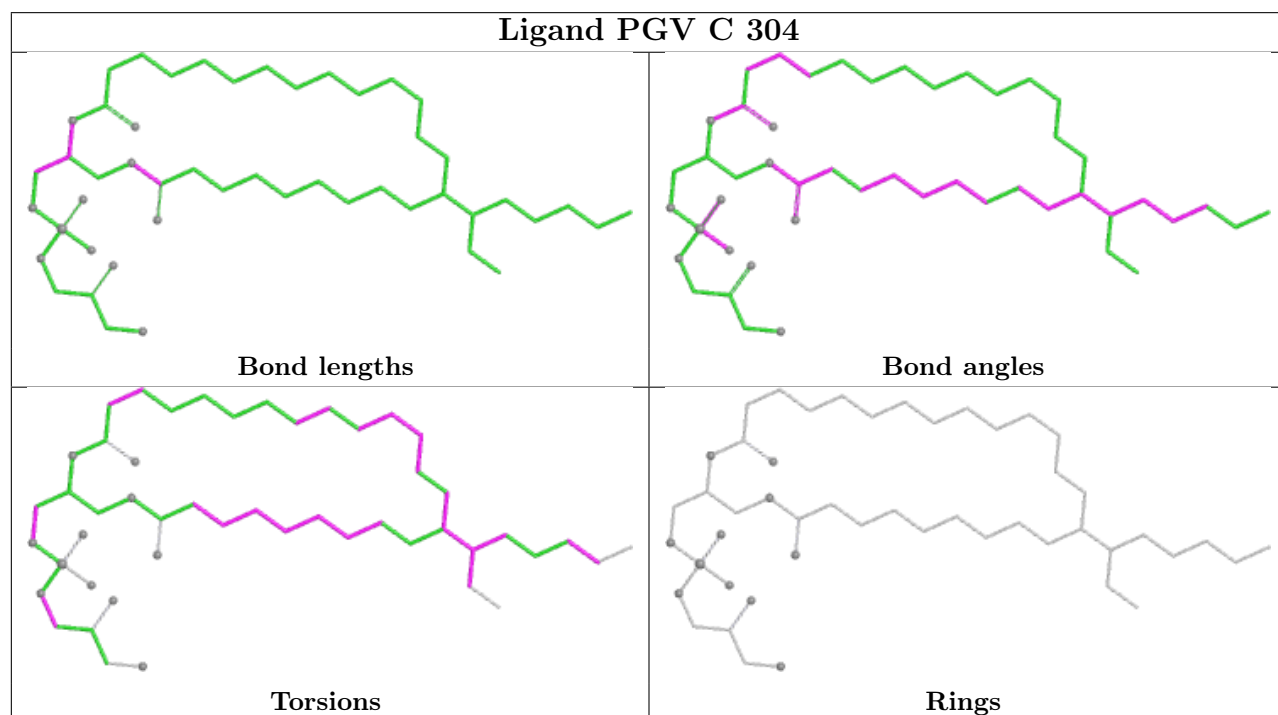
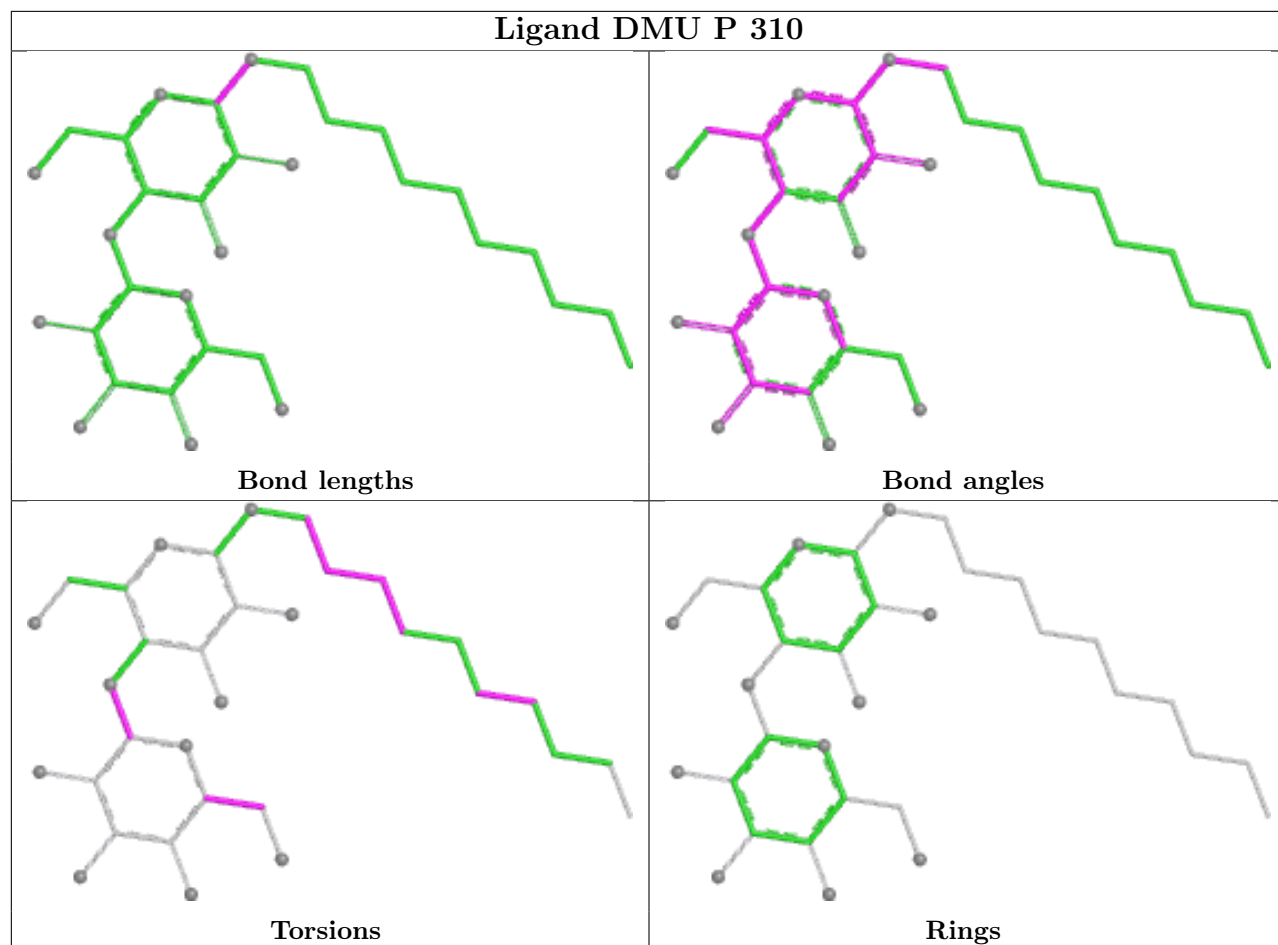


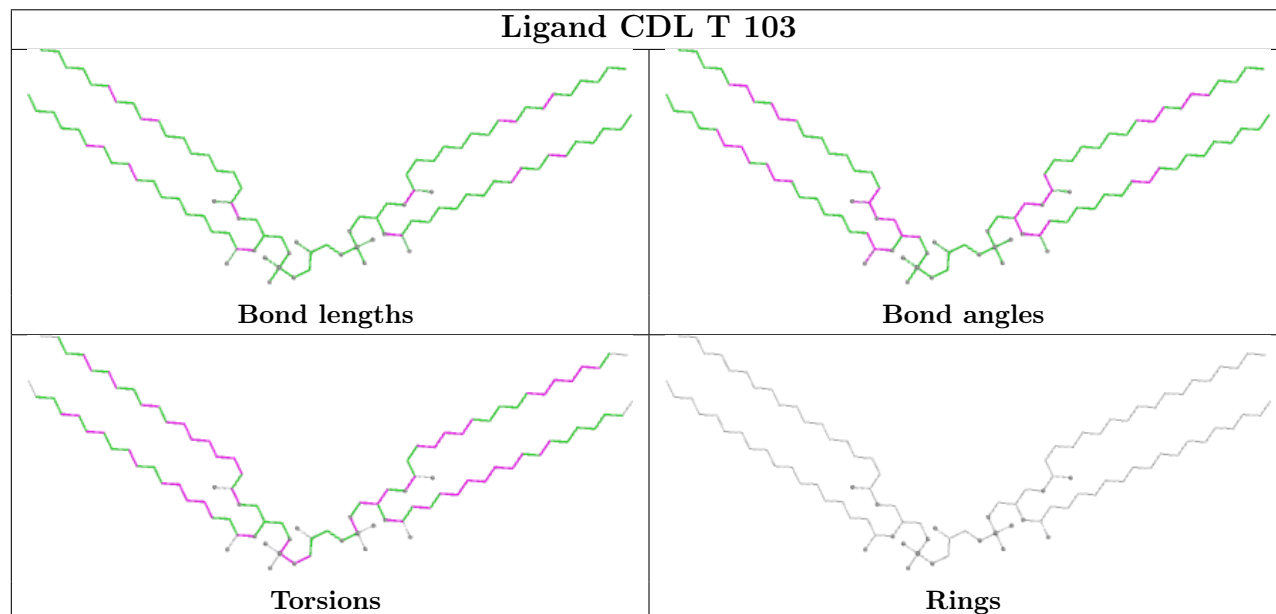
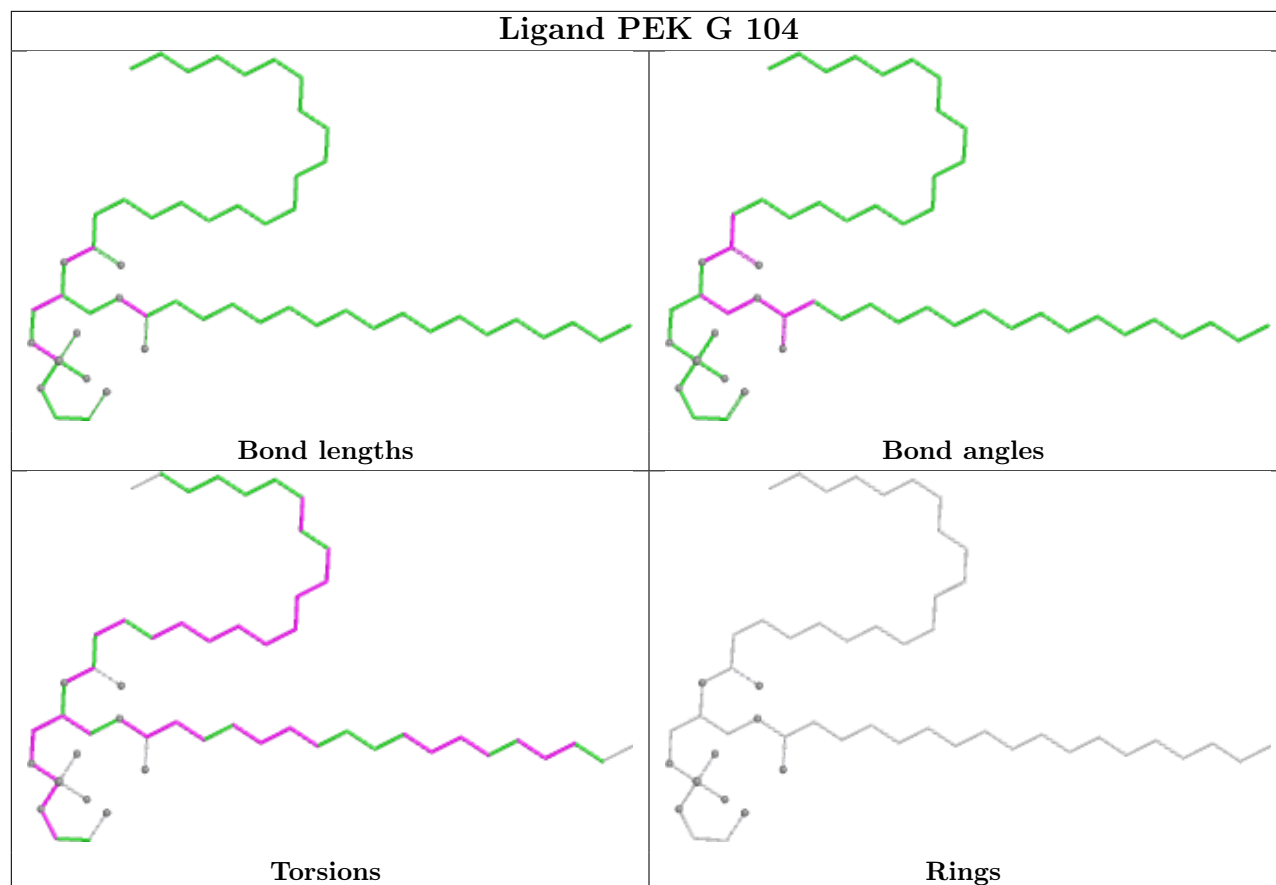


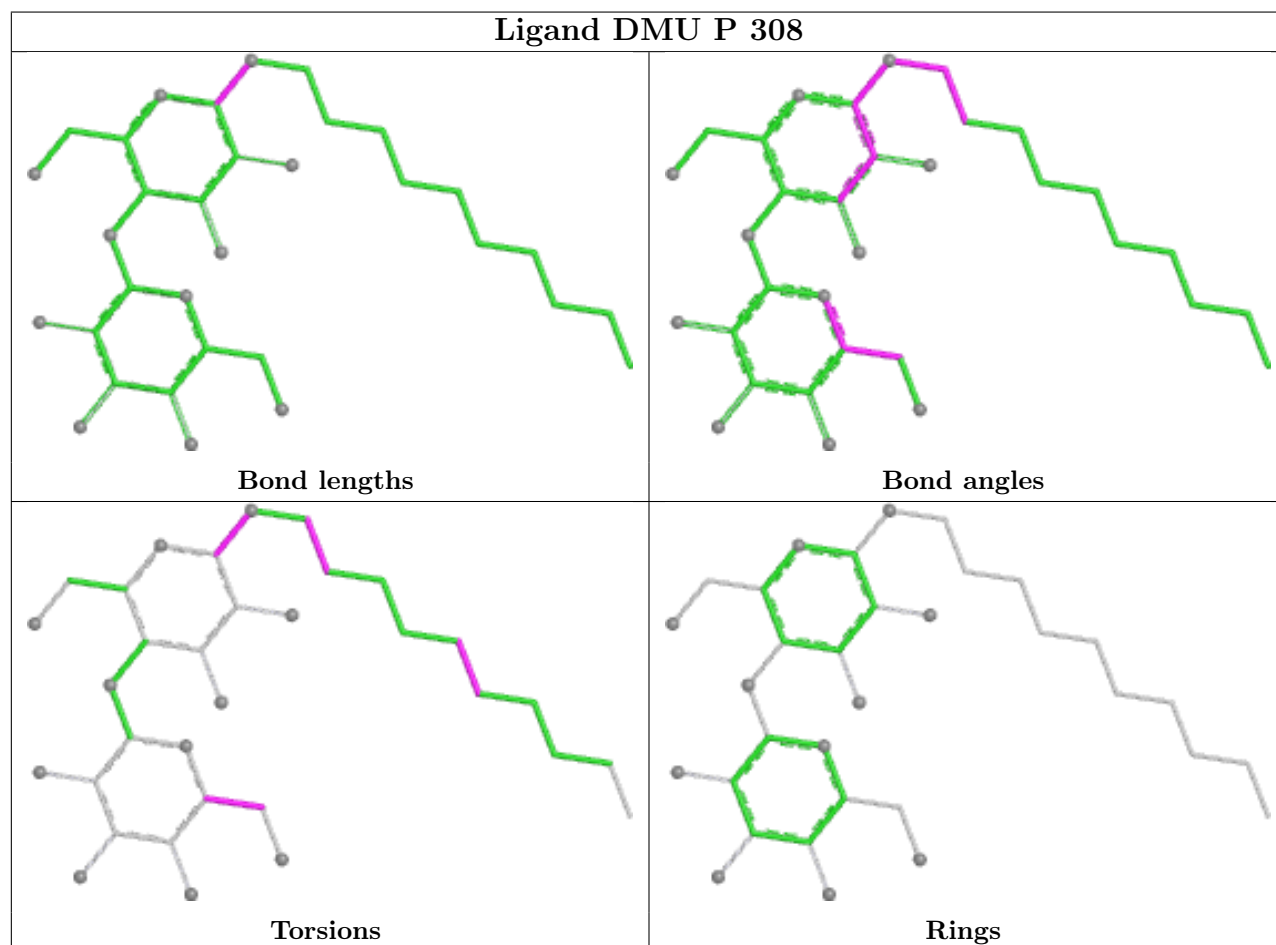
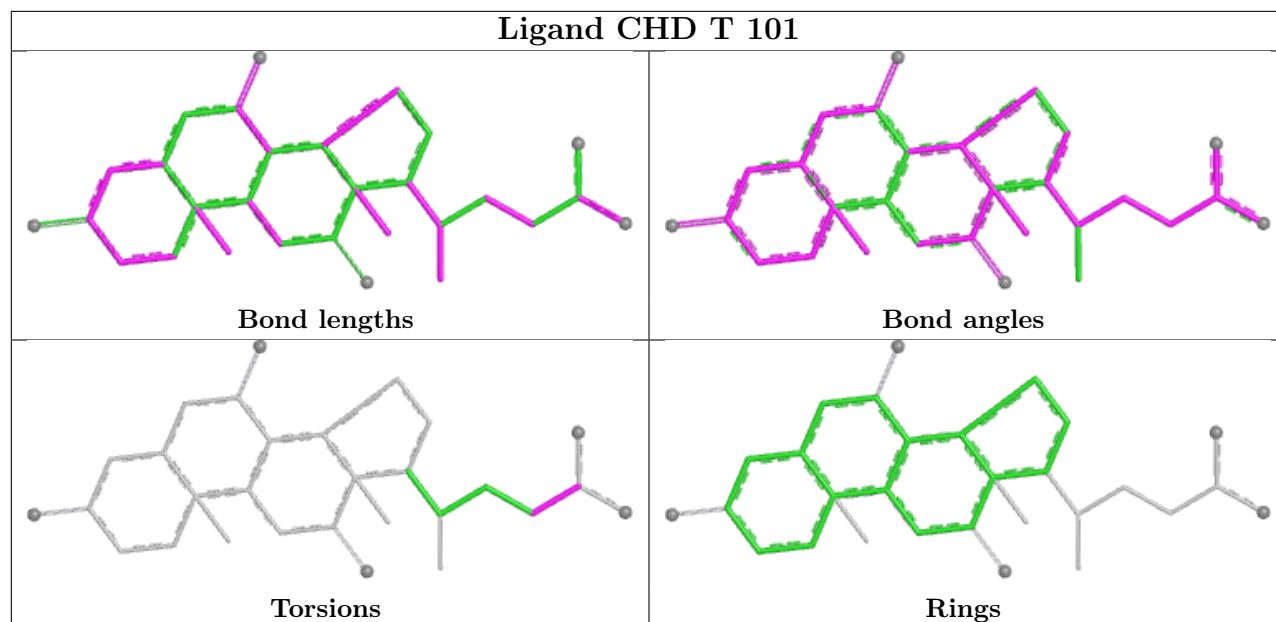




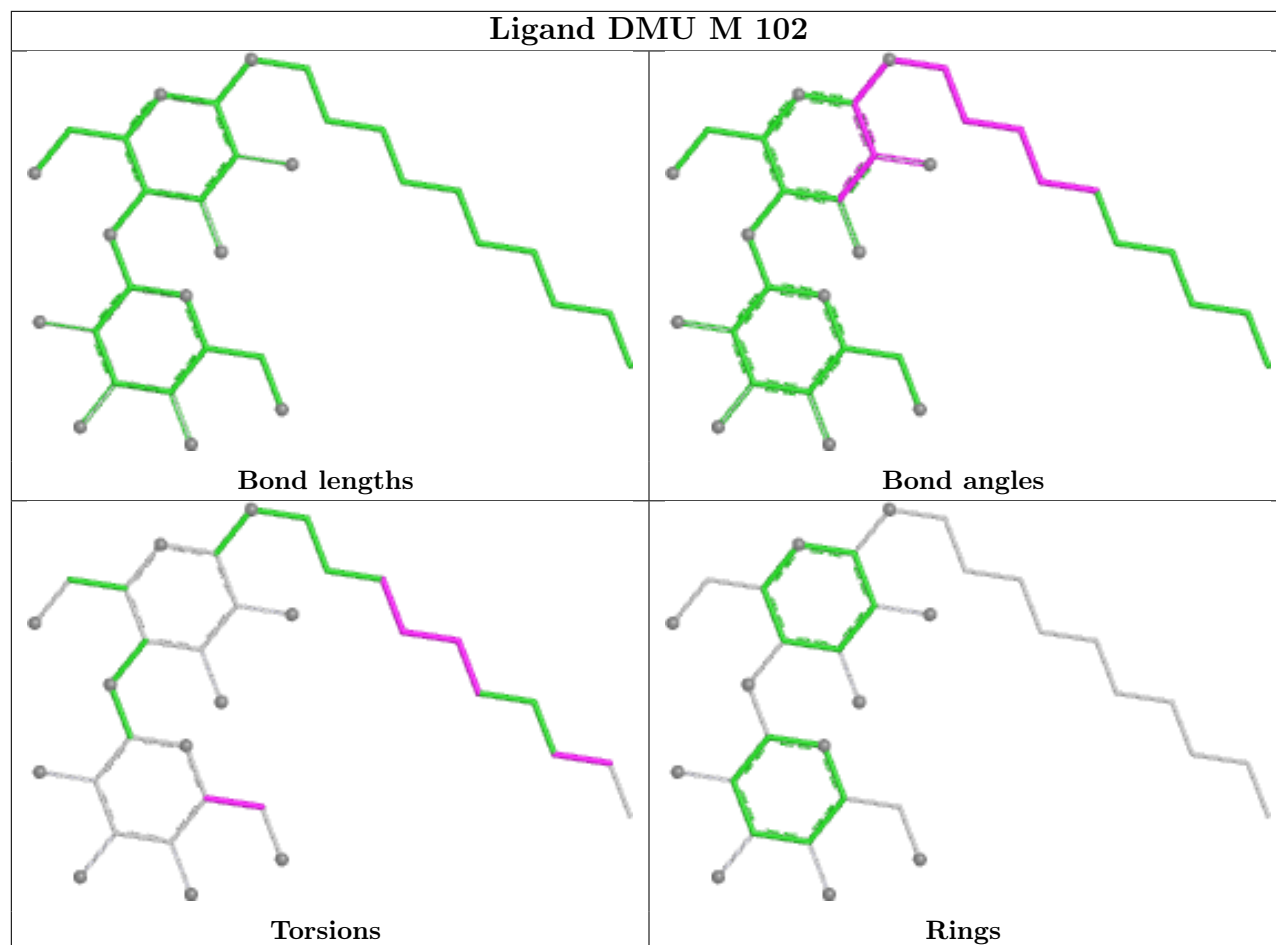




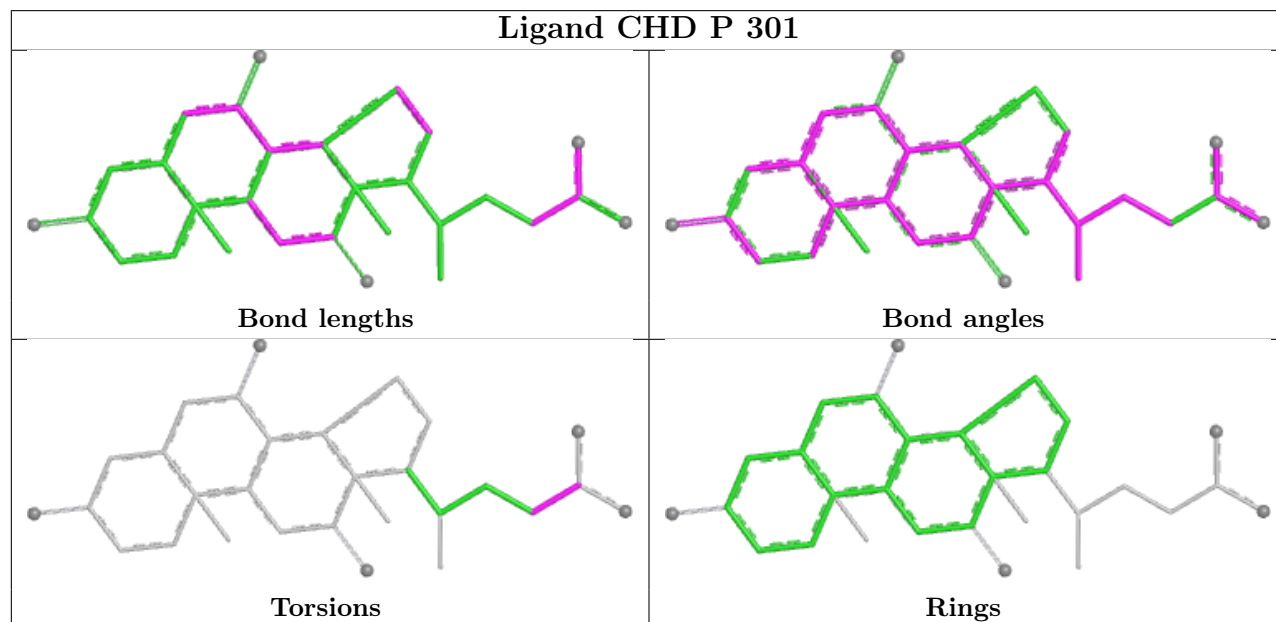


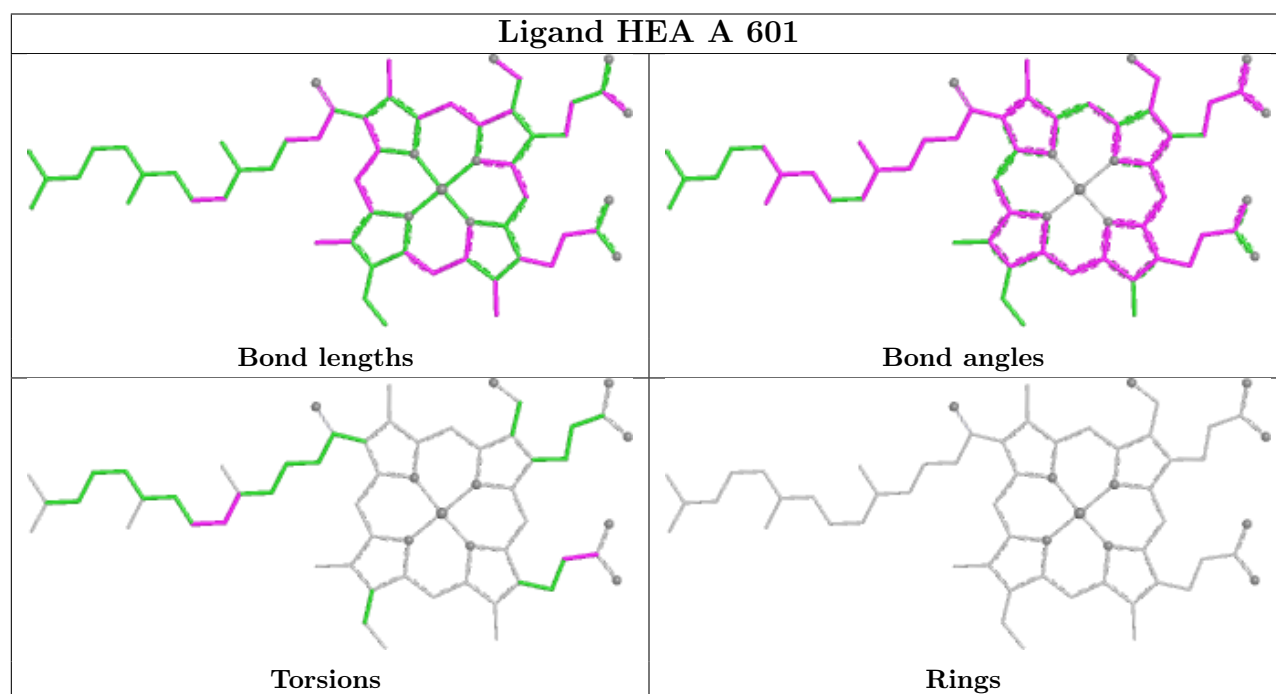
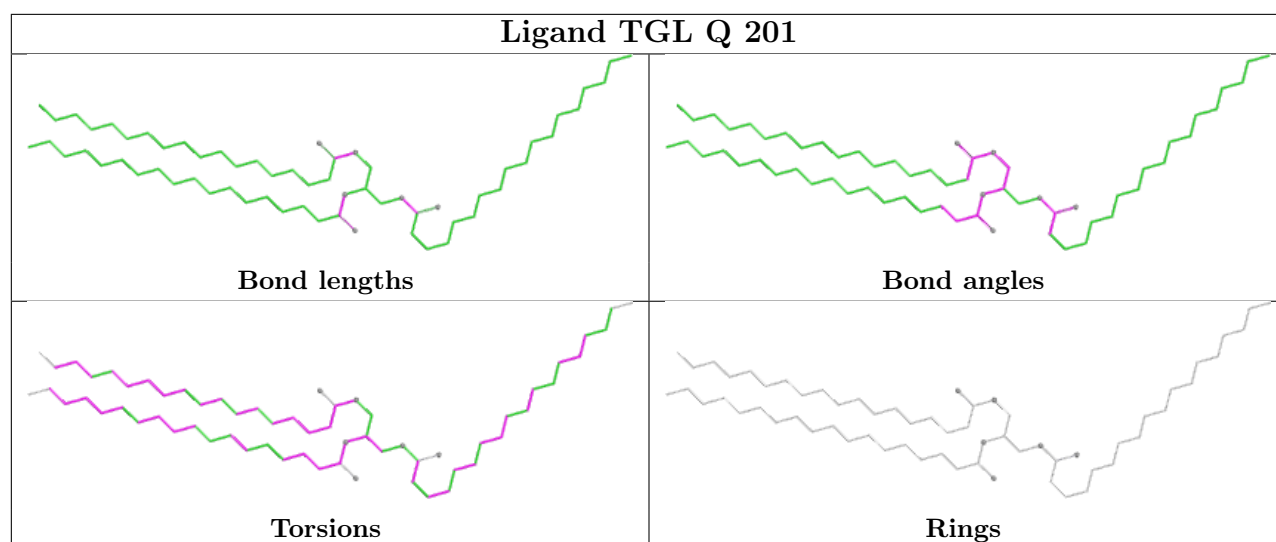


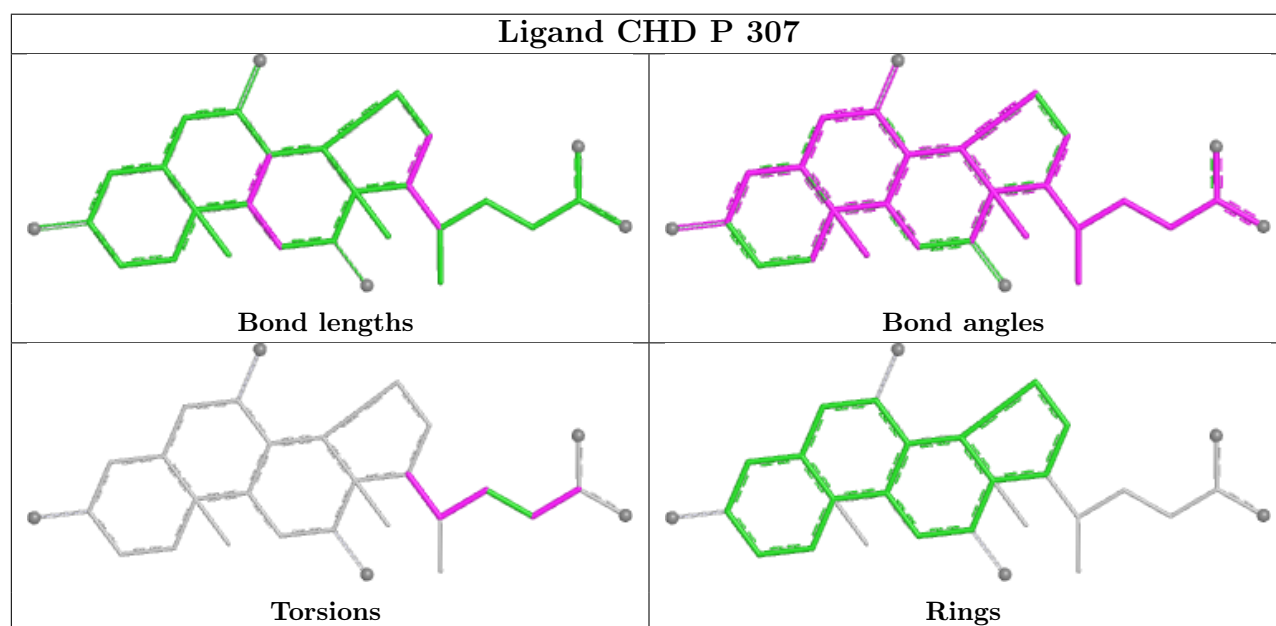
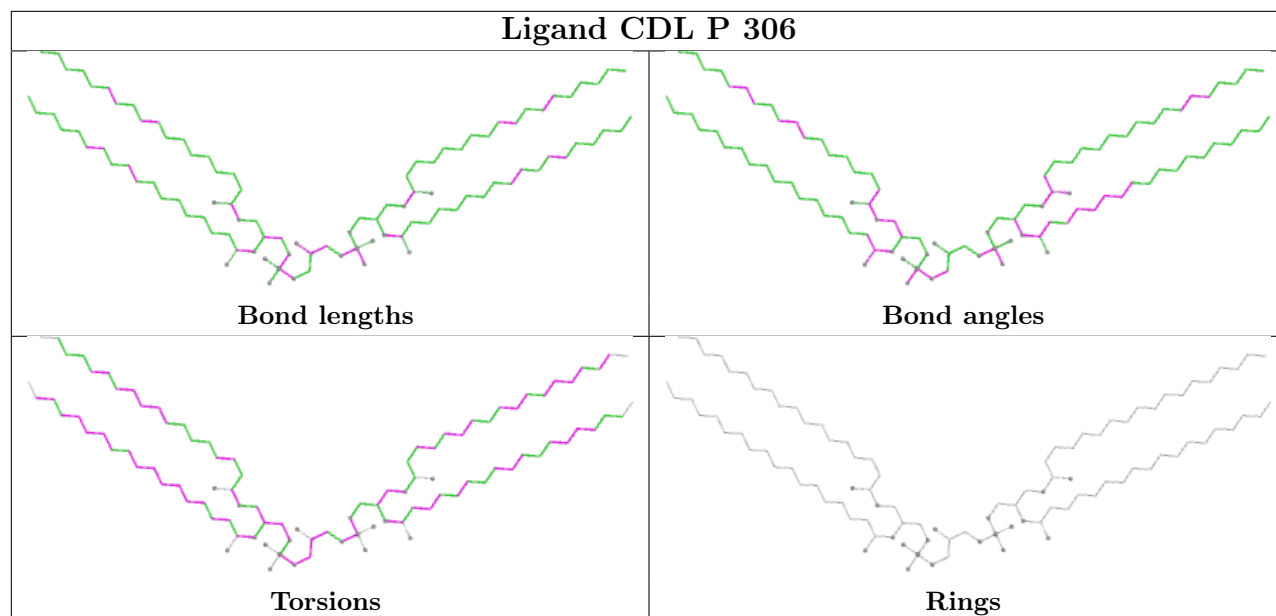
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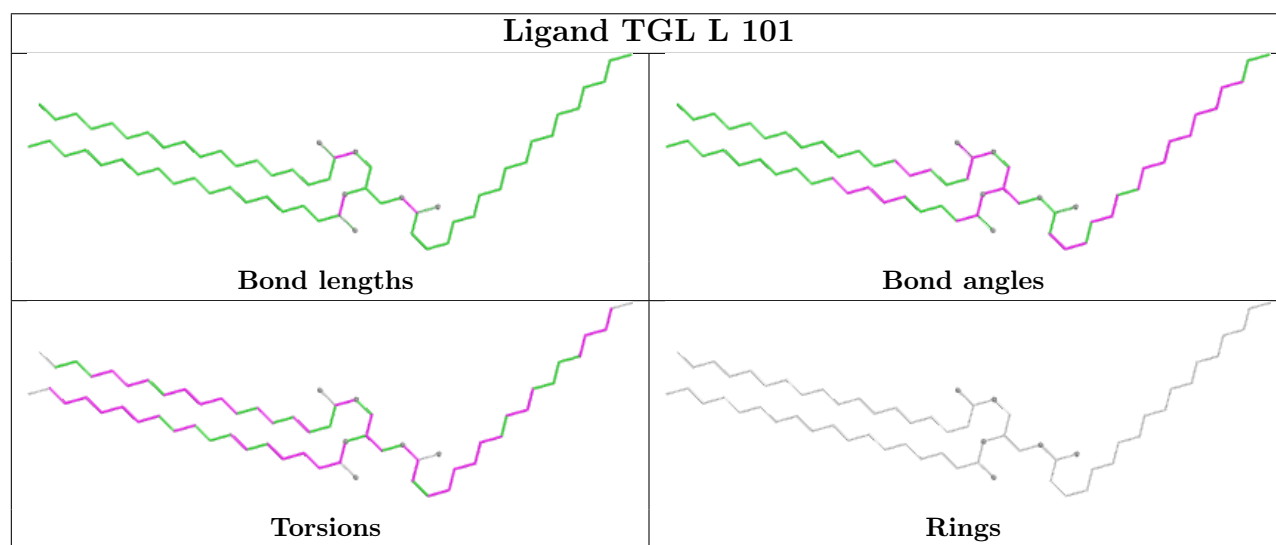
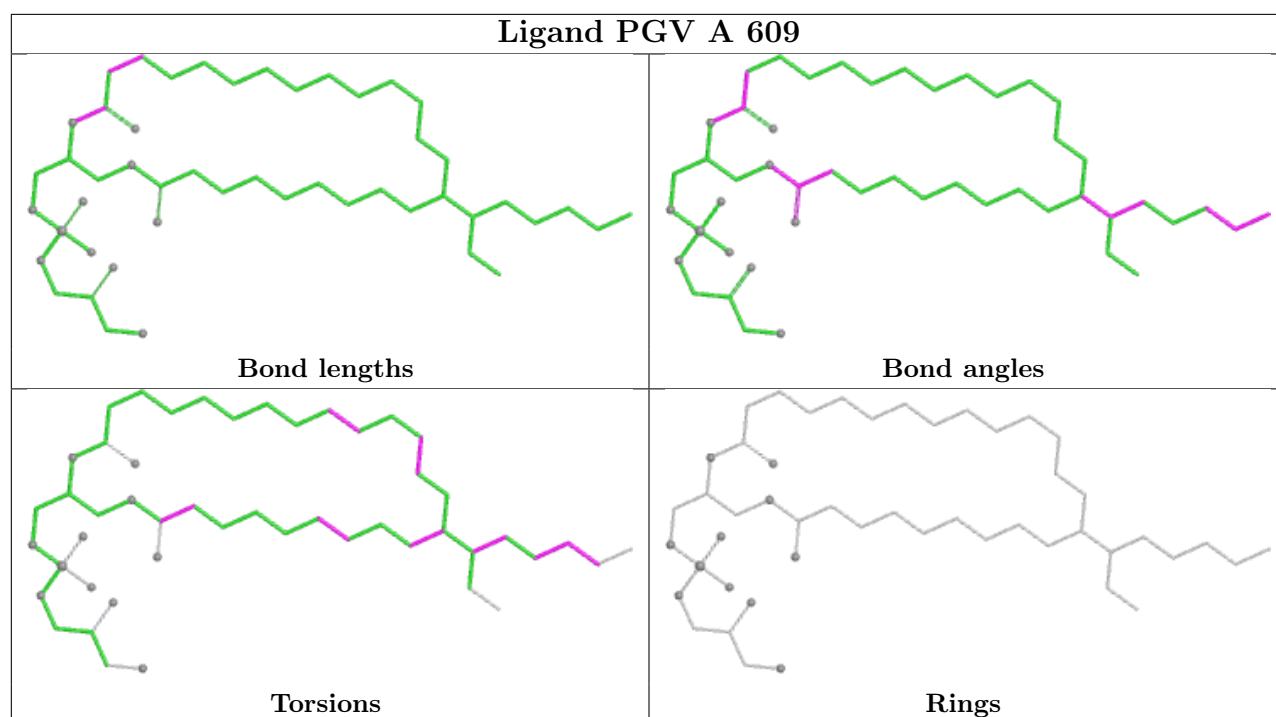


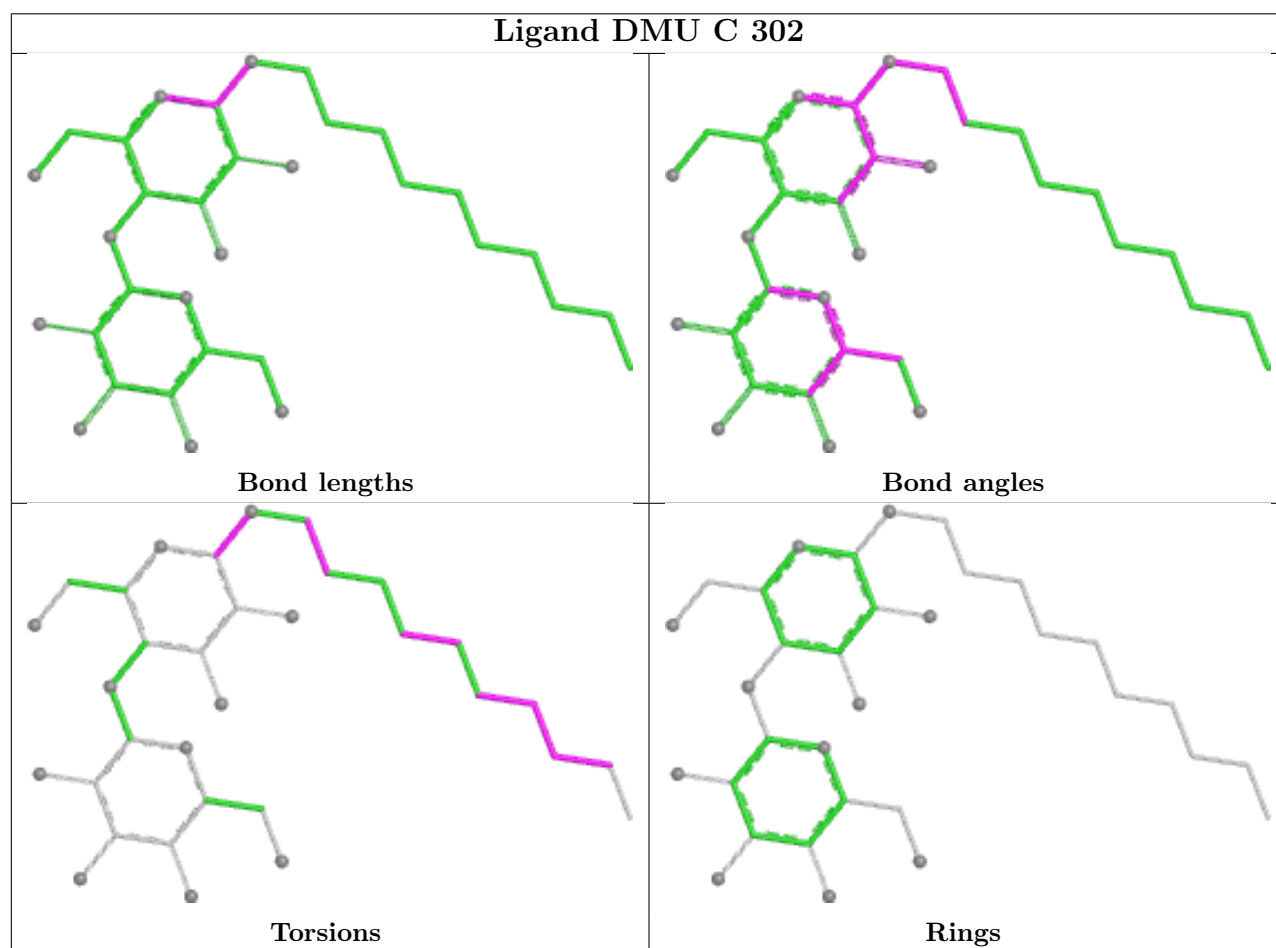
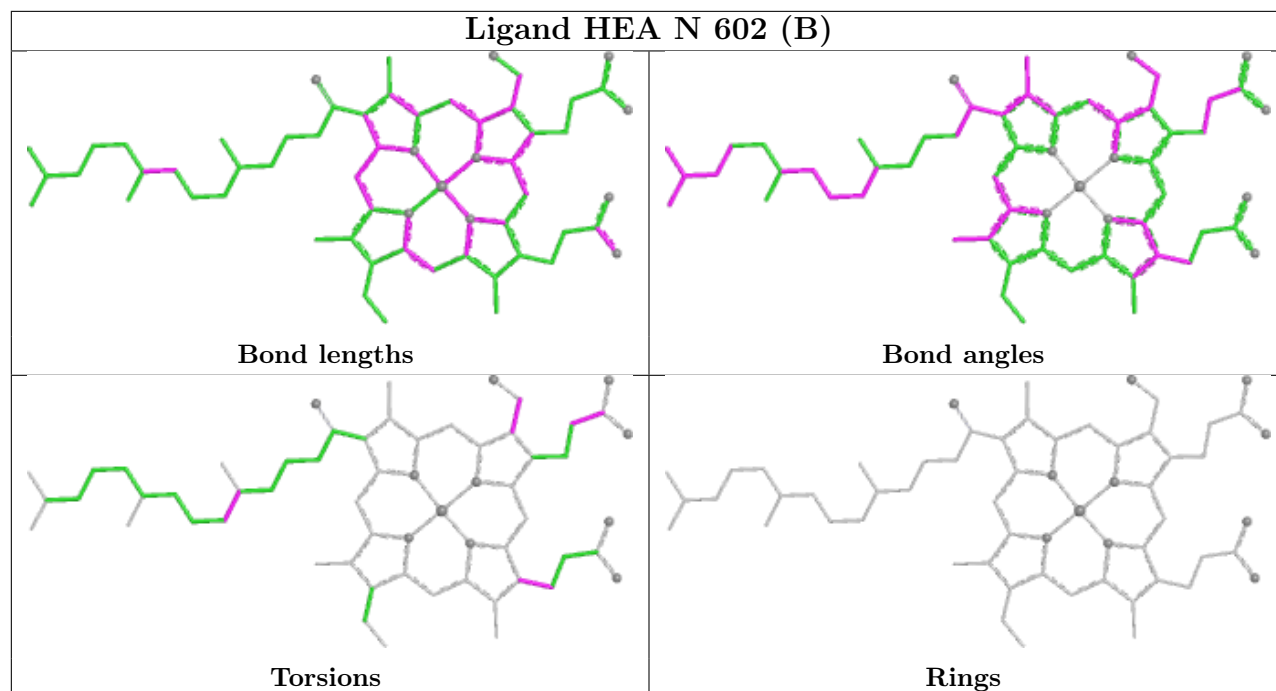
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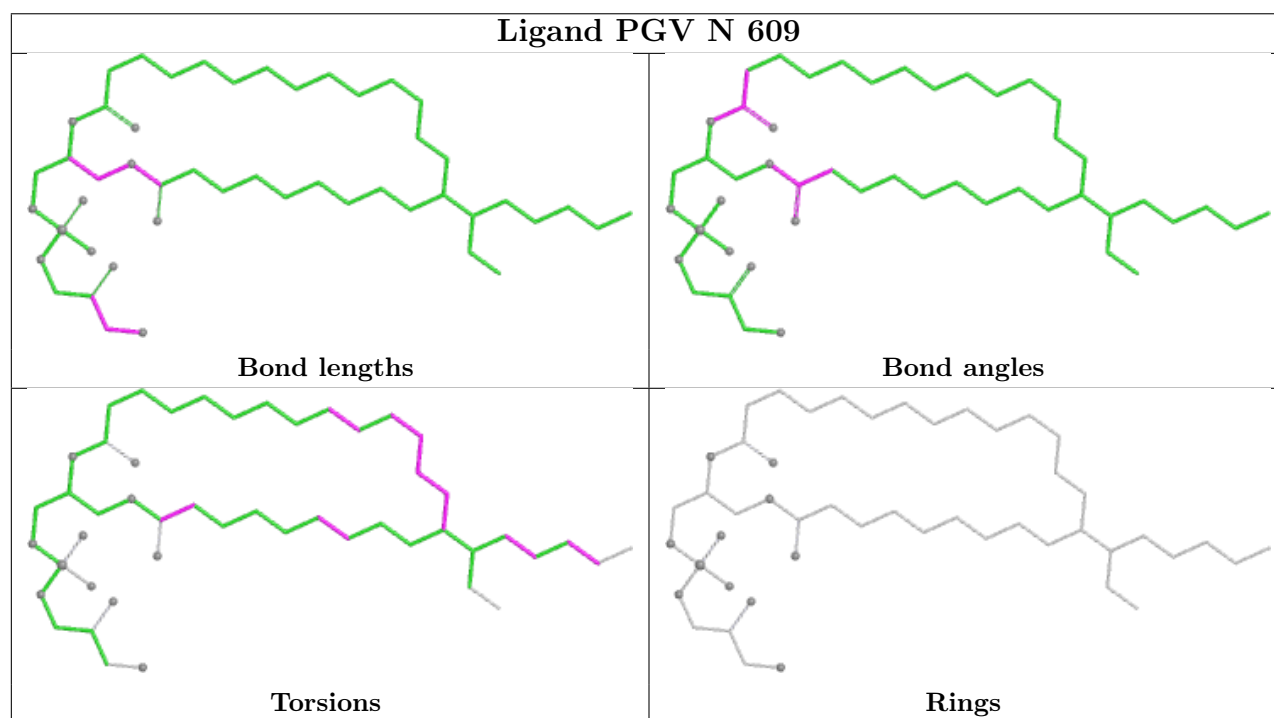
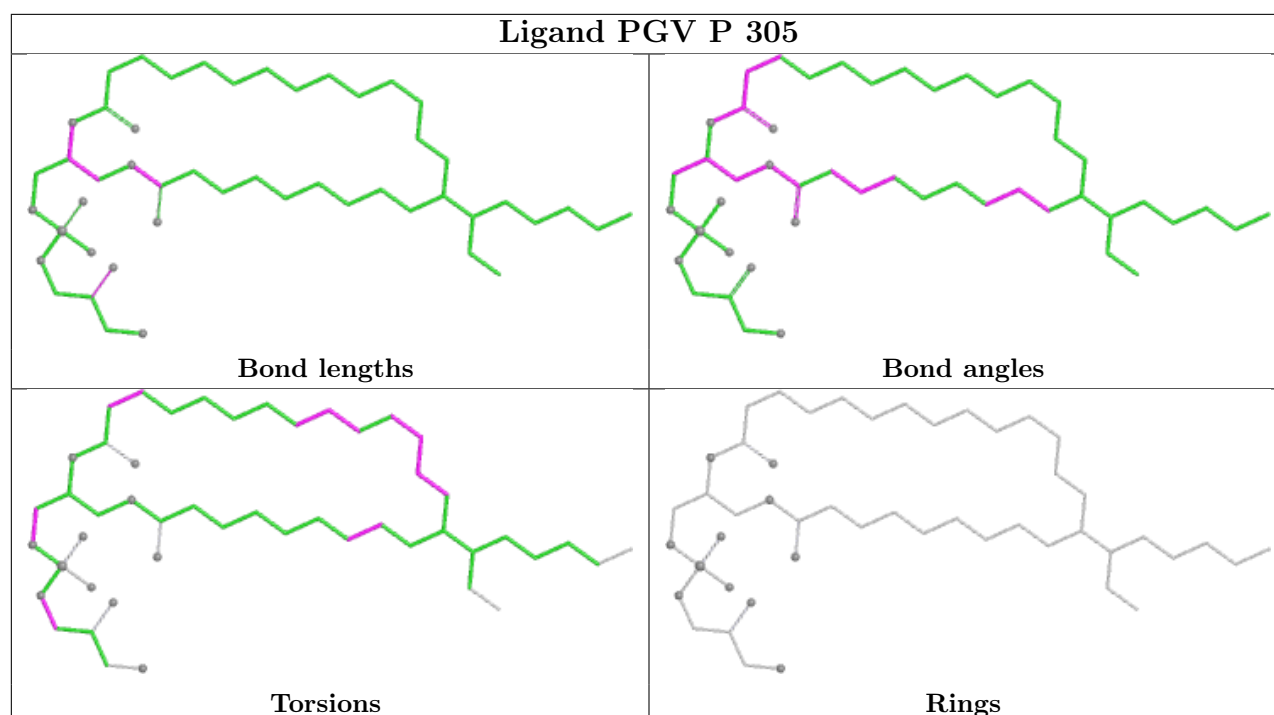


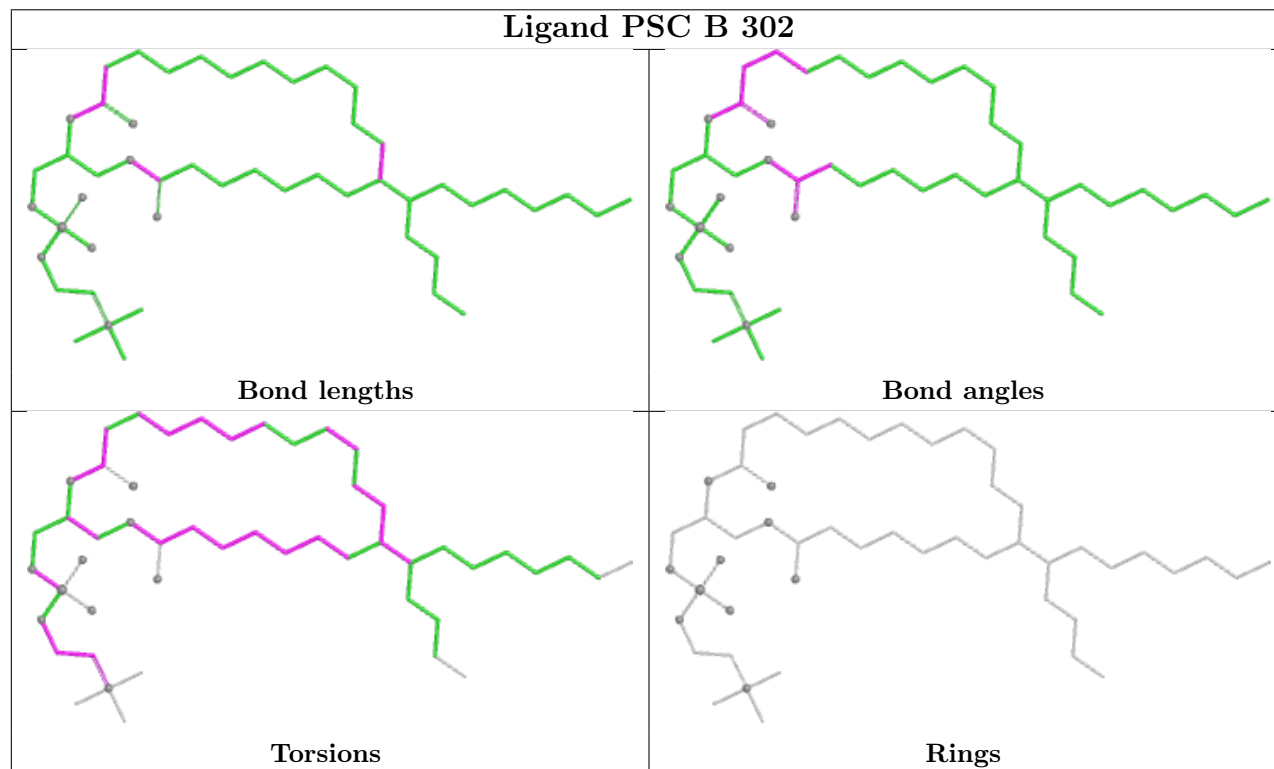
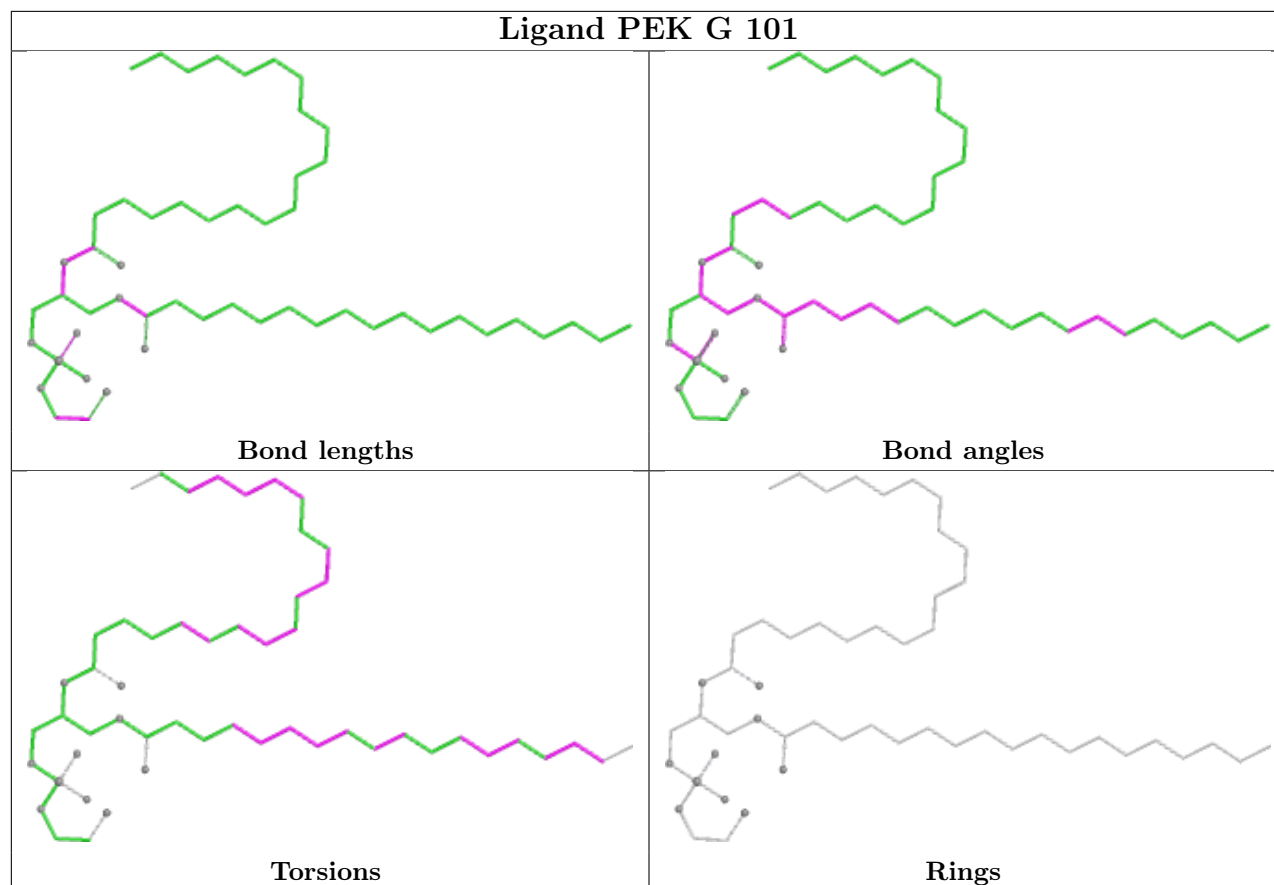


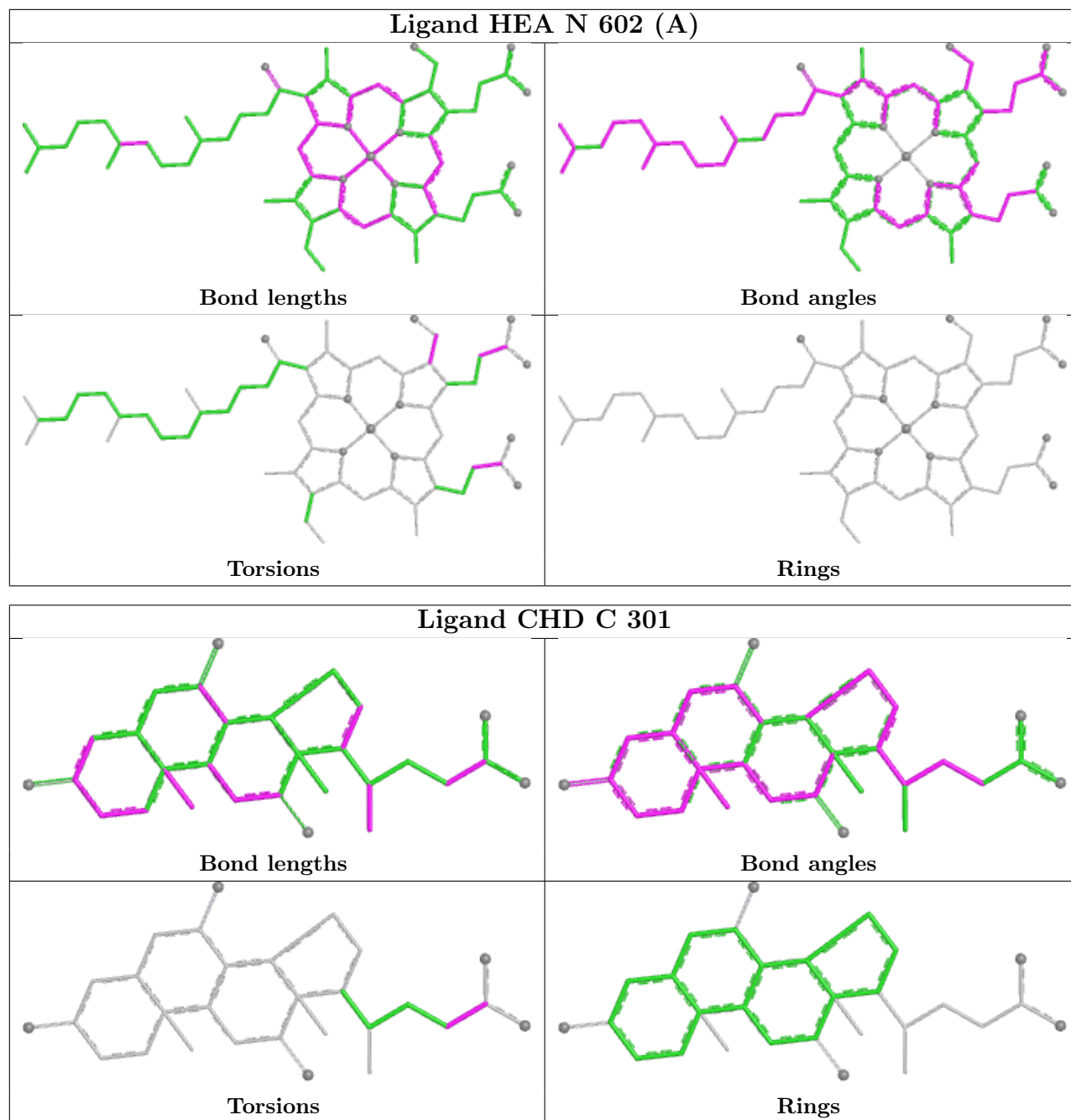


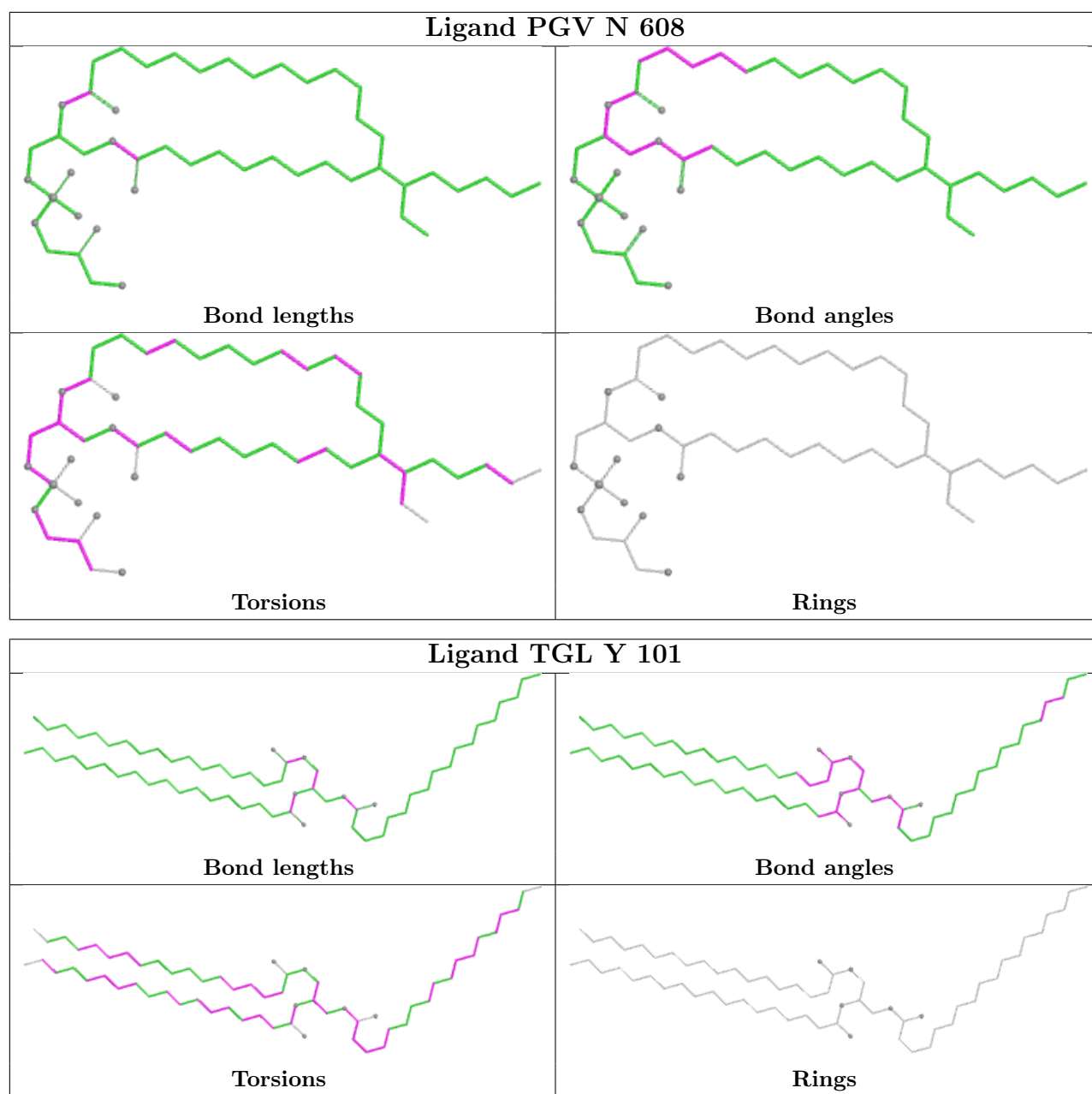




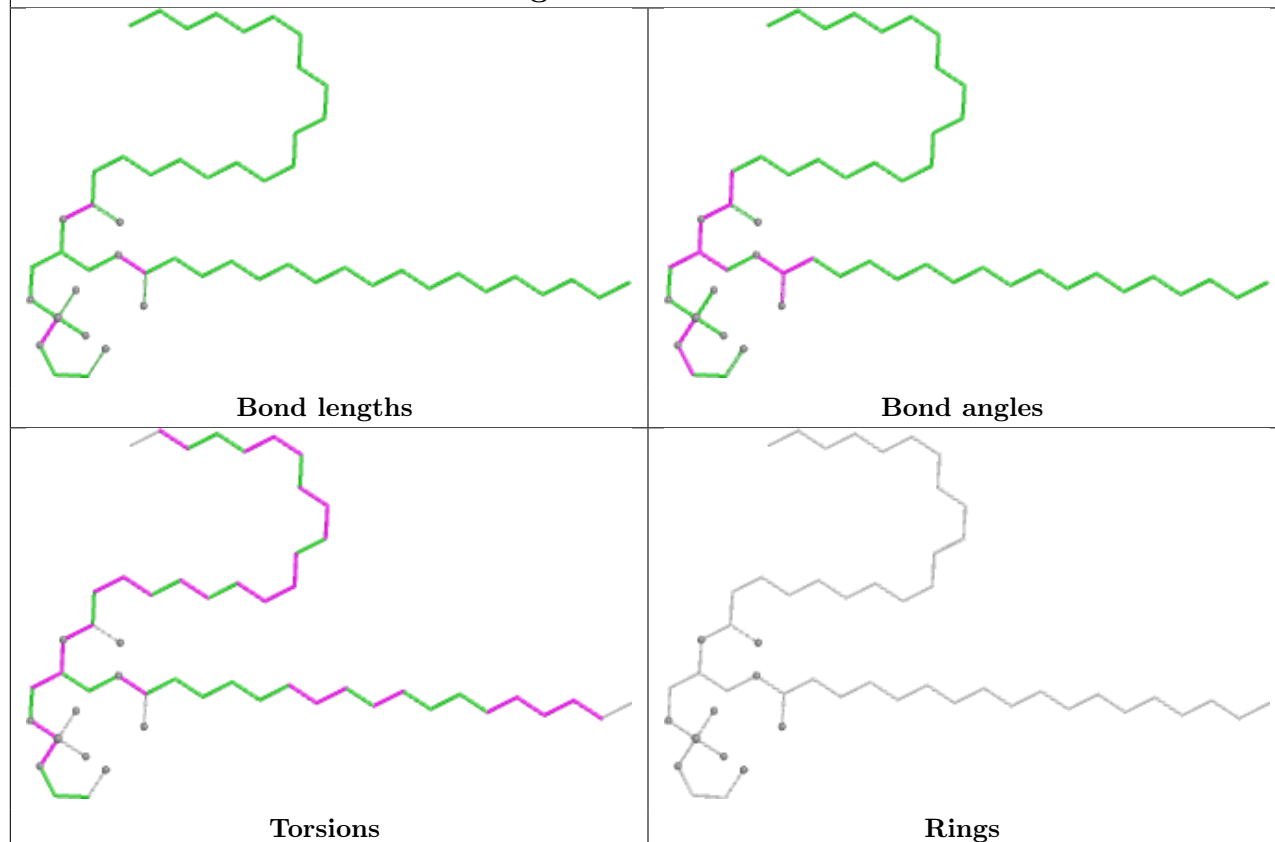




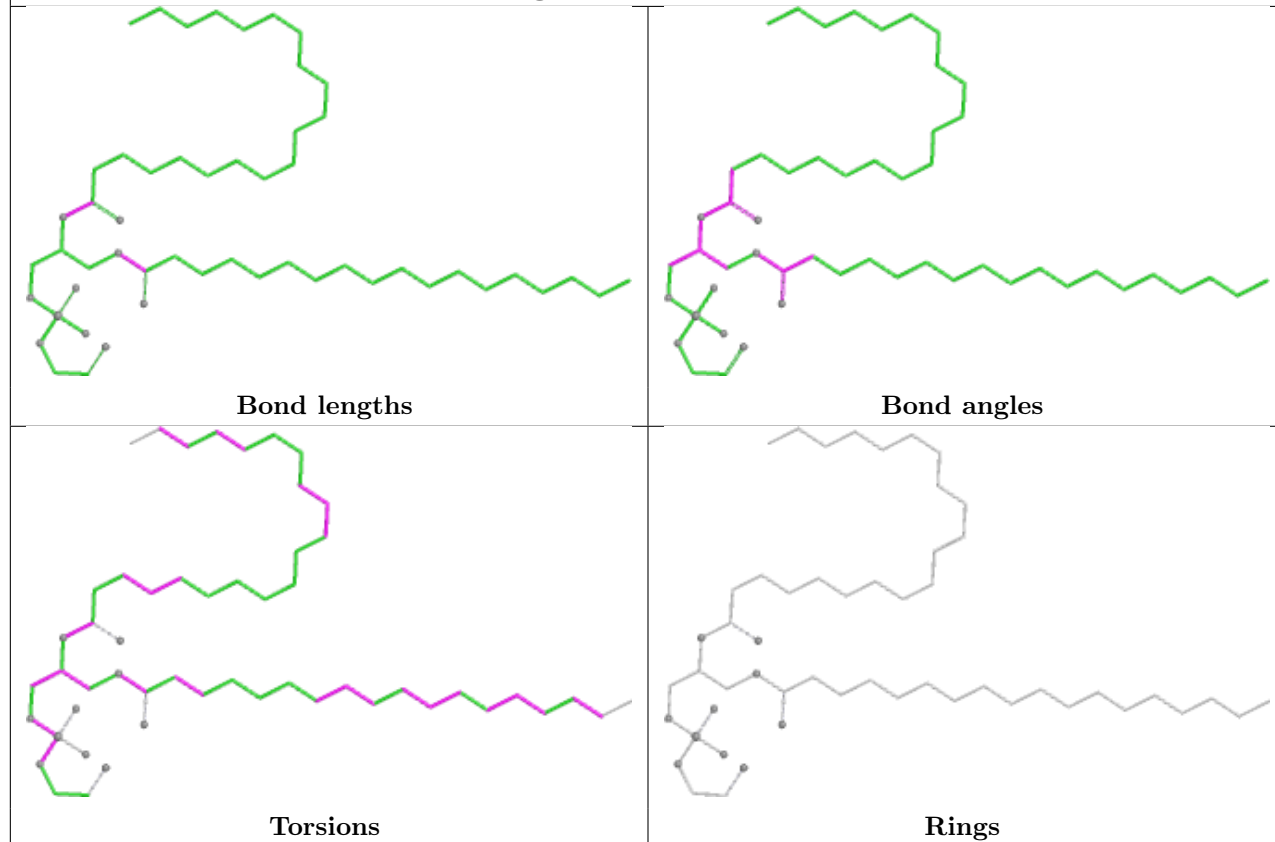


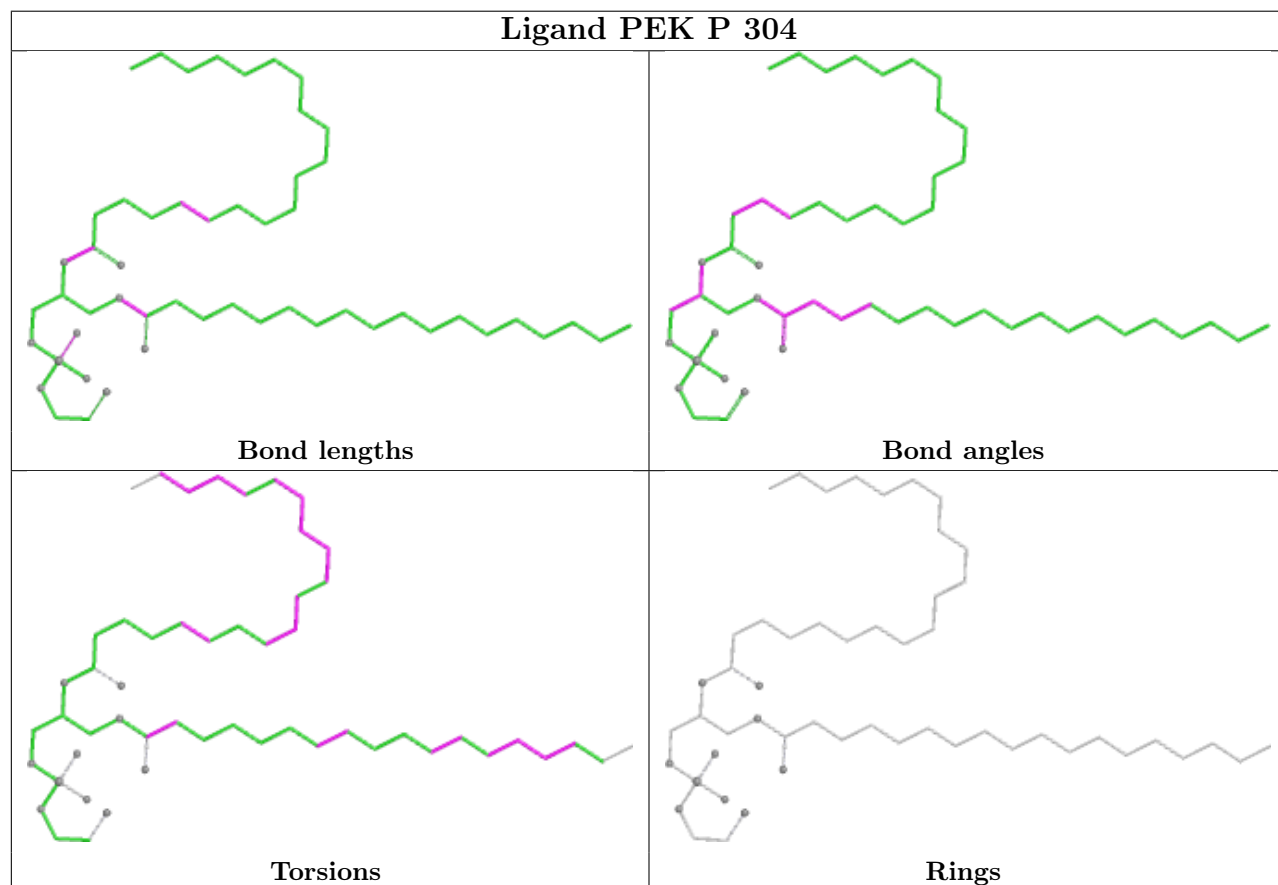
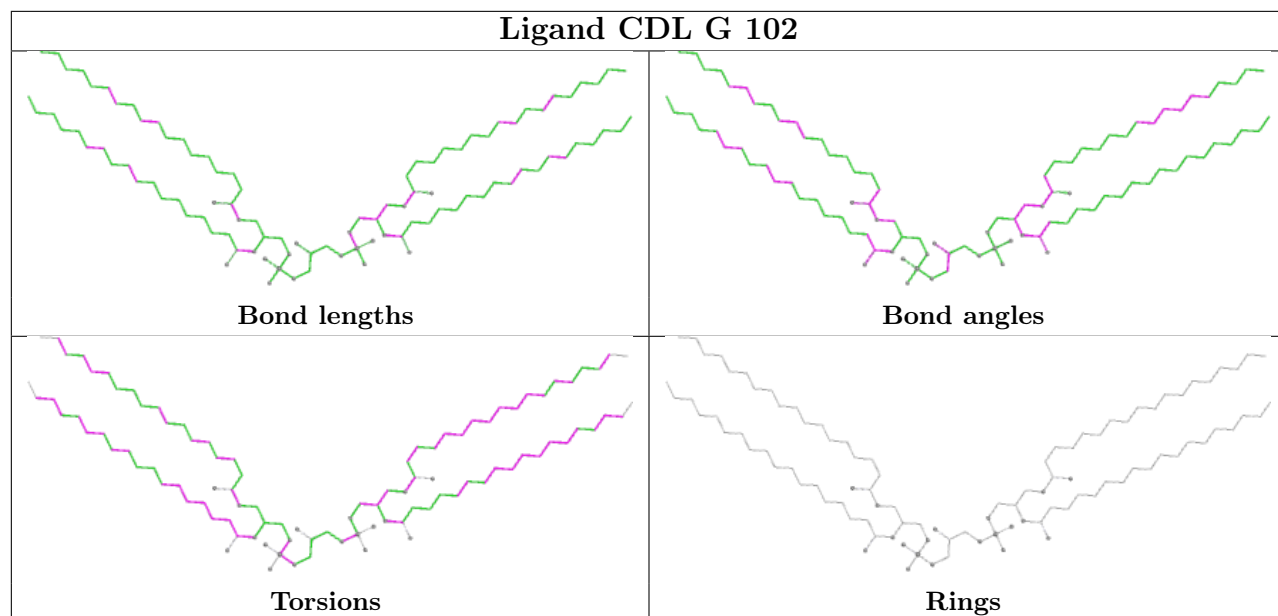


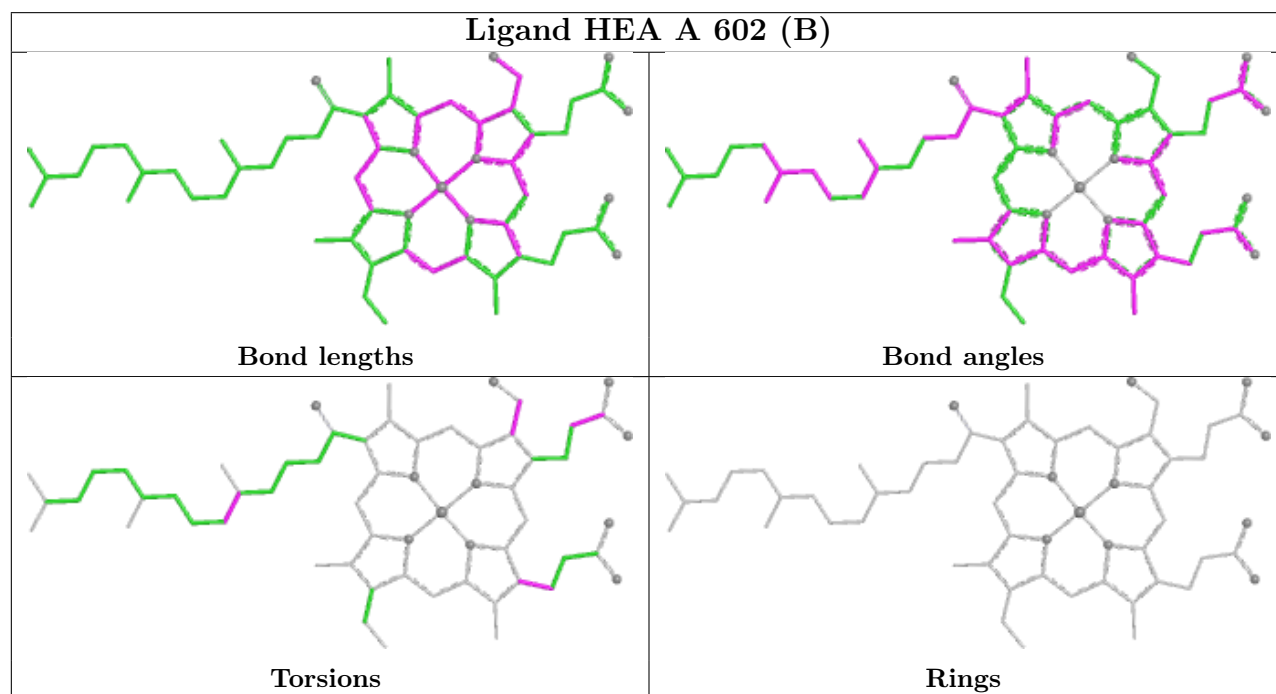
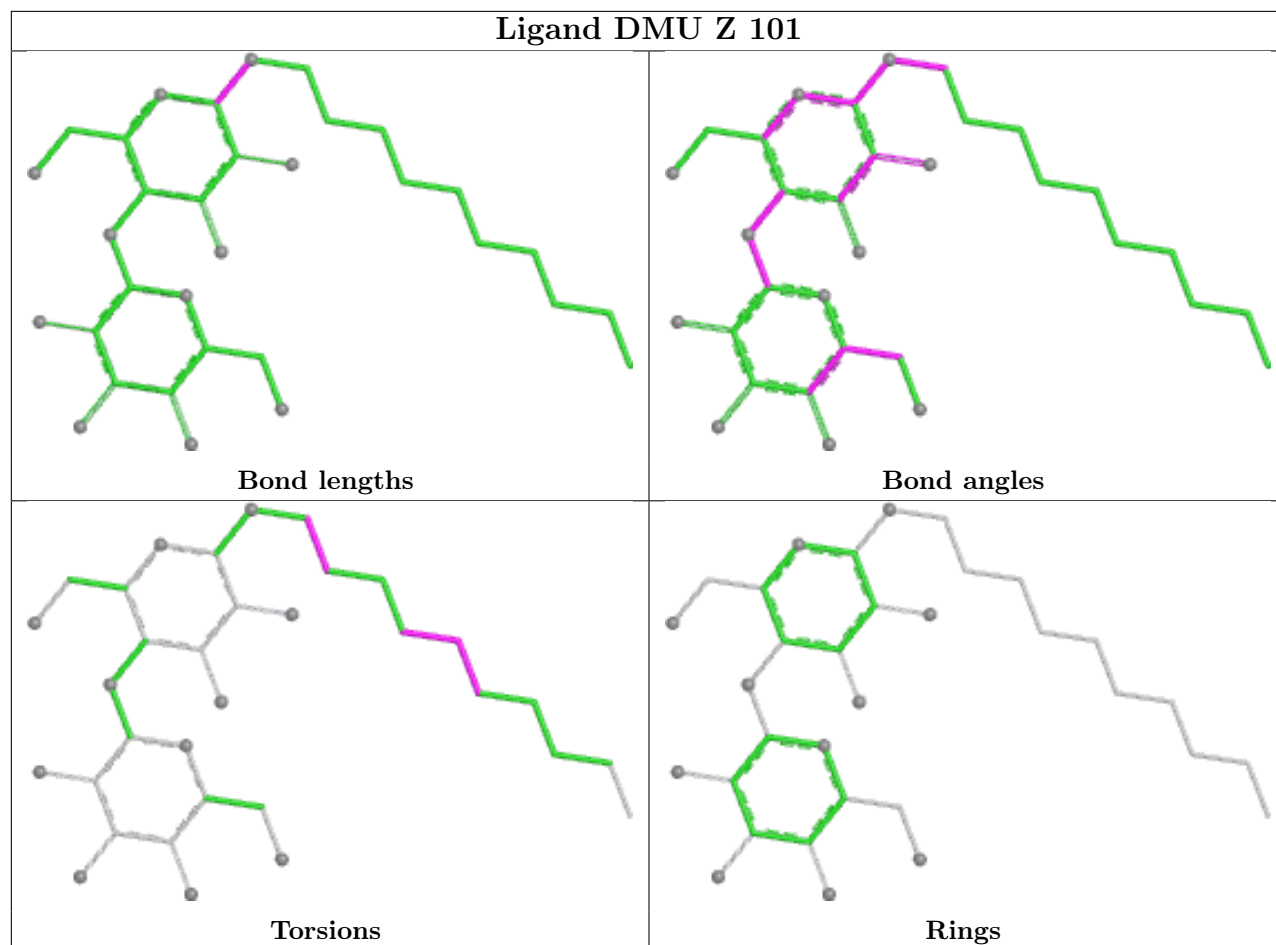
Ligand PEK C 307



Ligand PEK P 309







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%)	-0.59	3 (0%) 85 88	11, 25, 33, 69	18 (3%)
1	N	513/514 (99%)	-0.38	5 (0%) 79 83	13, 30, 40, 72	18 (3%)
2	B	226/227 (99%)	0.05	13 (5%) 29 31	20, 33, 53, 77	4 (1%)
2	O	226/227 (99%)	0.20	11 (4%) 35 38	20, 40, 65, 85	5 (2%)
3	C	259/261 (99%)	-0.37	4 (1%) 72 76	12, 29, 41, 79	7 (2%)
3	P	259/261 (99%)	-0.30	5 (1%) 66 70	12, 30, 42, 65	9 (3%)
4	D	144/147 (97%)	-0.03	4 (2%) 55 58	17, 34, 55, 77	1 (0%)
4	Q	144/147 (97%)	0.73	11 (7%) 20 21	19, 48, 74, 125	2 (1%)
5	E	105/109 (96%)	-0.15	2 (1%) 66 70	26, 33, 57, 116	0
5	R	105/109 (96%)	0.09	3 (2%) 53 57	32, 41, 60, 113	0
6	F	98/98 (100%)	0.38	9 (9%) 14 14	17, 37, 93, 130	3 (3%)
6	S	98/98 (100%)	0.64	9 (9%) 14 14	17, 37, 91, 133	2 (2%)
7	G	83/85 (97%)	0.97	19 (22%) 2 2	17, 38, 109, 137	1 (1%)
7	T	83/85 (97%)	1.12	21 (25%) 1 1	27, 40, 97, 133	0
8	H	79/85 (92%)	0.59	9 (11%) 10 9	30, 41, 90, 95	0
8	U	79/85 (92%)	0.68	13 (16%) 4 4	35, 45, 96, 124	0
9	I	72/73 (98%)	0.88	12 (16%) 4 4	31, 44, 75, 89	0
9	V	72/73 (98%)	0.87	8 (11%) 10 10	31, 53, 77, 99	0
10	J	58/59 (98%)	0.37	2 (3%) 48 52	29, 39, 65, 120	0
10	W	58/59 (98%)	0.41	2 (3%) 48 52	33, 44, 70, 110	0
11	K	49/56 (87%)	0.45	3 (6%) 27 29	33, 40, 55, 60	0
11	X	49/56 (87%)	0.74	5 (10%) 12 12	23, 50, 70, 78	1 (2%)
12	L	46/47 (97%)	-0.02	4 (8%) 16 16	26, 32, 50, 85	0
12	Y	46/47 (97%)	0.36	2 (4%) 40 43	23, 41, 58, 100	1 (2%)

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	M	43/46 (93%)	0.10	4 (9%) 14 14	27, 31, 66, 101	0
13	Z	43/46 (93%)	0.68	3 (6%) 22 24	39, 44, 76, 115	0
All	All	3550/3614 (98%)	0.04	186 (5%) 33 35	11, 34, 66, 137	72 (2%)

The worst 5 of 186 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
6	S	1	ALA	12.5
6	S	97	ALA	8.6
7	T	9	GLY	8.1
7	G	36	TRP	7.7
7	G	3	ALA	7.3

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
9	SAC	V	1	9/10	0.20	0.25	113,119,130,131	0
7	TPO	G	11	11/12	0.53	0.29	91,100,118,123	0
7	TPO	T	11	11/12	0.57	0.29	104,118,157,157	0
9	SAC	I	1	9/10	0.71	0.20	64,73,78,78	0
1	FME	A	1	10/11	0.93	0.12	36,45,72,87	0
1	FME	N	1	10/11	0.94	0.12	41,50,72,76	0
2	FME	B	1	10/11	0.97	0.08	29,30,41,66	0
2	FME	O	1	10/11	0.97	0.09	37,39,48,60	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
21	EDO	D	204	4/4	0.70	0.30	59,71,80,85	0
25	DMU	C	309	33/33	0.73	0.17	55,75,100,104	0
28	PEK	T	102	53/53	0.75	0.24	52,89,154,154	0
25	DMU	P	310	33/33	0.76	0.18	59,76,103,104	0
28	PEK	C	307	53/53	0.76	0.25	47,77,139,164	0
28	PEK	G	104	53/53	0.76	0.24	48,77,135,152	0
21	EDO	C	317	4/4	0.76	0.21	59,74,76,85	0
23	PSC	B	302	52/52	0.77	0.24	40,78,152,157	0
24	CHD	P	307	29/29	0.77	0.18	53,67,73,77	0
27	CDL	T	103	100/100	0.77	0.24	52,89,137,160	0
27	CDL	G	102	100/100	0.78	0.24	53,89,128,159	0
27	CDL	P	306	100/100	0.79	0.26	39,84,113,122	0
21	EDO	A	611	4/4	0.79	0.21	46,60,61,62	0
21	EDO	H	101	4/4	0.79	0.23	51,57,59,66	0
20	PGV	C	308	51/51	0.79	0.23	44,78,123,132	0
28	PEK	P	309	53/53	0.79	0.23	42,74,121,135	0
24	CHD	C	306	29/29	0.79	0.17	49,64,71,73	0
19	TGL	N	610	63/63	0.80	0.24	48,74,101,115	0
20	PGV	P	302	51/51	0.80	0.22	48,76,122,137	0
21	EDO	A	620	4/4	0.81	0.17	45,45,48,65	0
21	EDO	Q	204	4/4	0.81	0.18	64,68,71,77	0
19	TGL	Y	101	63/63	0.81	0.24	41,74,110,136	0
23	PSC	O	302	52/52	0.81	0.25	42,78,156,162	0
19	TGL	A	608	63/63	0.81	0.23	40,74,96,112	0
19	TGL	Q	201	63/63	0.82	0.24	48,73,95,107	0
25	DMU	P	308	33/33	0.82	0.14	45,74,113,121	0
20	PGV	N	608	51/51	0.82	0.21	47,78,115,127	0
21	EDO	Q	203	4/4	0.82	0.23	52,54,57,67	0
19	TGL	D	201	63/63	0.82	0.23	33,61,85,90	0
25	DMU	P	311	33/33	0.83	0.16	62,82,105,108	0
20	PGV	M	101	51/51	0.83	0.21	31,63,99,110	0
21	EDO	C	315	4/4	0.84	0.16	35,45,47,50	0
21	EDO	S	103	4/4	0.84	0.20	61,68,69,70	0
27	CDL	C	305	100/100	0.84	0.20	37,77,110,115	0
19	TGL	L	101	63/63	0.84	0.22	32,59,93,112	0
21	EDO	C	318	4/4	0.84	0.15	74,101,105,116	0
25	DMU	C	310	33/33	0.85	0.16	53,80,105,115	0
21	EDO	V	101	4/4	0.85	0.20	63,67,68,70	0
25	DMU	C	302	33/33	0.86	0.14	32,72,98,105	0
21	EDO	A	613	4/4	0.86	0.22	59,63,69,80	0
21	EDO	A	616	4/4	0.87	0.19	47,50,53,66	0
21	EDO	O	305	4/4	0.87	0.18	72,74,75,82	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
21	EDO	C	314	4/4	0.87	0.16	55,65,65,66	0
21	EDO	D	206	4/4	0.87	0.23	49,59,60,73	0
21	EDO	S	104	4/4	0.88	0.12	39,54,65,66	0
21	EDO	C	316	4/4	0.88	0.16	61,63,71,83	0
21	EDO	R	201	4/4	0.88	0.21	56,62,64,83	0
21	EDO	R	204	4/4	0.88	0.12	52,55,58,60	0
21	EDO	A	615	4/4	0.88	0.17	41,42,45,45	0
21	EDO	U	101	4/4	0.89	0.16	62,65,66,72	0
21	EDO	N	618	4/4	0.89	0.17	42,43,51,55	0
21	EDO	C	312	4/4	0.89	0.14	79,82,85,92	0
25	DMU	Z	101	33/33	0.89	0.12	47,56,73,77	0
21	EDO	E	203	4/4	0.90	0.13	48,52,54,60	0
21	EDO	R	205	4/4	0.90	0.17	42,49,53,53	0
21	EDO	G	105	4/4	0.90	0.19	51,62,74,87	0
21	EDO	A	619	4/4	0.90	0.14	55,58,58,60	0
21	EDO	A	618	4/4	0.90	0.14	43,51,65,69	0
26	UNX	C	303	1/1	0.90	0.14	30,30,30,30	0
21	EDO	R	202	4/4	0.90	0.15	49,58,63,63	0
21	EDO	N	611	4/4	0.91	0.15	65,65,66,69	0
25	DMU	M	102	33/33	0.91	0.12	39,47,56,66	0
21	EDO	D	202	4/4	0.91	0.13	39,45,50,65	0
21	EDO	N	619	4/4	0.91	0.14	45,46,55,56	0
21	EDO	A	617	4/4	0.91	0.16	44,66,70,75	0
21	EDO	E	205	4/4	0.92	0.14	52,55,58,67	0
21	EDO	A	612	4/4	0.92	0.16	31,32,37,74	0
21	EDO	N	621	4/4	0.92	0.16	40,45,52,57	0
21	EDO	N	614	4/4	0.92	0.17	58,61,62,72	0
21	EDO	P	312	4/4	0.92	0.15	32,33,45,52	0
26	UNX	P	303	1/1	0.92	0.20	29,29,29,29	0
21	EDO	P	315	4/4	0.92	0.15	46,54,57,72	0
21	EDO	C	319	4/4	0.93	0.18	32,42,46,58	0
21	EDO	C	313	4/4	0.93	0.12	56,60,64,81	0
21	EDO	N	612	4/4	0.93	0.18	30,37,45,57	0
21	EDO	D	203	4/4	0.93	0.12	54,54,66,72	0
21	EDO	F	104	4/4	0.93	0.12	47,48,53,55	0
21	EDO	A	614	4/4	0.93	0.11	19,25,27,39	0
21	EDO	E	201	4/4	0.94	0.10	41,41,47,56	0
18	AZI	N	607	3/3	0.94	0.15	32,32,34,38	0
21	EDO	P	314	4/4	0.94	0.13	48,56,63,65	0
21	EDO	O	304	4/4	0.94	0.11	61,62,65,76	0
24	CHD	C	301	29/29	0.95	0.08	27,31,34,38	0
21	EDO	N	615	4/4	0.95	0.12	33,33,39,43	0

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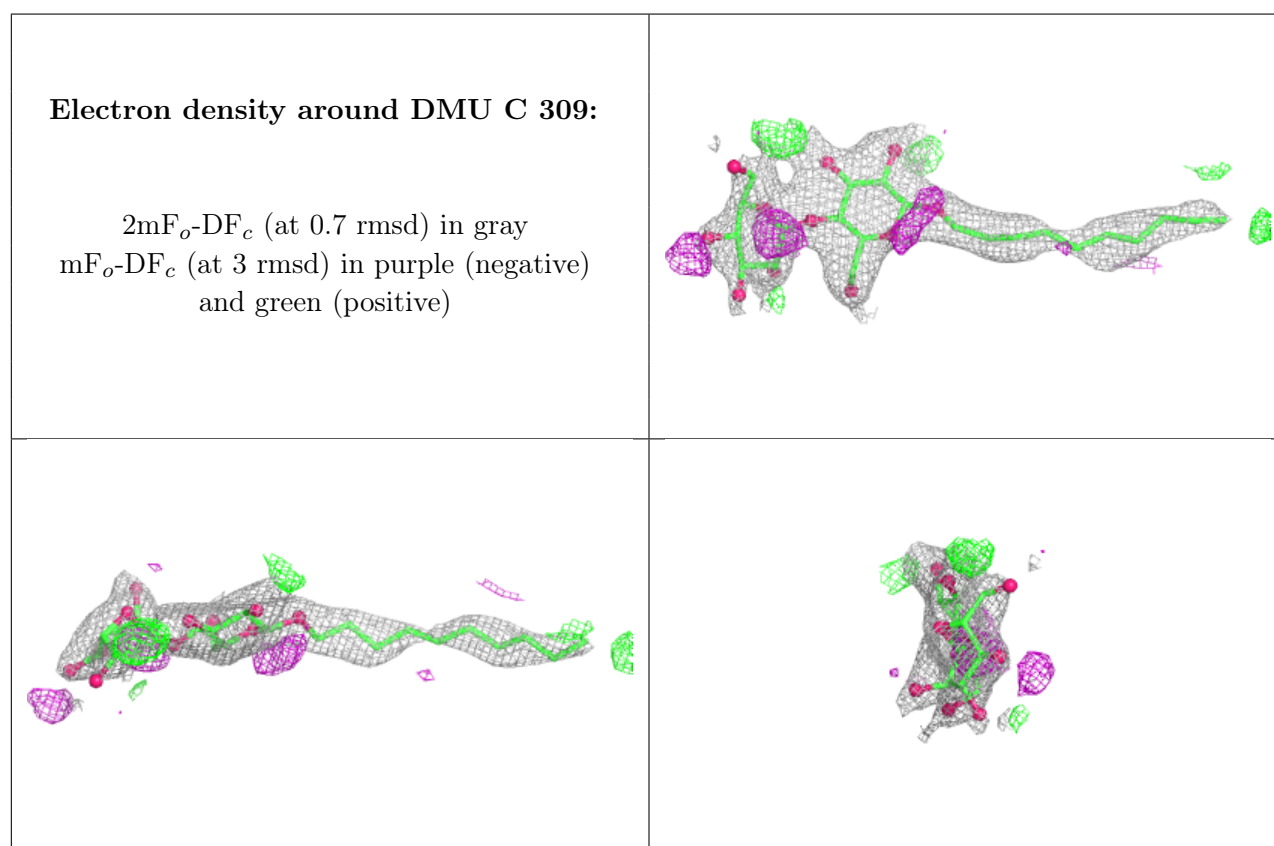
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
21	EDO	T	104	4/4	0.95	0.12	34,36,39,42	0
21	EDO	N	617	4/4	0.95	0.12	41,42,44,52	0
21	EDO	Q	202	4/4	0.95	0.13	58,58,64,64	0
28	PEK	P	304	53/53	0.95	0.14	29,47,85,94	0
21	EDO	N	613	4/4	0.95	0.13	36,44,45,49	0
21	EDO	C	311	4/4	0.95	0.10	38,39,39,41	0
20	PGV	N	609	51/51	0.96	0.10	27,35,62,66	0
21	EDO	E	204	4/4	0.96	0.08	37,39,42,46	0
16	MG	N	604	1/1	0.96	0.05	31,31,31,31	0
28	PEK	G	101	53/53	0.96	0.12	27,44,79,97	0
21	EDO	P	313	4/4	0.96	0.09	41,42,43,46	0
21	EDO	N	620	4/4	0.96	0.11	42,55,55,65	0
21	EDO	F	103	4/4	0.96	0.08	36,37,37,39	0
21	EDO	O	303	4/4	0.96	0.09	37,37,37,38	0
20	PGV	A	609	51/51	0.97	0.09	23,32,62,65	0
20	PGV	P	305	51/51	0.97	0.09	26,35,72,76	0
21	EDO	R	203	4/4	0.97	0.08	44,44,46,46	0
24	CHD	P	301	29/29	0.97	0.07	27,32,36,40	0
21	EDO	A	610	4/4	0.97	0.10	36,38,39,45	0
24	CHD	T	101	29/29	0.97	0.06	26,28,31,39	0
21	EDO	B	303	4/4	0.97	0.07	28,28,34,34	0
21	EDO	N	616	4/4	0.97	0.07	26,31,33,35	0
21	EDO	D	205	4/4	0.97	0.09	35,36,53,55	0
20	PGV	C	304	51/51	0.97	0.10	26,32,80,91	0
21	EDO	G	106	4/4	0.97	0.08	29,34,38,40	0
18	AZI	A	607	3/3	0.97	0.15	30,30,31,34	0
21	EDO	E	202	4/4	0.97	0.08	37,39,43,45	0
21	EDO	S	102	4/4	0.98	0.07	28,29,30,31	0
17	NA	A	605	1/1	0.98	0.04	28,28,28,28	0
18	AZI	A	606	3/3	0.98	0.06	28,28,28,28	0
14	HEA	A	602[B]	60/60	0.98	0.06	18,22,33,37	60
24	CHD	G	103	29/29	0.98	0.05	26,30,33,38	0
14	HEA	N	601	60/60	0.98	0.07	25,28,48,50	0
14	HEA	A	602[A]	60/60	0.98	0.06	18,22,25,27	60
14	HEA	A	601	60/60	0.99	0.06	19,23,45,45	0
14	HEA	N	602[A]	60/60	0.99	0.06	21,25,29,31	60
17	NA	N	605	1/1	0.99	0.03	35,35,35,35	0
14	HEA	N	602[B]	60/60	0.99	0.06	21,26,39,42	60
21	EDO	F	102	4/4	0.99	0.05	24,25,26,28	0
16	MG	A	604	1/1	0.99	0.03	24,24,24,24	0
18	AZI	N	606	3/3	0.99	0.07	33,33,34,35	0
15	CU	N	603	1/1	1.00	0.01	28,28,28,28	0

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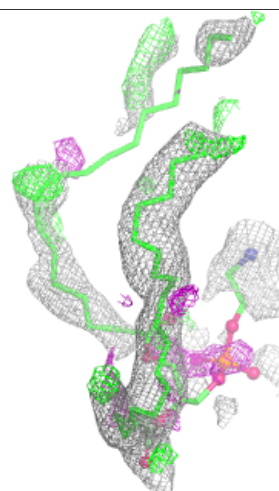
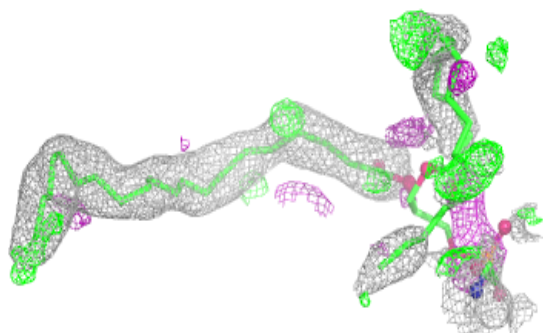
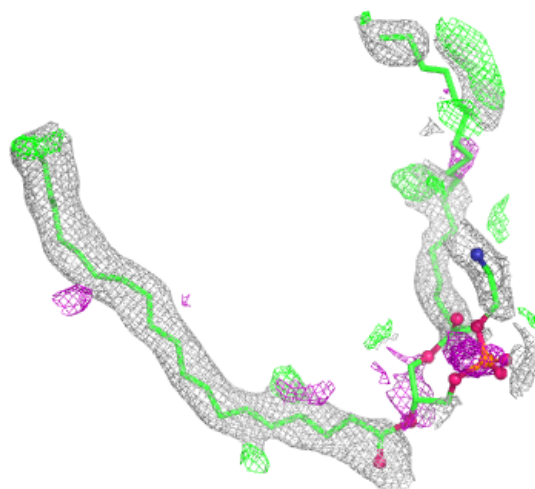
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
15	CU	A	603	1/1	1.00	0.02	25,25,25,25	0
22	CUA	B	301	2/2	1.00	0.02	25,25,25,26	0
22	CUA	O	301	2/2	1.00	0.02	31,31,31,32	0
29	ZN	F	101	1/1	1.00	0.02	30,30,30,30	0
29	ZN	S	101	1/1	1.00	0.03	32,32,32,32	0

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.



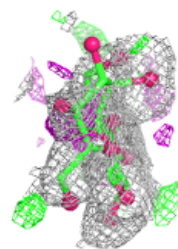
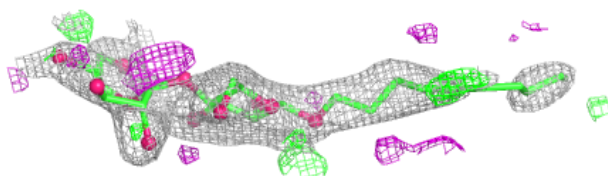
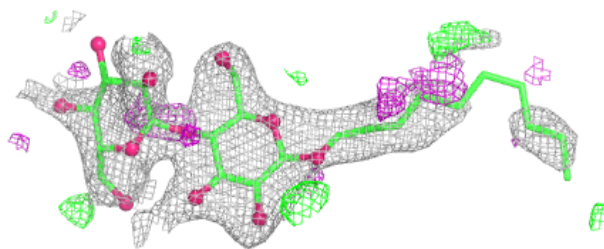
Electron density around PEK T 102:

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



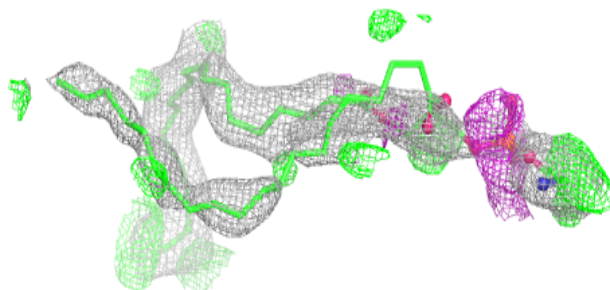
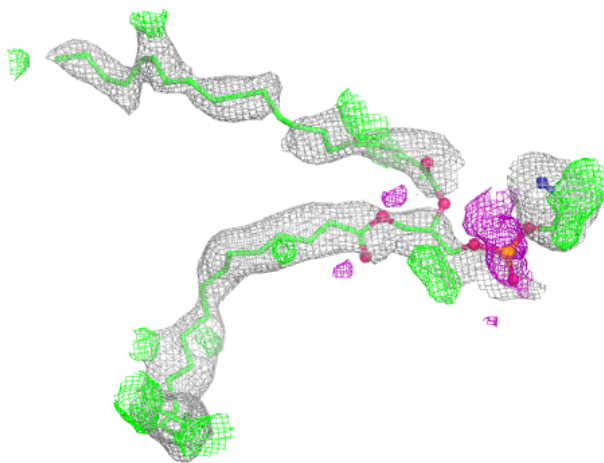
Electron density around DMU P 310:

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



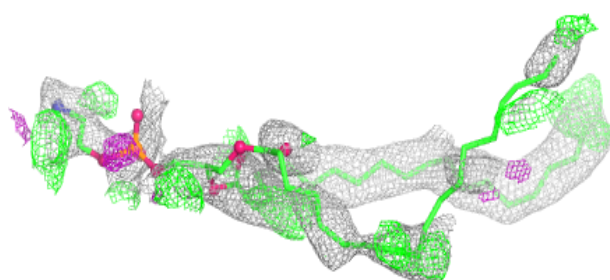
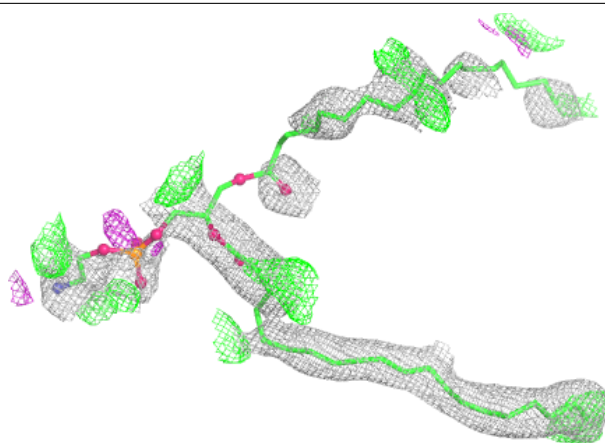
Electron density around PEK C 307:

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

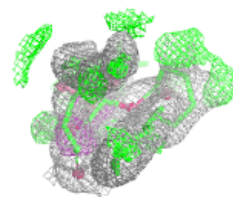
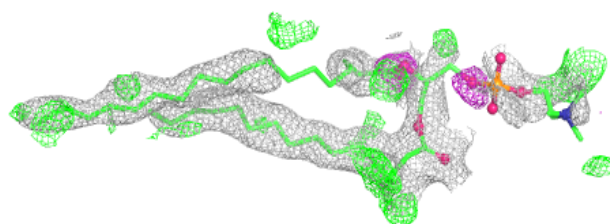
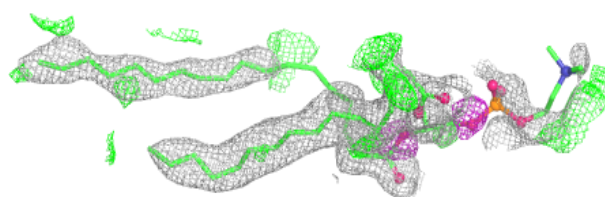


Electron density around PEK G 104:

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

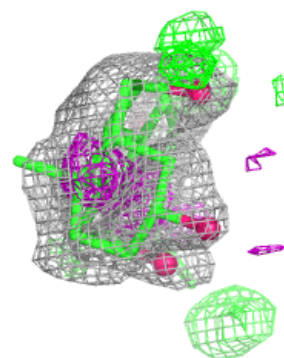
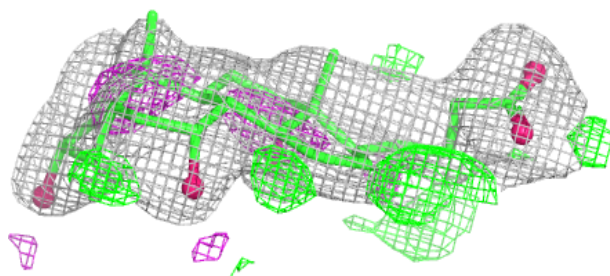
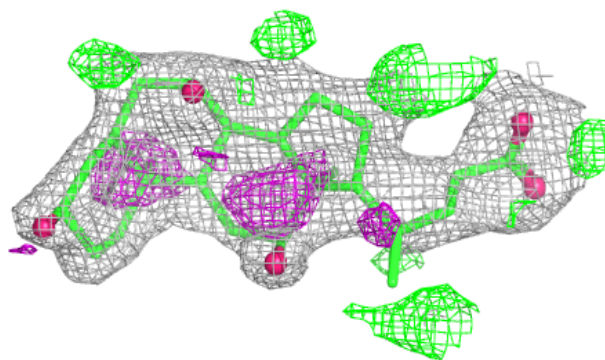
**Electron density around PSC B 302:**

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

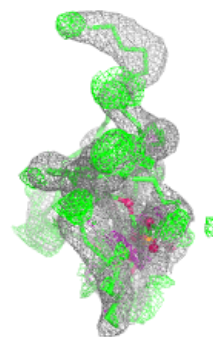
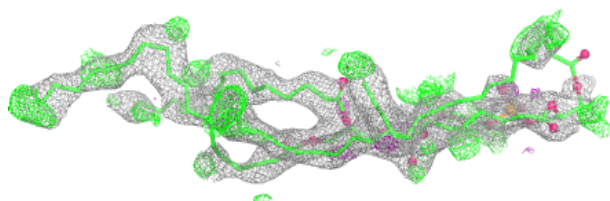
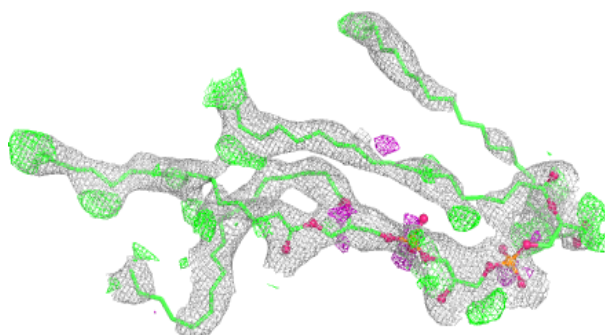


Electron density around CHD P 307:

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

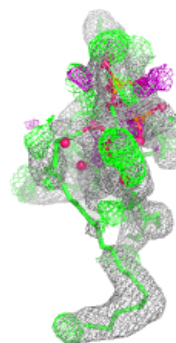
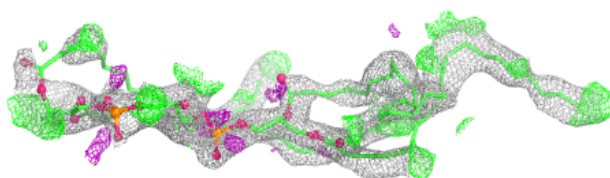
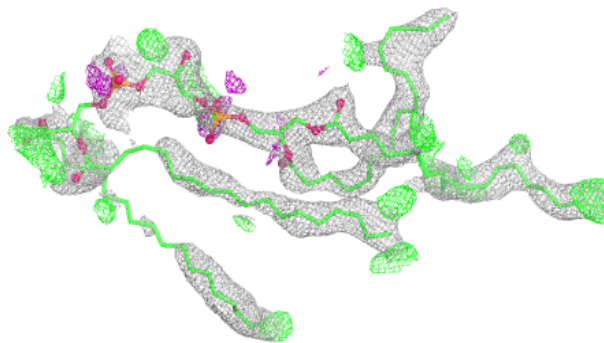
**Electron density around CDL T 103:**

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



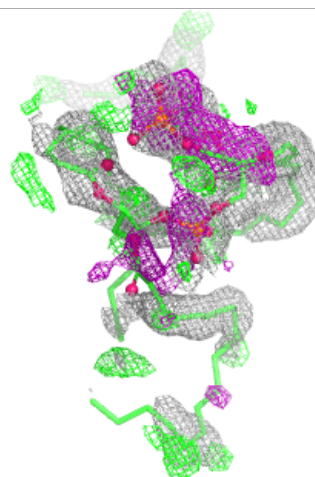
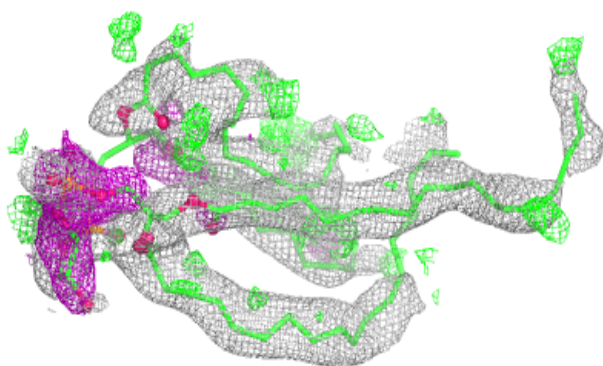
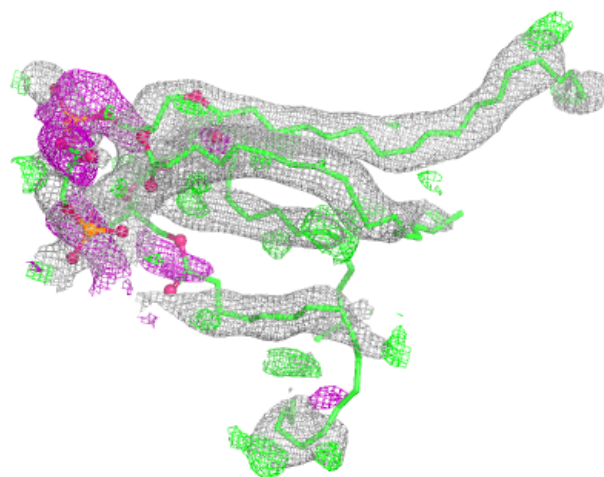
Electron density around CDL G 102:

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



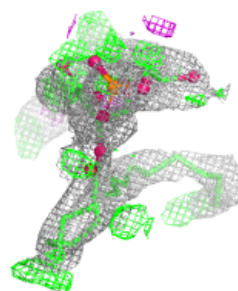
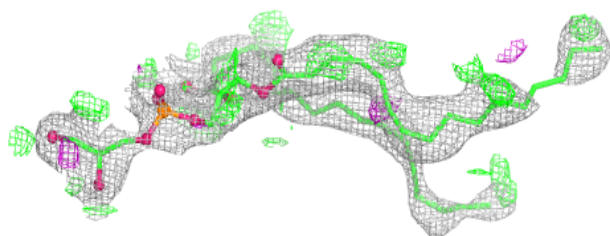
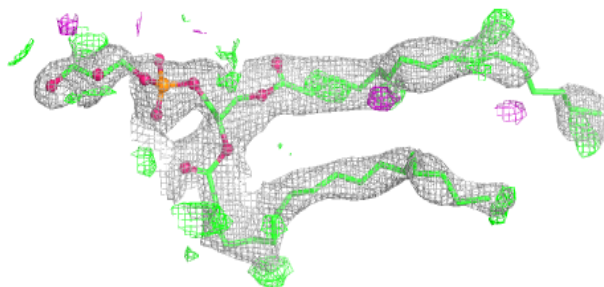
Electron density around CDL P 306:

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

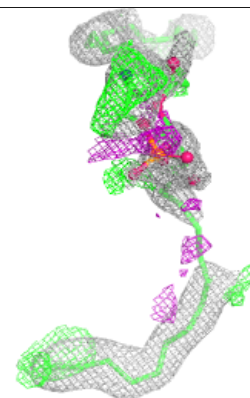
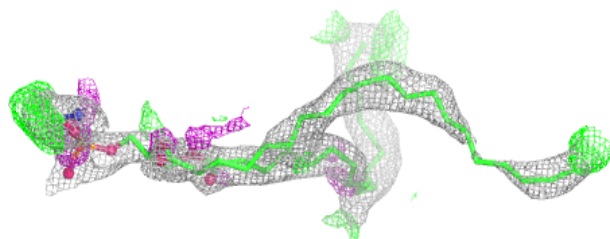
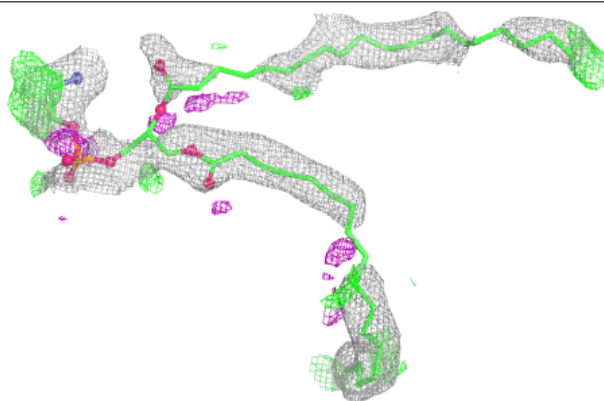


Electron density around PGV C 308:

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

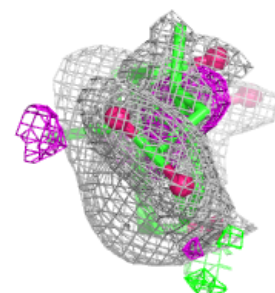
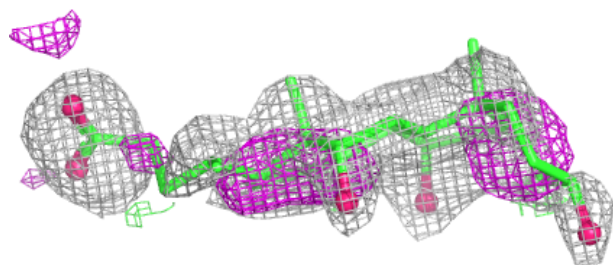
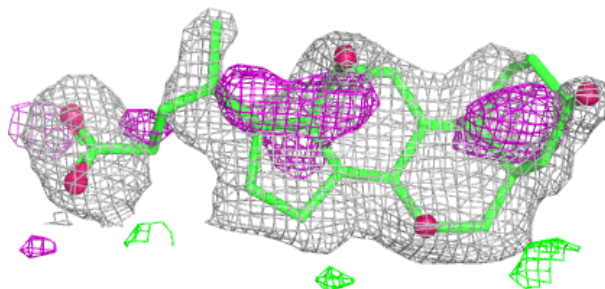
**Electron density around PEK P 309:**

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

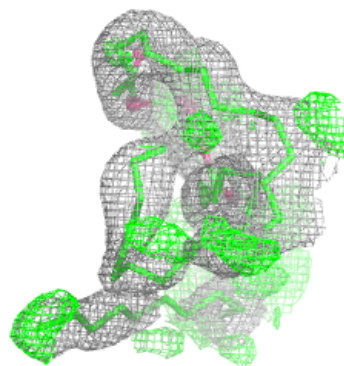
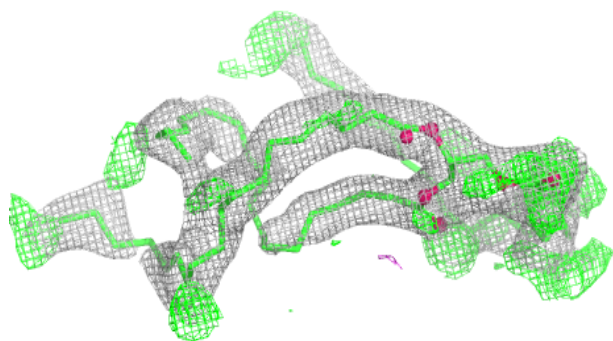
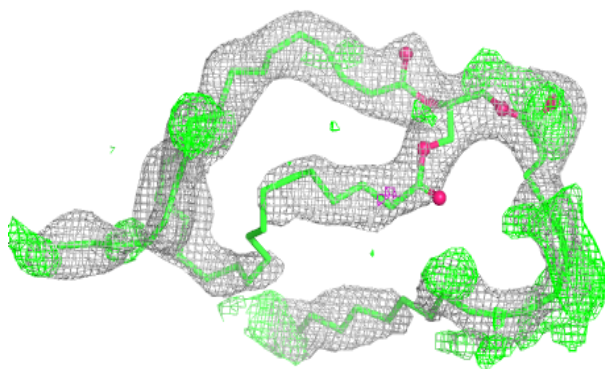


Electron density around CHD C 306:

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

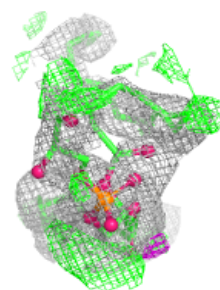
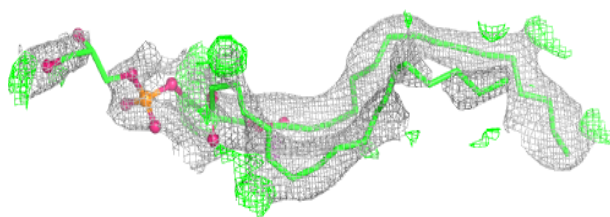
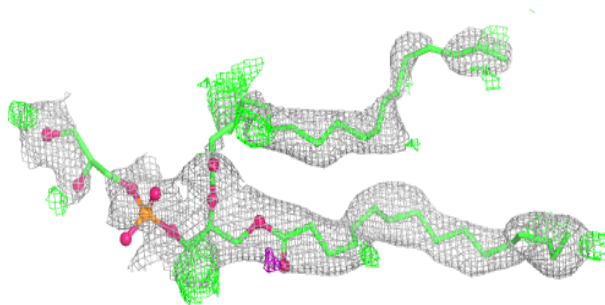
**Electron density around TGL N 610:**

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



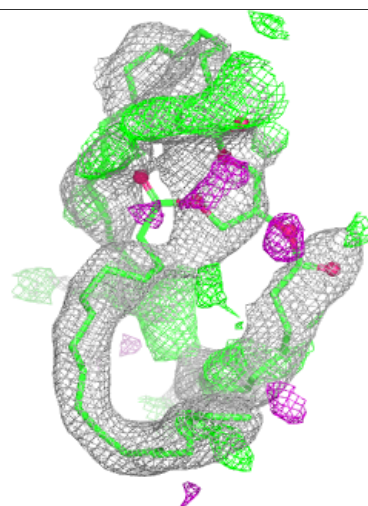
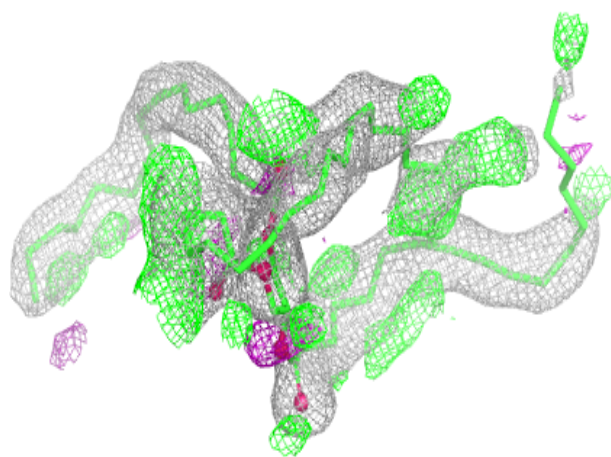
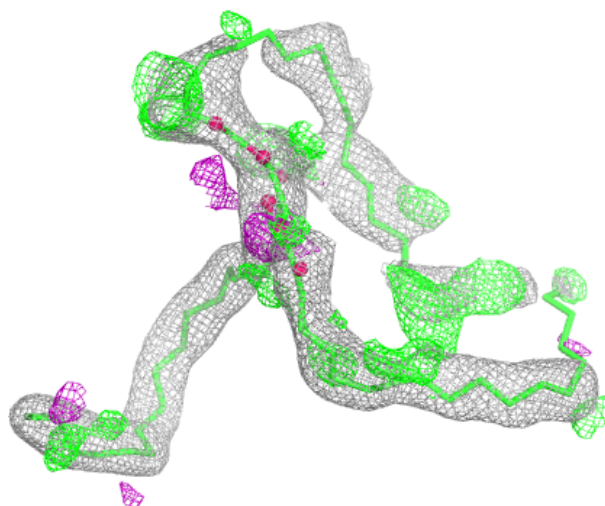
Electron density around PGV P 302:

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



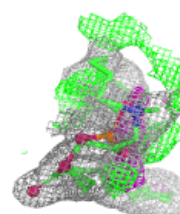
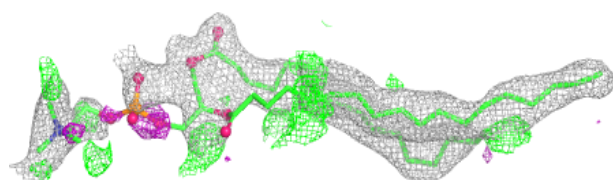
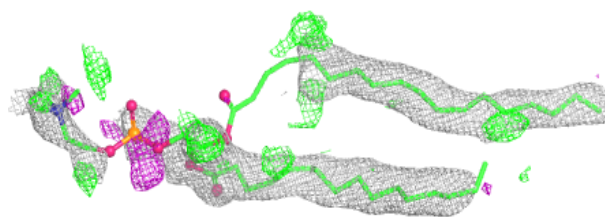
Electron density around TGL Y 101:

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

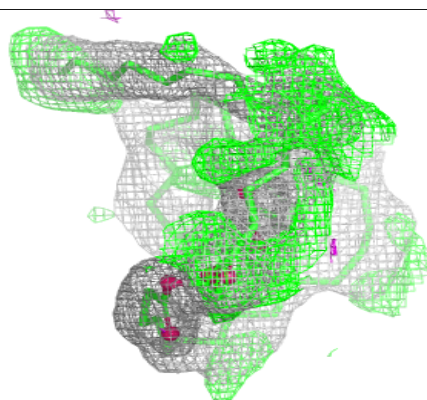
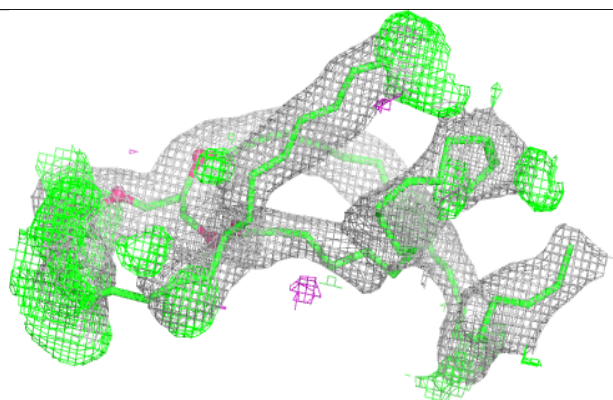
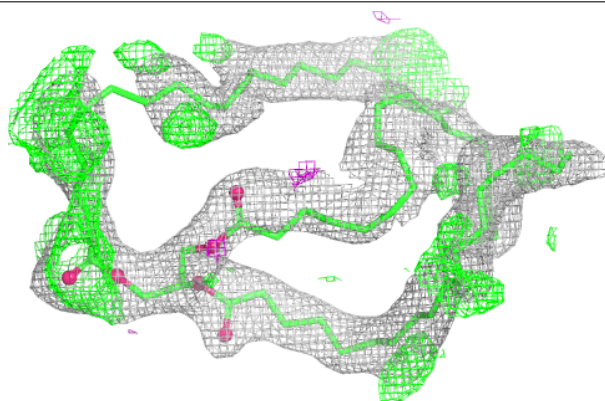


Electron density around PSC O 302:

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

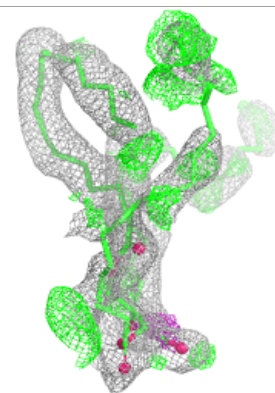
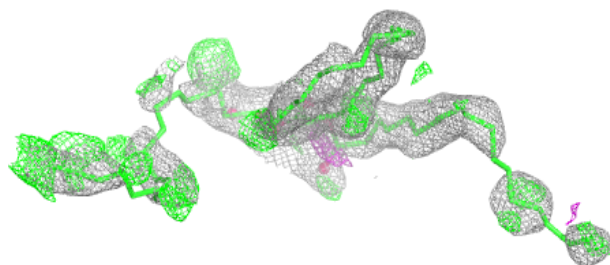
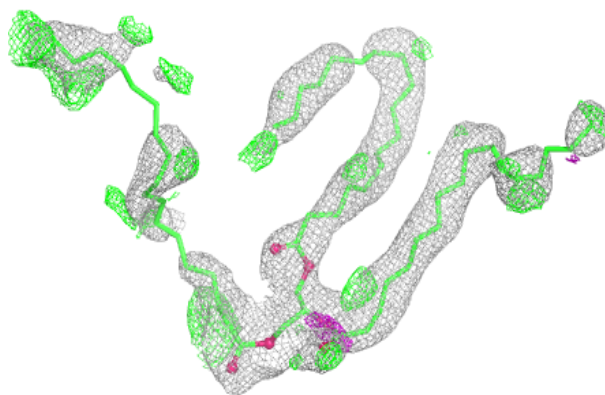
**Electron density around TGL A 608:**

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

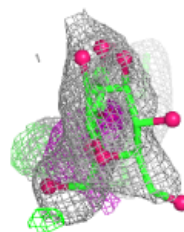
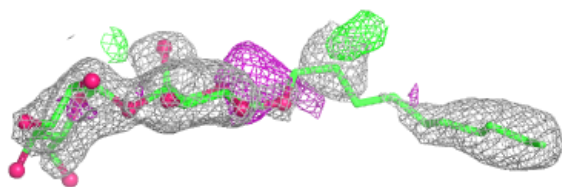
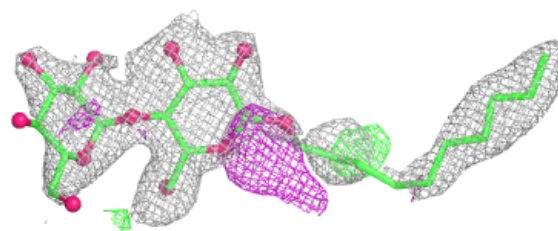


Electron density around TGL Q 201:

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

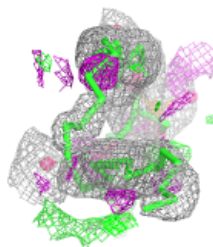
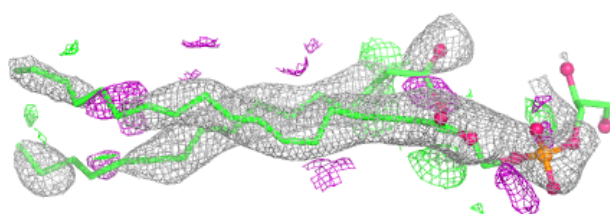
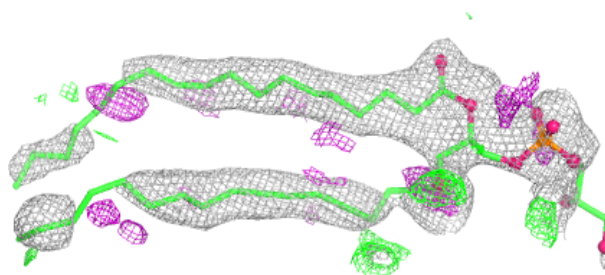
**Electron density around DMU P 308:**

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

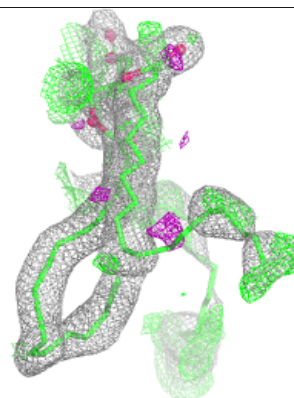
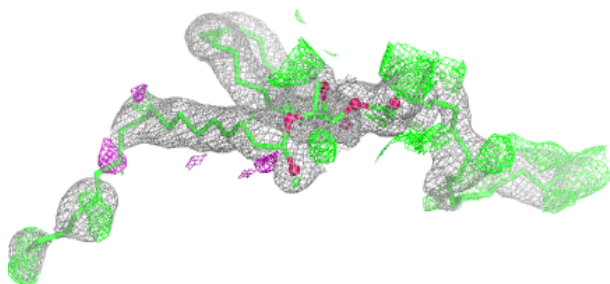
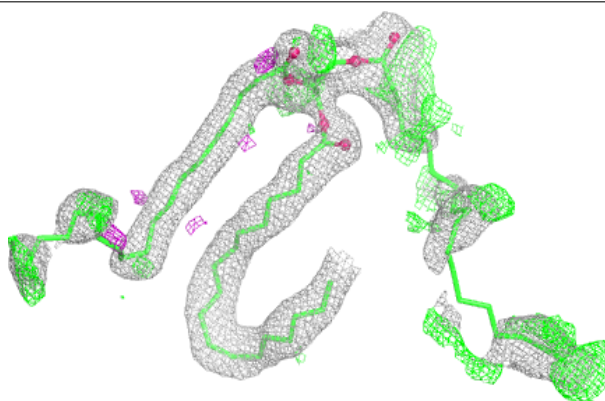


Electron density around PGV N 608:

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

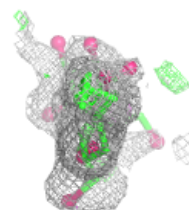
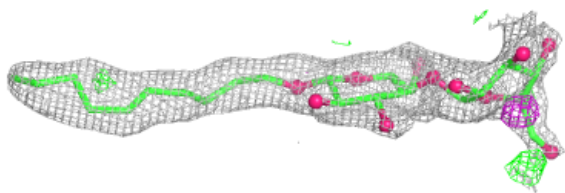
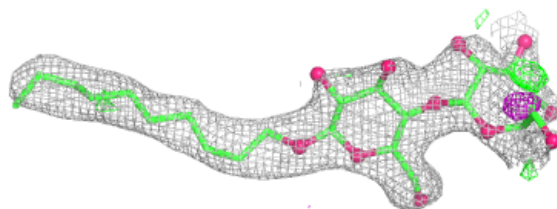
**Electron density around TGL D 201:**

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

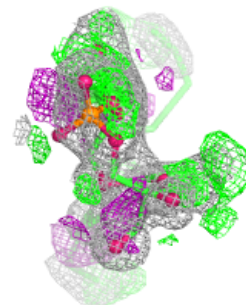
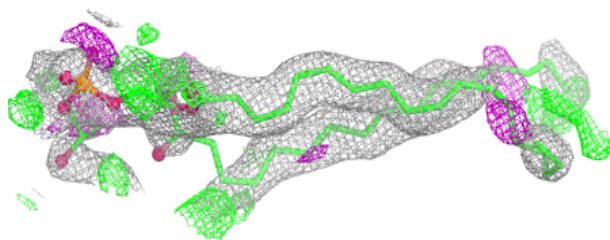
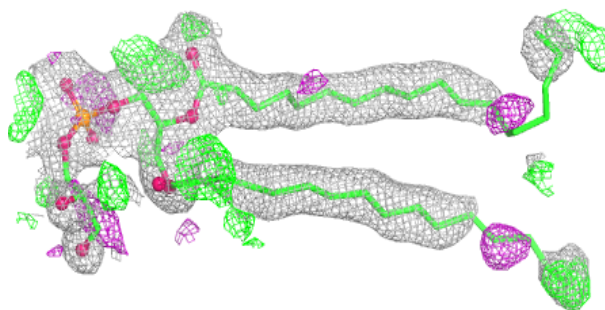


Electron density around DMU P 311:

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

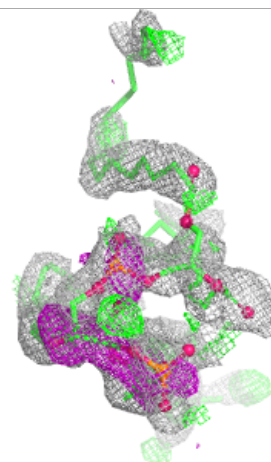
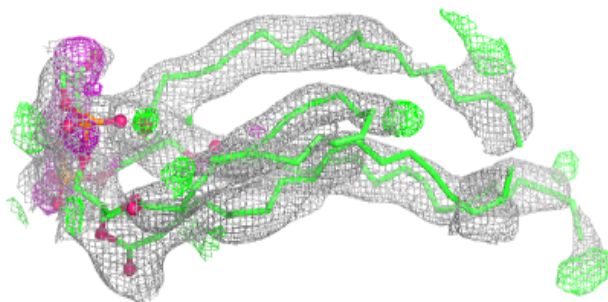
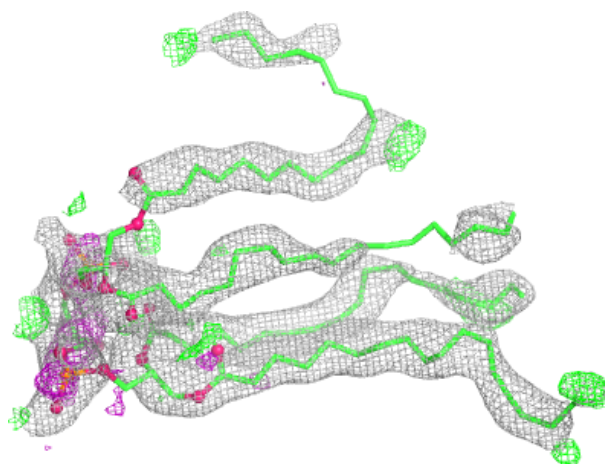
**Electron density around PGV M 101:**

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



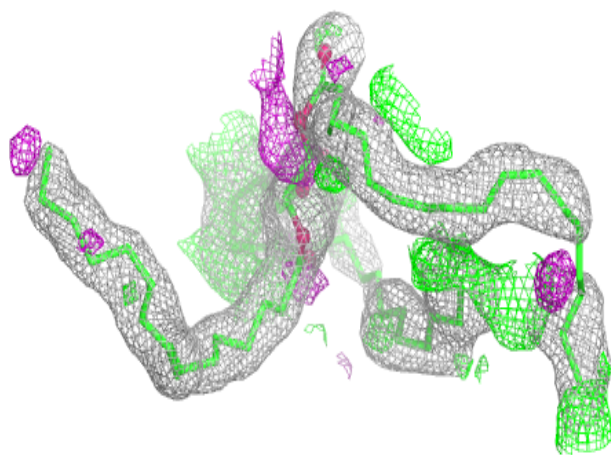
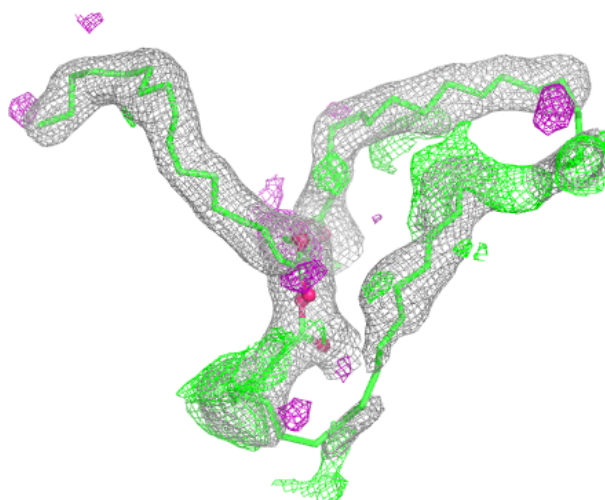
Electron density around CDL C 305:

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



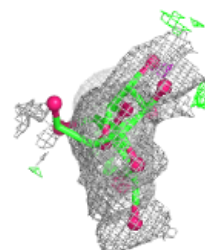
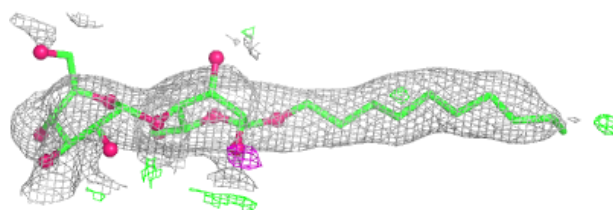
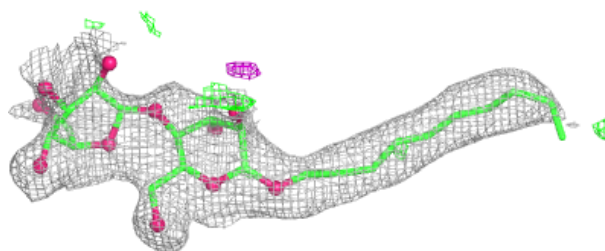
Electron density around TGL L 101:

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

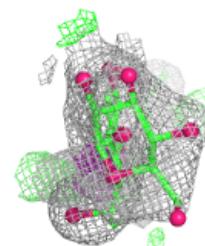
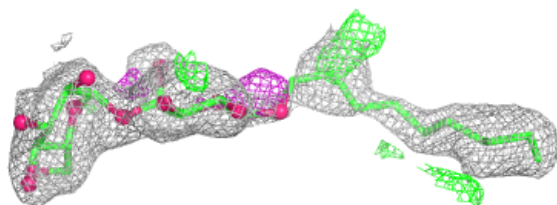
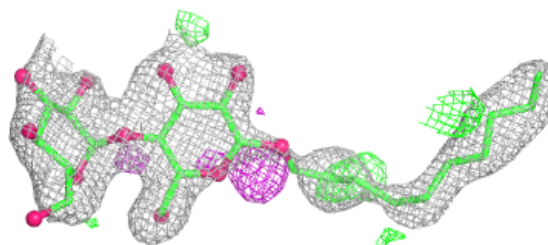


Electron density around DMU C 310:

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

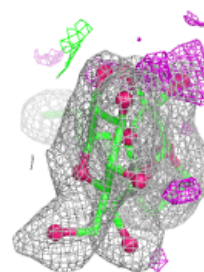
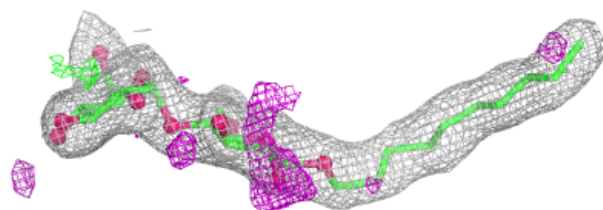
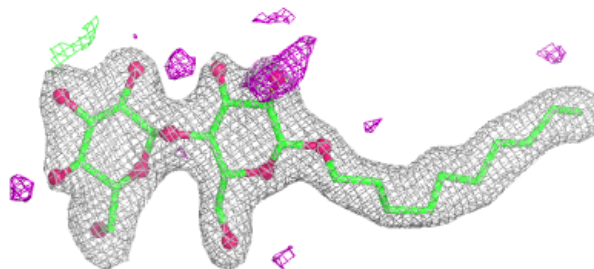
**Electron density around DMU C 302:**

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

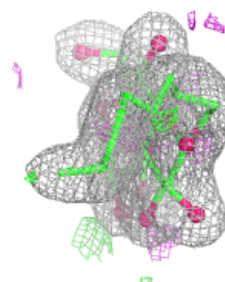
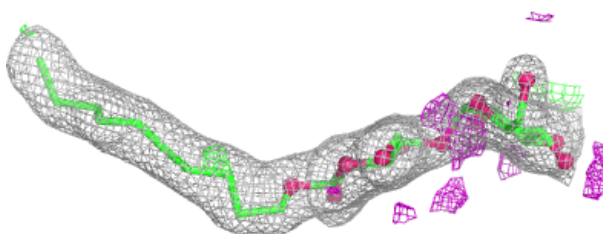
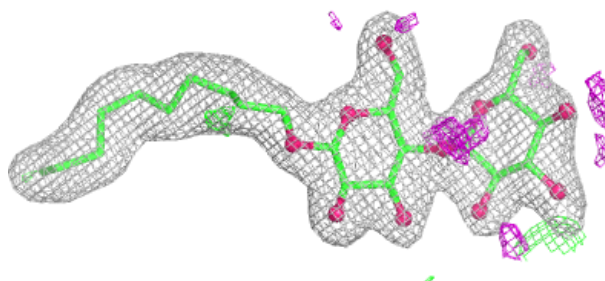


Electron density around DMU Z 101:

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

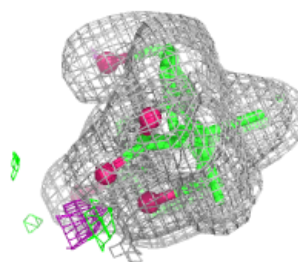
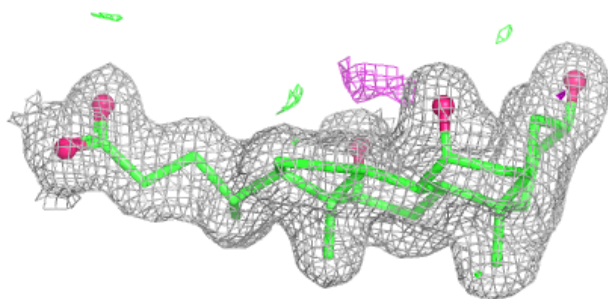
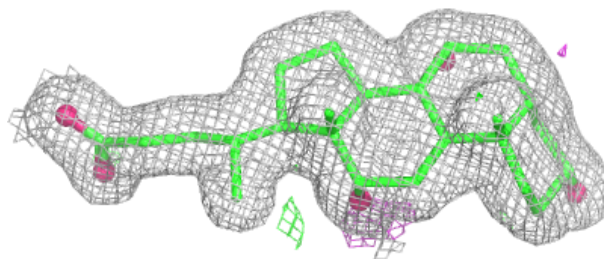
**Electron density around DMU M 102:**

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

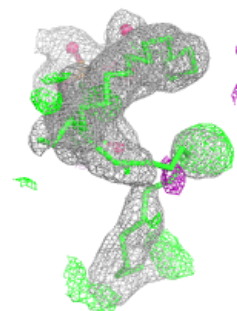
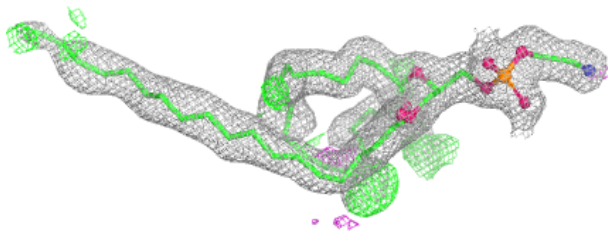
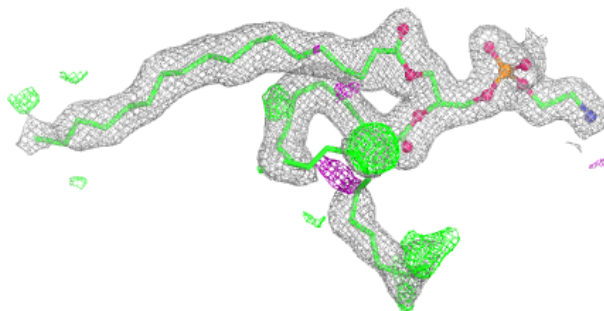


Electron density around CHD C 301:

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

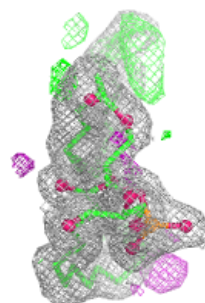
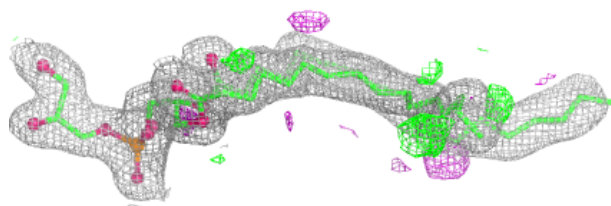
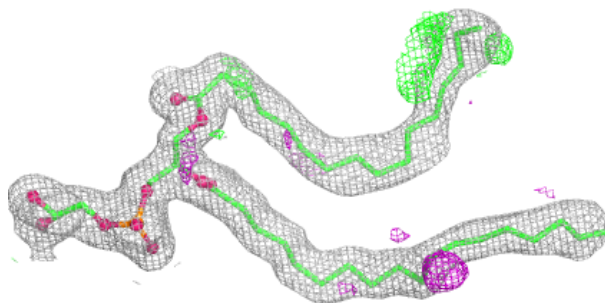
**Electron density around PEK P 304:**

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

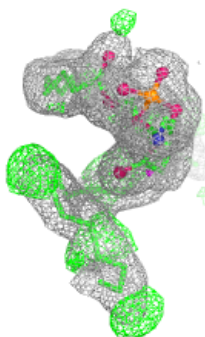
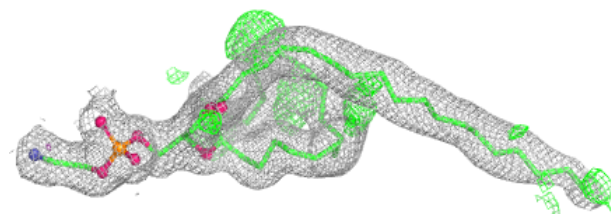
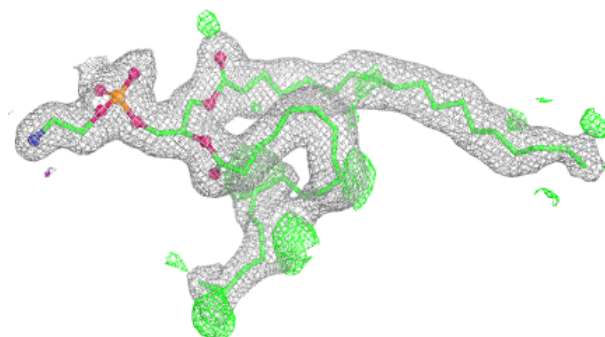


Electron density around PGV N 609:

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

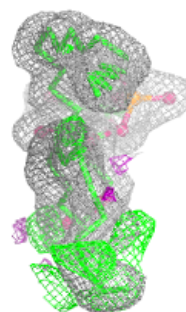
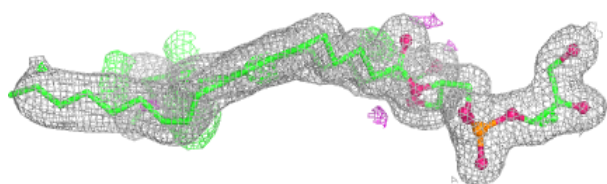
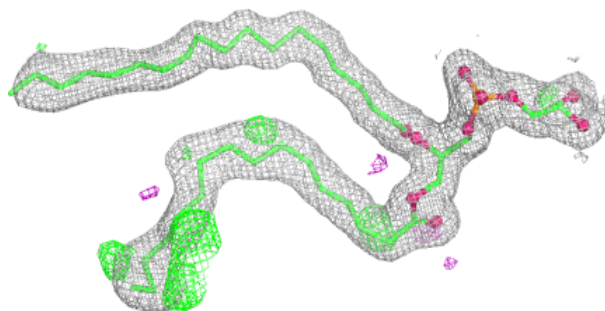
**Electron density around PEK G 101:**

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

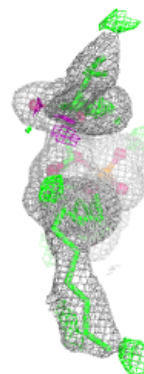
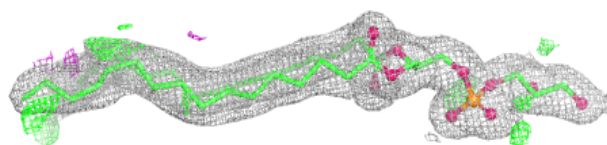
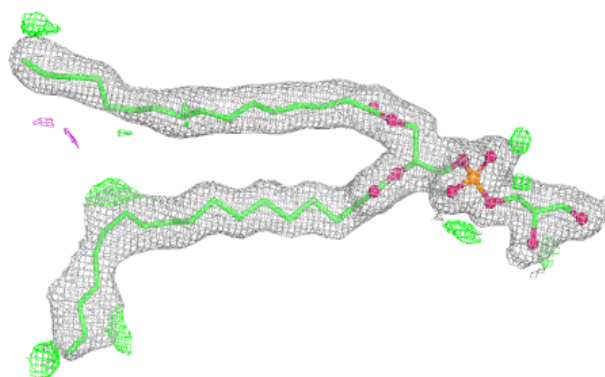


Electron density around PGV A 609:

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

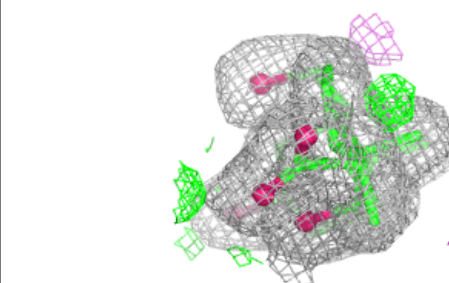
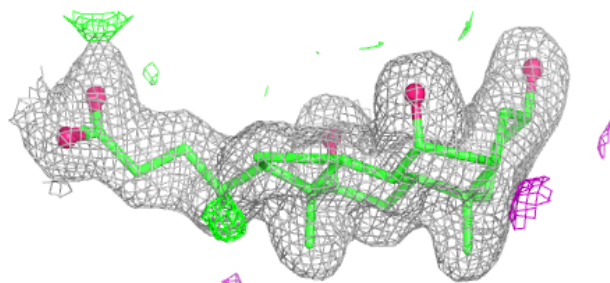
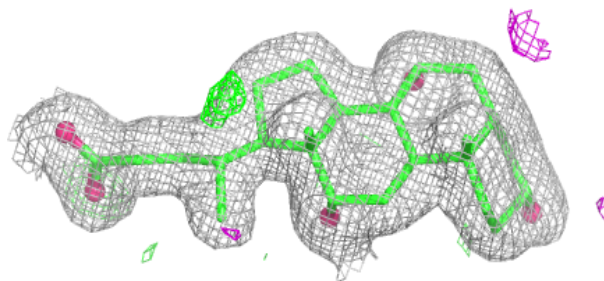
**Electron density around PGV P 305:**

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

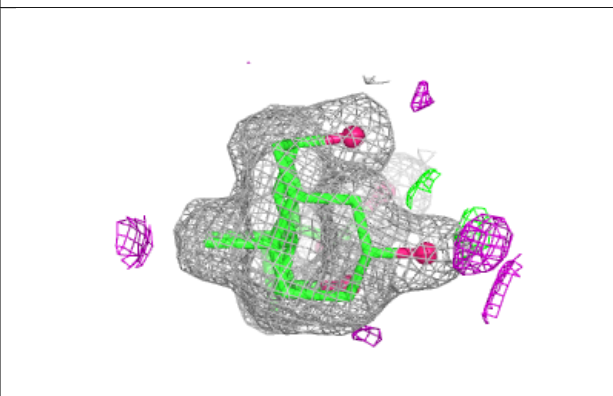
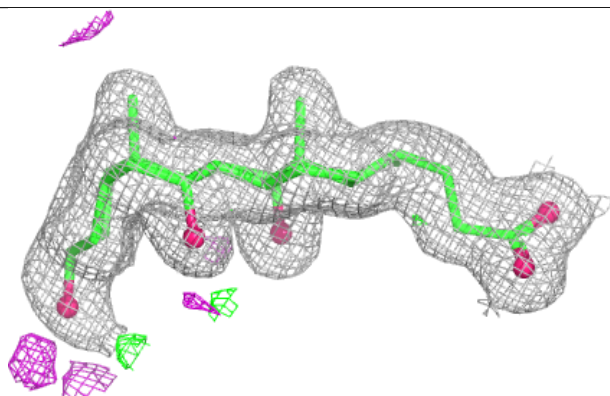
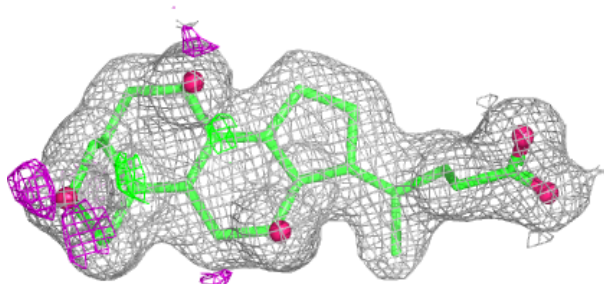


Electron density around CHD P 301:

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

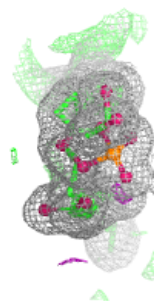
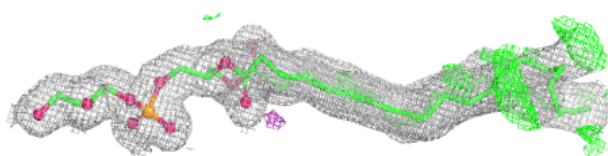
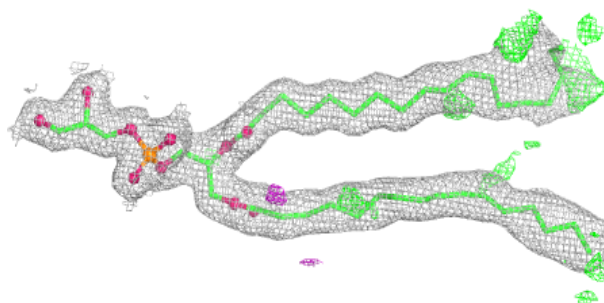
**Electron density around CHD T 101:**

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

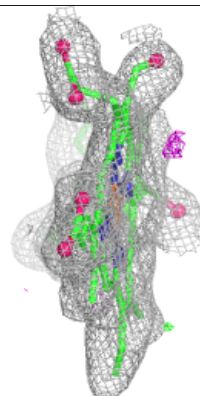
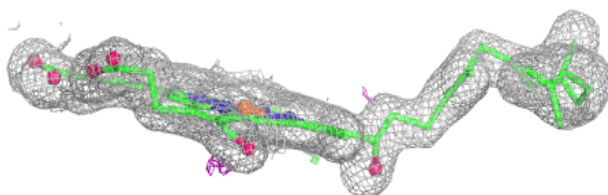
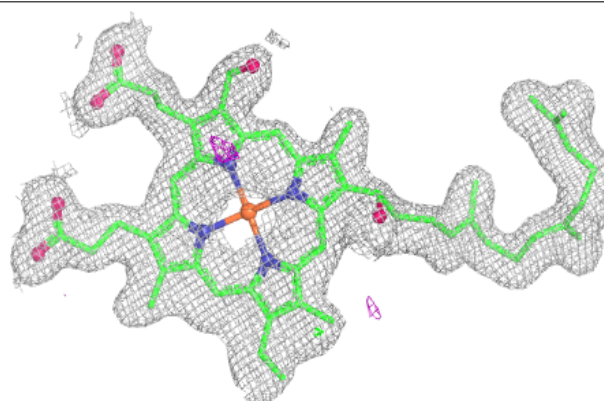


Electron density around PGV C 304:

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

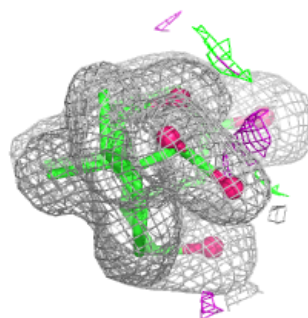
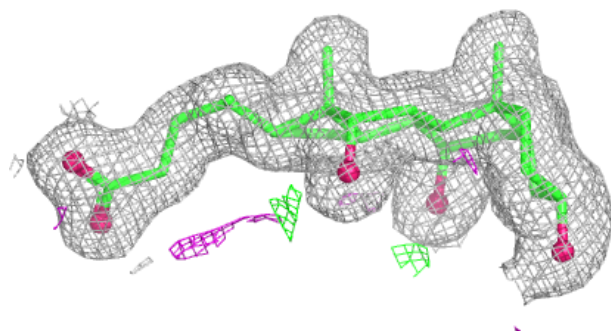
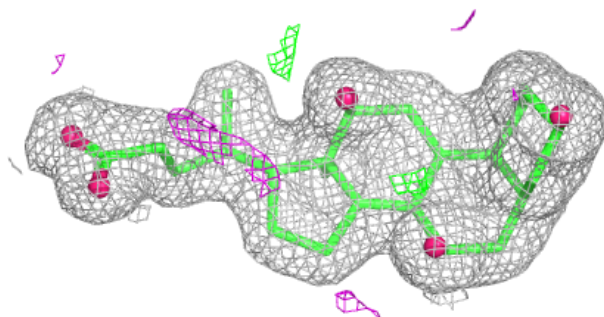
**Electron density around HEA A 602 (B):**

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

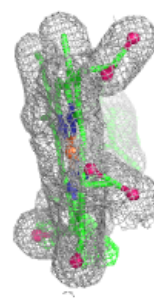
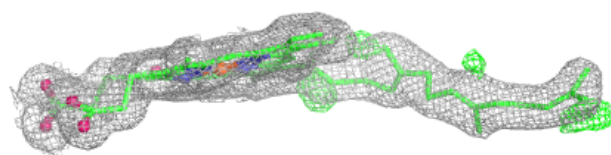
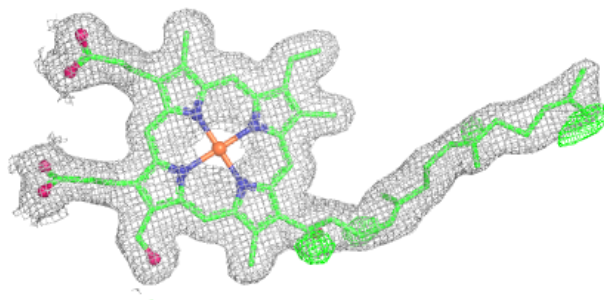


Electron density around CHD G 103:

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

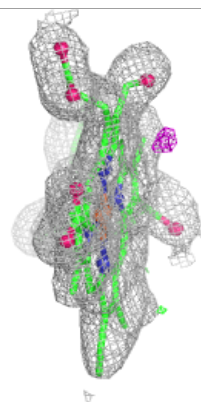
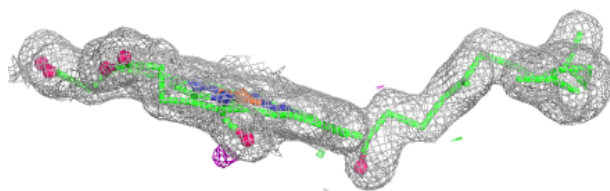
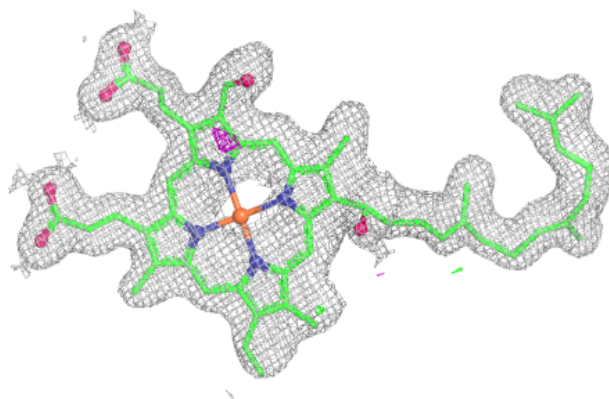
**Electron density around HEA N 601:**

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

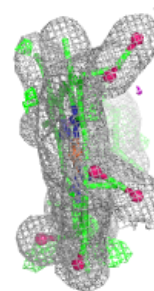
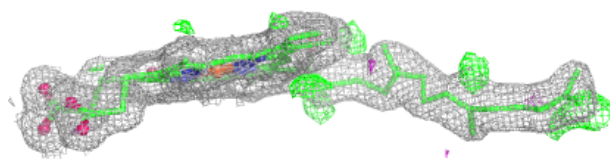
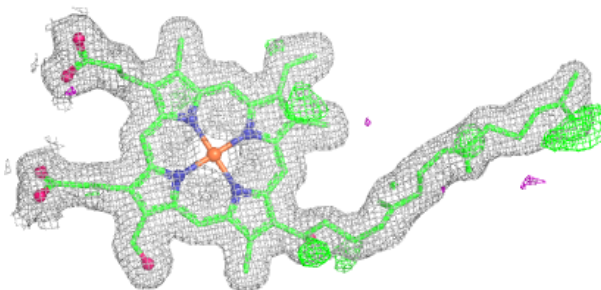


Electron density around HEA A 602 (A):

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

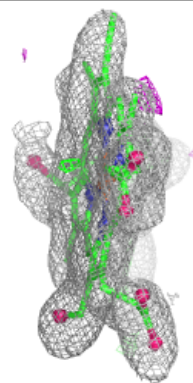
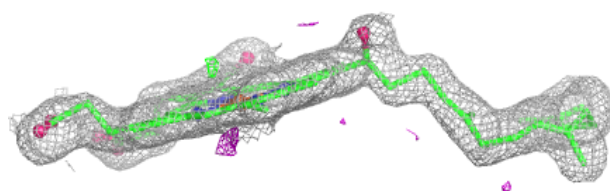
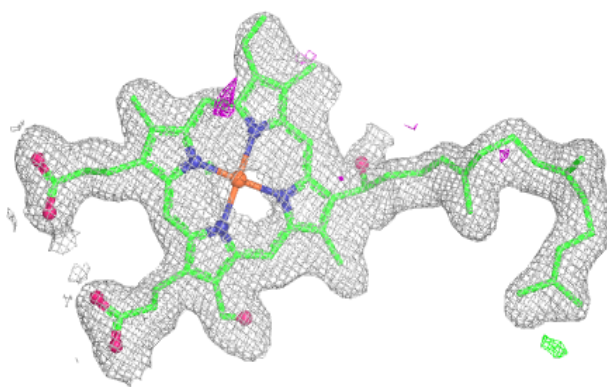
**Electron density around HEA A 601:**

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

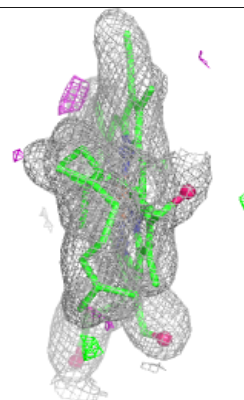
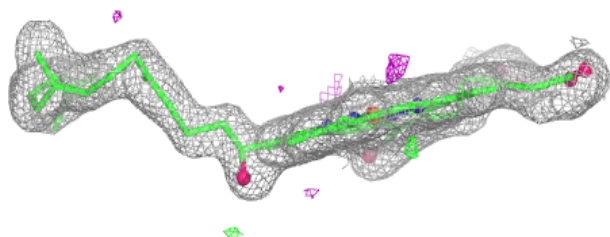
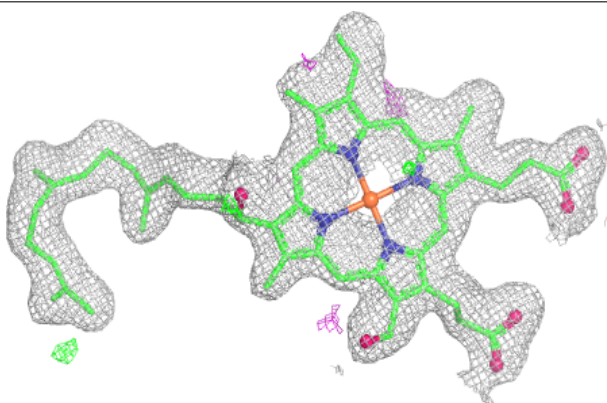


Electron density around HEA N 602 (A):

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

**Electron density around HEA N 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.