



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jun 22, 2024 – 04:27 PM EDT

PDB ID : 5B1A  
Title : Bovine heart cytochrome c oxidase in the fully oxidized state at 1.5 angstrom resolution  
Authors : Yano, N.; Muramoto, K.; Shimada, A.; Takemura, S.; Baba, J.; Fujisawa, H.; Mochizuki, M.; Shinzawa-Itoh, K.; Yamashita, E.; Tsukihara, T.; Yoshikawa, S.  
Deposited on : 2015-12-01  
Resolution : 1.50 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](mailto: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>.

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The following versions of software and data (see [references](#) ①) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

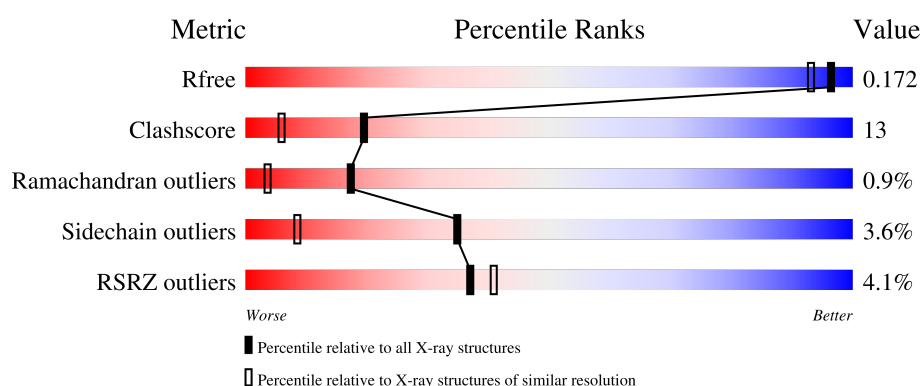
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.50 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.






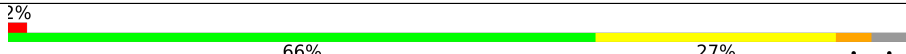
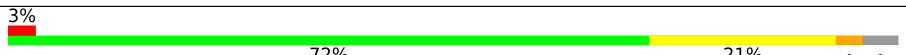
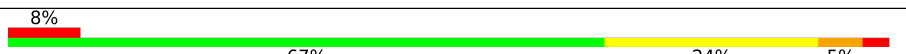
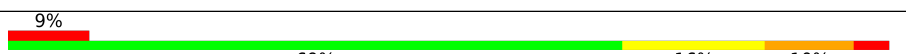
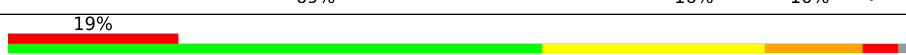

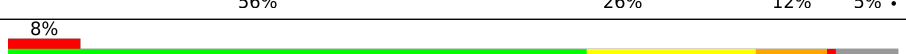
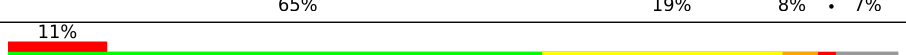

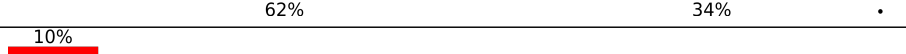
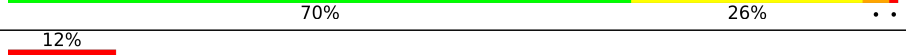





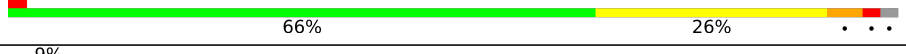
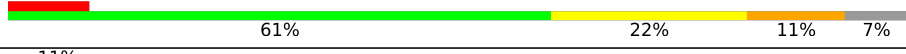
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	2936 (1.50-1.50)
Clashscore	141614	3144 (1.50-1.50)
Ramachandran outliers	138981	3066 (1.50-1.50)
Sidechain outliers	138945	3064 (1.50-1.50)
RSRZ outliers	127900	2884 (1.50-1.50)

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  $\geq 3$ , 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	<div> <div>73%</div> <div>22%</div> <div>.</div> </div>
1	N	514	<div> <div>72%</div> <div>24%</div> <div>.</div> </div>
2	B	227	<div> <div>2%</div> <div>65%</div> <div>29%</div> <div>.</div> </div>
2	O	227	<div> <div>2%</div> <div>69%</div> <div>26%</div> <div>.</div> </div>
3	C	261	<div> <div>72%</div> <div>25%</div> <div>.</div> </div>

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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
2	FME	B	1	-	X	X	-
20	TGL	D	201	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
20	TGL	Y	101	-	X	X	-
22	CHD	J	102	-	-	-	X
22	CHD	W	101	-	-	-	X
25	CDL	C	303	-	-	X	-
25	CDL	G	102	-	-	X	-
25	CDL	P	304	-	-	X	-
25	CDL	T	103	-	-	X	-
9	SAC	V	1	-	X	-	-

## 2 Entry composition

There are 29 unique types of molecules in this entry. The entry contains 35054 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
			4168	2778	645	704	41			
1	N	514	Total	C	N	O	S	0	16	0
			4154	2771	643	699	41			

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

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

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

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

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	144	Total	C	N	O	S	0	5	0
			1242	809	206	223	4			
4	Q	144	Total	C	N	O	S	0	3	0
			1224	797	202	221	4			

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

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

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

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

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

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

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

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

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

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

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

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

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

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

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

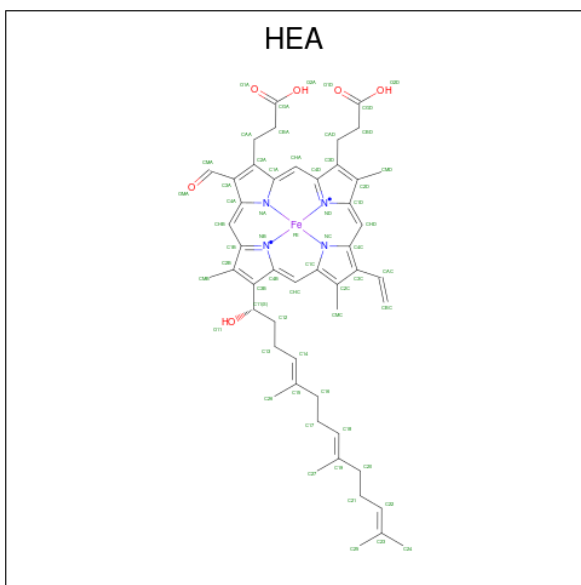
- Molecule 12 is a protein called Cytochrome c oxidase subunit 7C, mitochondrial.

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

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

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

- Molecule 14 is HEME-A (three-letter code: HEA) (formula: C<sub>49</sub>H<sub>56</sub>FeN<sub>4</sub>O<sub>6</sub>).



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 60	C 49	Fe 1	N 4	O 6	0	0
14	N	1	Total 60	C 49	Fe 1	N 4	O 6	0	0
14	N	1	Total 60	C 49	Fe 1	N 4	O 6	0	0

- Molecule 15 is COPPER (II) ION (three-letter code: 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		

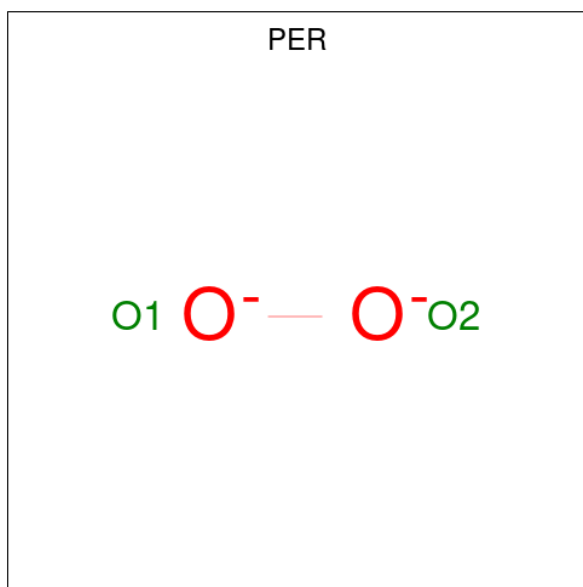
- Molecule 16 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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 (three-letter code: NA) (formula: Na).

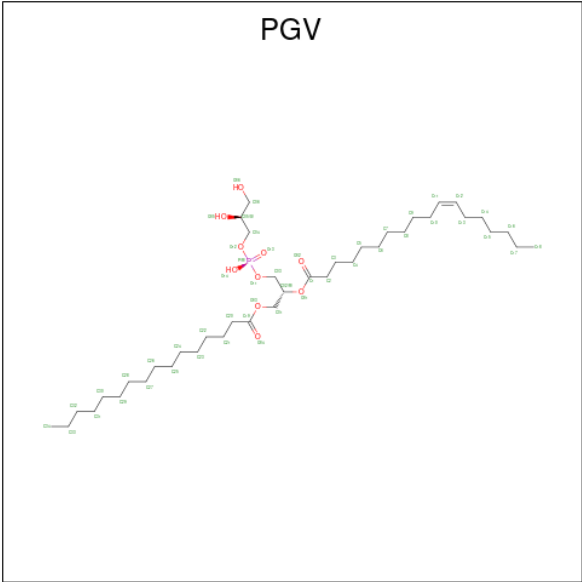
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
17	A	1	Total	Na	0	0
			1	1		
17	N	1	Total	Na	0	0
			1	1		

- Molecule 18 is PEROXIDE ION (three-letter code: PER) (formula: O<sub>2</sub>).



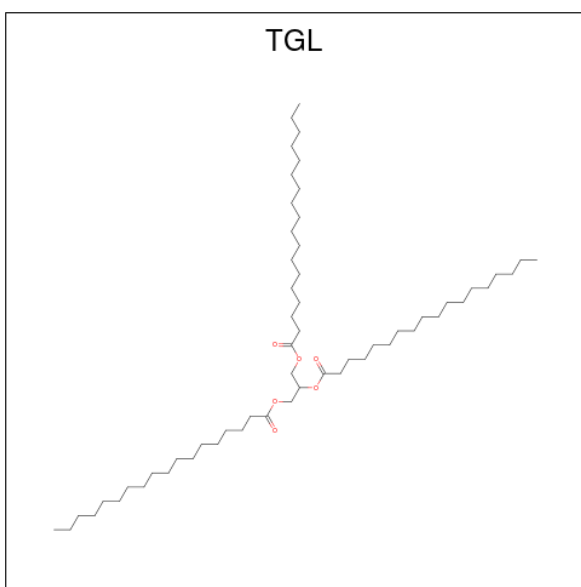
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
18	A	1	Total	O	0	0
			2	2		
18	N	1	Total	O	0	0
			2	2		

- Molecule 19 is (1R)-2-{{[(2S)-2,3-DIHYDROXYPROPYL]OXY}(HYDROXY)PHOSPHORYL]OXY}-1-[(PALMITOYLOXY)METHYL]ETHYL (11E)-OCTADEC-11-ENOATE (three-letter code: PGV) (formula: C<sub>40</sub>H<sub>77</sub>O<sub>10</sub>P).



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

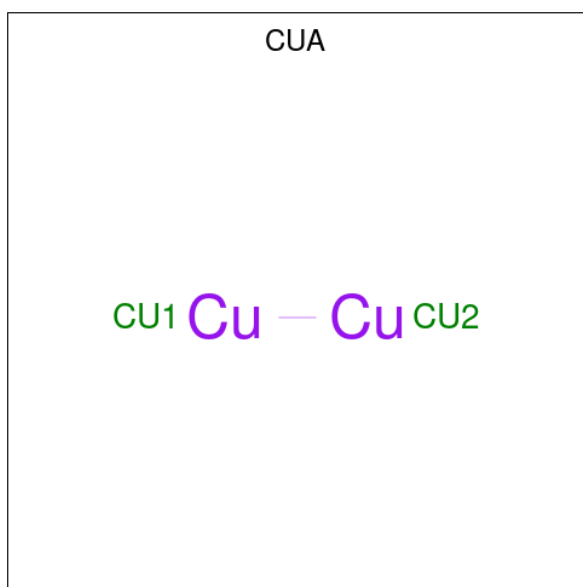
- Molecule 20 is TRISTEAROYLGLYCEROL (three-letter code: TGL) (formula: C<sub>57</sub>H<sub>110</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
20	B	1	Total	C	O	0	0
			63	57	6		
20	D	1	Total	C	O	0	0
			63	57	6		
20	L	1	Total	C	O	0	0
			63	57	6		
20	N	1	Total	C	O	0	0
			63	57	6		
20	Q	1	Total	C	O	0	0
			63	57	6		
20	Y	1	Total	C	O	0	0
			63	57	6		

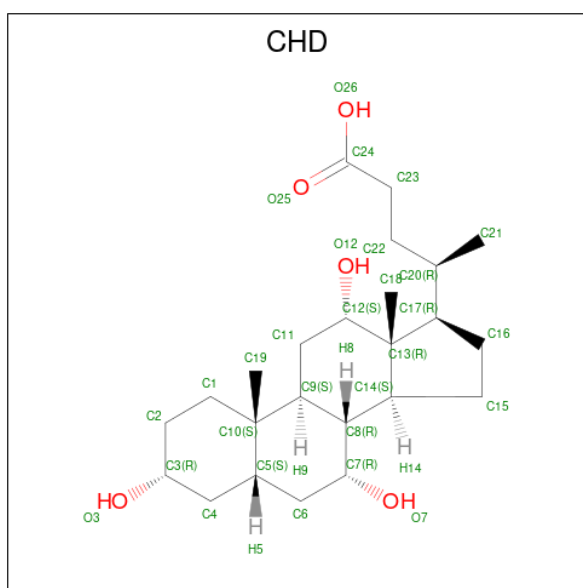
- Molecule 21 is DINUCLEAR COPPER ION (three-letter code: CUA) (formula: Cu<sub>2</sub>).





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

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



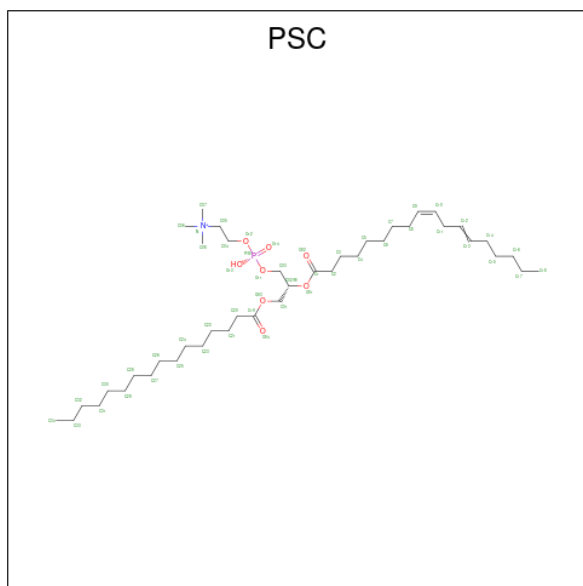
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
22	B	1	Total	C	O	0	0
			29	24	5		
22	C	1	Total	C	O	0	0
			29	24	5		

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

- Molecule 23 is (7R,17E,20E)-4-HYDROXY-N,N,N-TRIMETHYL-9-OXO-7-[(PALMITOYLOXY)METHYL]-3,5,8-TRIOXA-4-PHOSPHAHEXACOSA-17,20-DIEN-1-AMINIUM 4-OXIDE (three-letter code: PSC) (formula: C<sub>42</sub>H<sub>81</sub>NO<sub>8</sub>P).

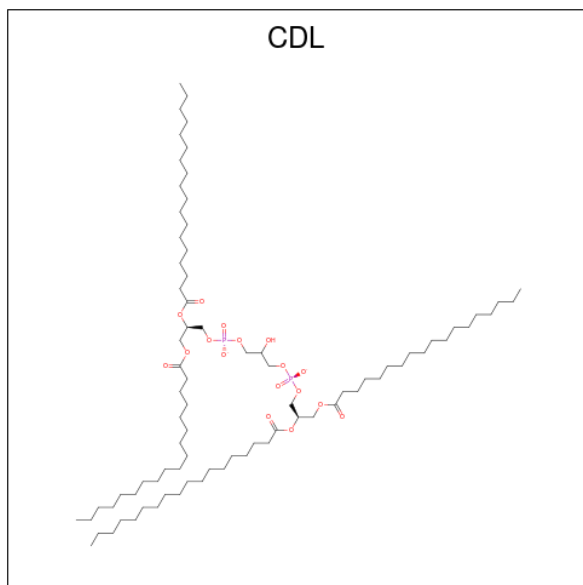


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
23	B	1	Total	C	N	O	P	0	0
			52	42	1	8	1		
23	O	1	Total	C	N	O	P	0	0
			52	42	1	8	1		

- Molecule 24 is UNKNOWN ATOM OR ION (three-letter code: UNX) (formula: X).

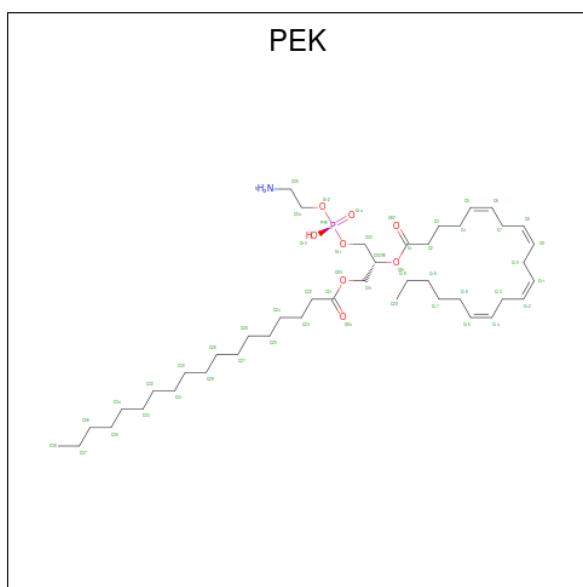
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
24	C	1	Total X 1 1	0	0
24	P	1	Total X 1 1	0	0

- Molecule 25 is CARDIOLIPIN (three-letter code: CDL) (formula:  $C_{81}H_{156}O_{17}P_2$ ).



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

- Molecule 26 is (1S)-2-{[(2-AMINOETHOXY)(HYDROXY)PHOSPHORYL]OXY}-1-[(STEAROYLOXY)METHYL]ETHYL (5E,8E,11E,14E)-ICOSA-5,8,11,14-TETRAENOATE (three-letter code: PEK) (formula:  $C_{43}H_{78}NO_8P$ ).

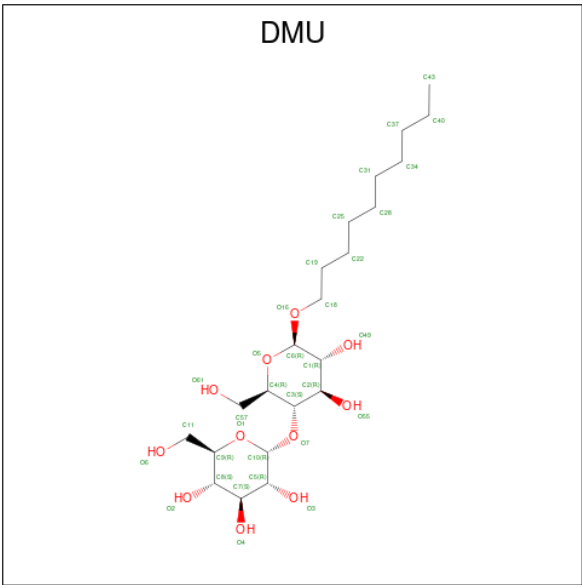


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
26	C	1	Total	C	N	O	P	0	0
			53	43	1	8	1		
26	G	1	Total	C	N	O	P	0	0
			53	43	1	8	1		
26	G	1	Total	C	N	O	P	0	0
			53	43	1	8	1		
26	P	1	Total	C	N	O	P	0	0
			53	43	1	8	1		
26	T	1	Total	C	N	O	P	0	0
			53	43	1	8	1		
26	T	1	Total	C	N	O	P	0	0
			53	43	1	8	1		

- Molecule 27 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 28 is DECYL-BETA-D-MALTOPYRANOSIDE (three-letter code: DMU) (formula: C<sub>22</sub>H<sub>42</sub>O<sub>11</sub>).



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

- Molecule 29 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
29	A	297	Total	O	0	0
			297	297		
29	B	273	Total	O	0	1
			274	274		
29	C	176	Total	O	0	0
			176	176		
29	D	266	Total	O	0	0
			266	266		
29	E	178	Total	O	0	0
			178	178		
29	F	199	Total	O	0	0
			199	199		
29	G	100	Total	O	0	0
			100	100		
29	H	122	Total	O	0	0
			122	122		

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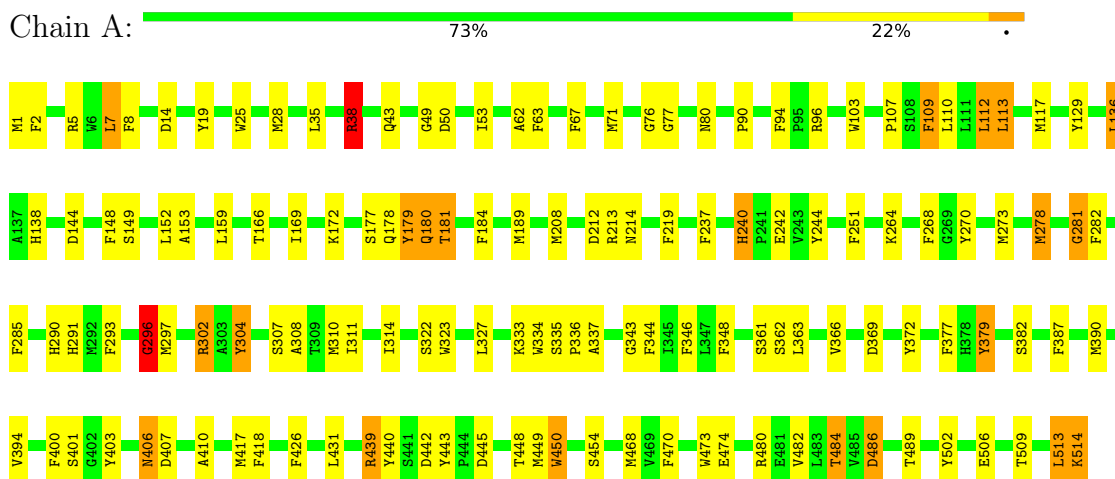
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
29	I	88	Total 88	O 88	0	0
29	J	63	Total 63	O 63	0	0
29	K	69	Total 69	O 69	0	0
29	L	48	Total 48	O 48	0	0
29	M	47	Total 47	O 47	0	0
29	N	290	Total 290	O 290	0	0
29	O	242	Total 243	O 243	0	1
29	P	173	Total 173	O 173	0	0
29	Q	164	Total 164	O 164	0	0
29	R	151	Total 151	O 151	0	0
29	S	186	Total 186	O 186	0	0
29	T	94	Total 94	O 94	0	0
29	U	110	Total 110	O 110	0	0
29	V	71	Total 71	O 71	0	0
29	W	58	Total 58	O 58	0	0
29	X	57	Total 57	O 57	0	0
29	Y	40	Total 40	O 40	0	0
29	Z	37	Total 37	O 37	0	0

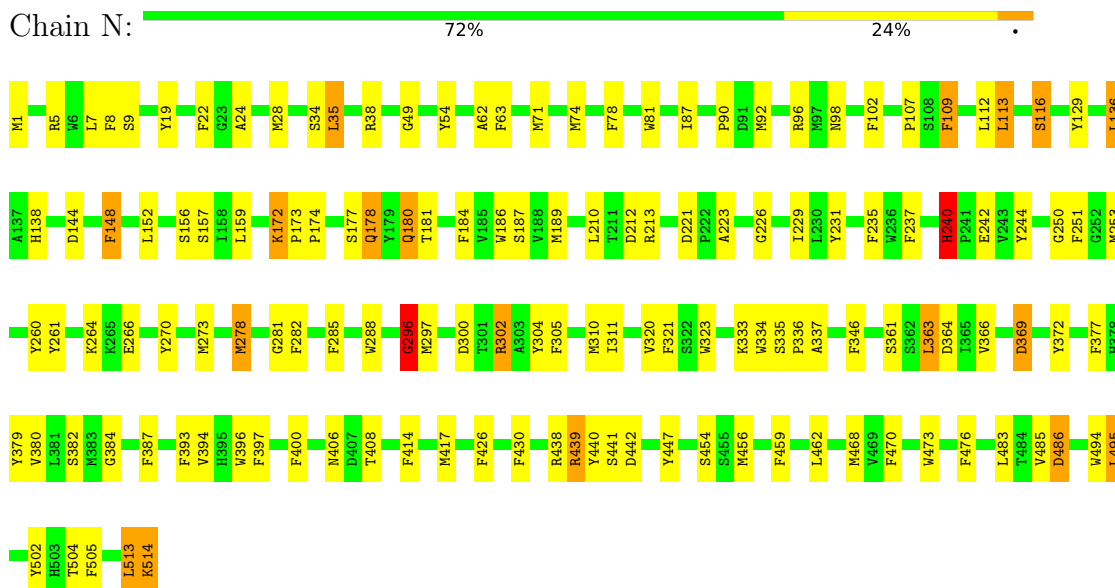
### 3 Residue-property plots

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.

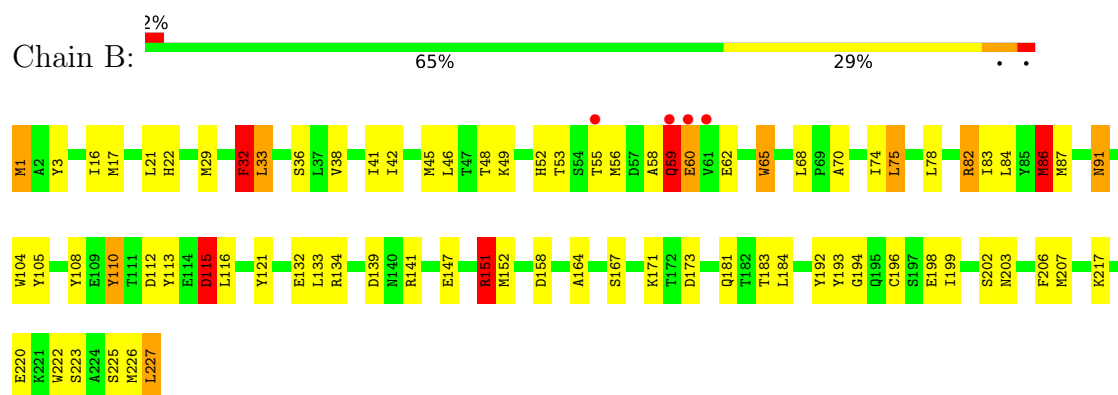
#### • Molecule 1: Cytochrome c oxidase subunit 1



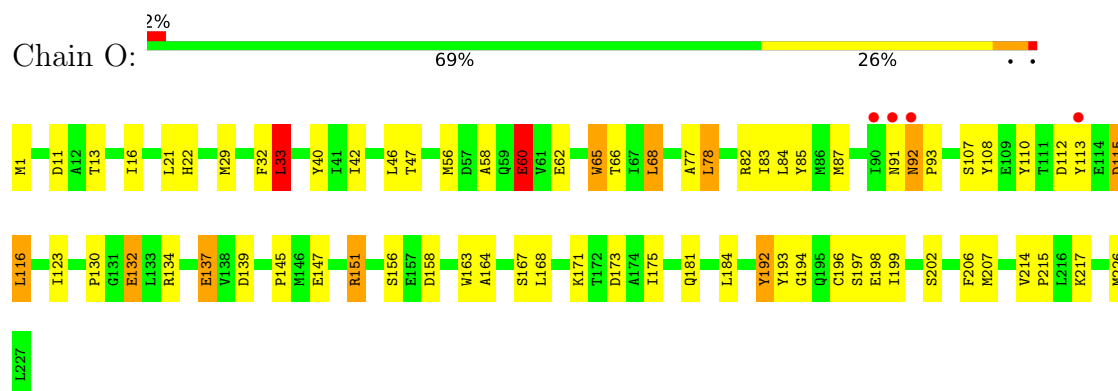
#### • Molecule 1: Cytochrome c oxidase subunit 1



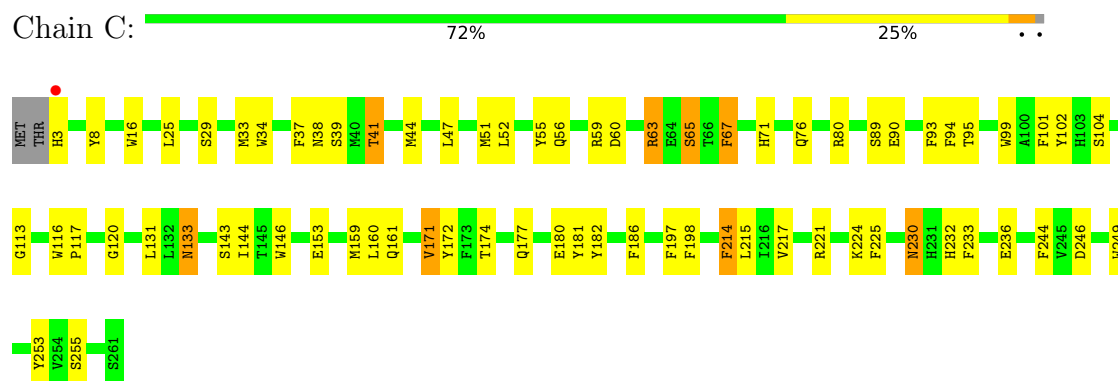
#### • Molecule 2: Cytochrome c oxidase subunit 2



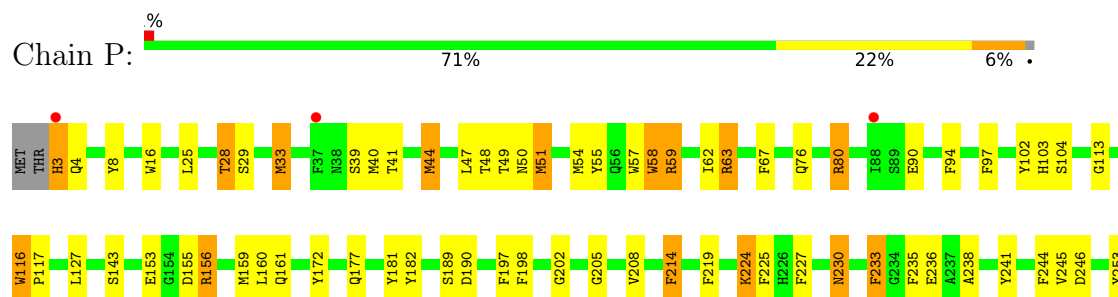
• Molecule 2: Cytochrome c oxidase subunit 2



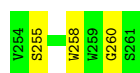
• Molecule 3: Cytochrome c oxidase subunit 3



• Molecule 3: Cytochrome c oxidase subunit 3



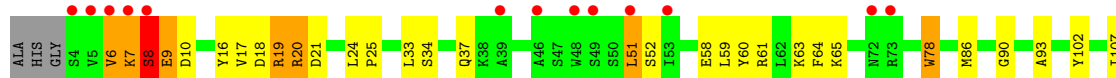




- Molecule 4: Cytochrome c oxidase subunit 4 isoform 1, mitochondrial



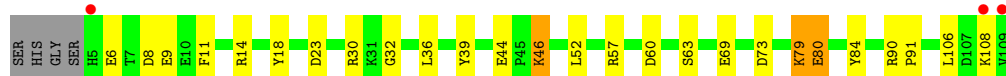
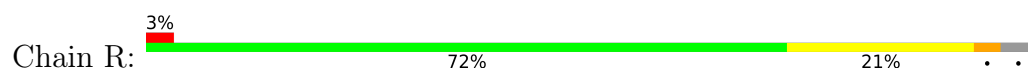
- Molecule 4: Cytochrome c oxidase subunit 4 isoform 1, mitochondrial



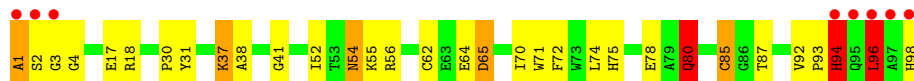
- Molecule 5: Cytochrome c oxidase subunit 5A, mitochondrial



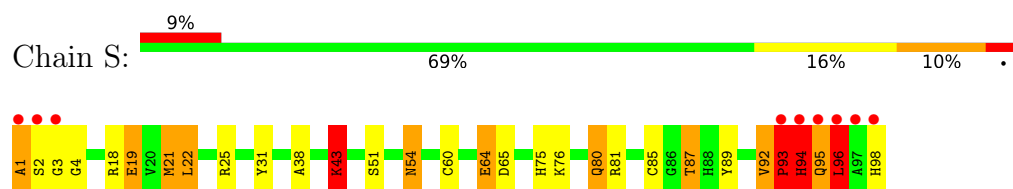
- Molecule 5: Cytochrome c oxidase subunit 5A, mitochondrial



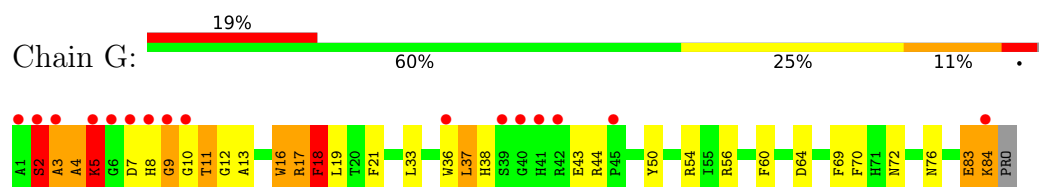
- Molecule 6: Cytochrome c oxidase subunit 5B, mitochondrial



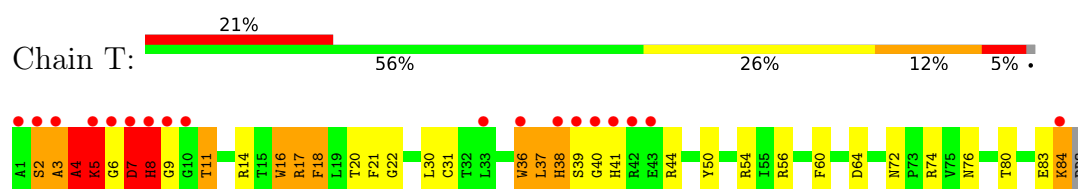
- Molecule 6: Cytochrome c oxidase subunit 5B, mitochondrial



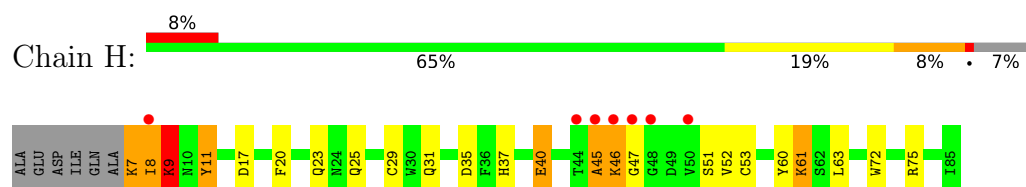
- Molecule 7: Cytochrome c oxidase subunit 6A2, mitochondrial



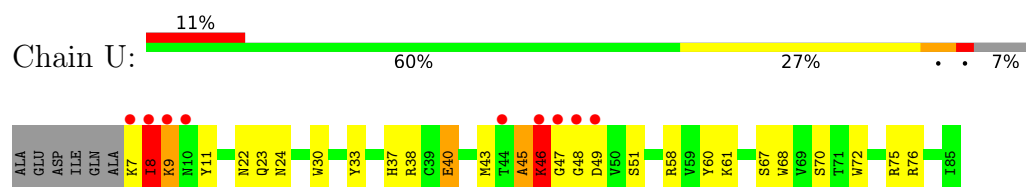
- Molecule 7: Cytochrome c oxidase subunit 6A2, mitochondrial



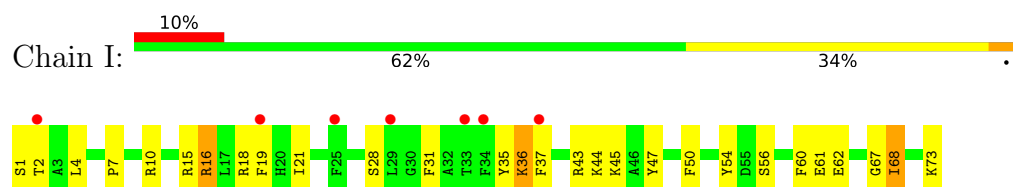
- Molecule 8: Cytochrome c oxidase subunit 6B1



- Molecule 8: Cytochrome c oxidase subunit 6B1

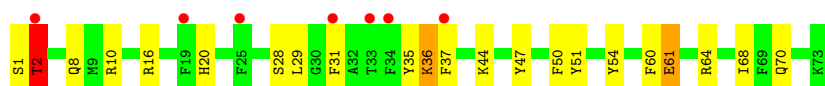


- Molecule 9: Cytochrome c oxidase subunit 6C



- Molecule 9: Cytochrome c oxidase subunit 6C

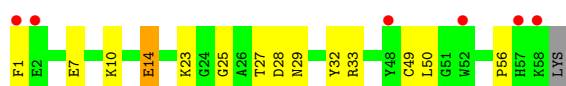
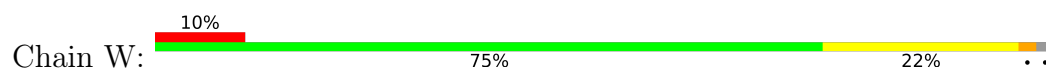




- Molecule 10: Cytochrome c oxidase subunit 7A1, mitochondrial



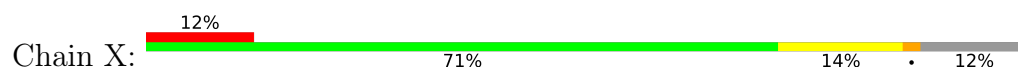
- Molecule 10: Cytochrome c oxidase subunit 7A1, mitochondrial



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



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



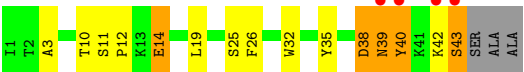
- Molecule 12: Cytochrome c oxidase subunit 7C, mitochondrial



- 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, $\alpha$ , $\beta$ , $\gamma$	181.94Å 204.40Å 177.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 1.50 89.10 – 1.40	Depositor EDS
% Data completeness (in resolution range)	98.3 (40.00-1.50) 98.2 (89.10-1.40)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.86 (at 1.40Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, $R_{free}$	0.149 , 0.172 0.149 , 0.172	Depositor DCC
$R_{free}$ test set	63174 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	20.6	Xtriage
Anisotropy	0.642	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 66.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.001 for l,-k,h	Xtriage
$F_o, F_c$ correlation	0.98	EDS
Total number of atoms	35054	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.68% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: PSC, PGV, HEA, TGL, CUA, CDL, ZN, NA, TPO, DMU, MG, PER, UNX, PEK, CU, SAC, FME, CHD

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	1.84	54/4297 (1.3%)	1.89	89/5864 (1.5%)
1	N	1.85	57/4283 (1.3%)	1.76	84/5845 (1.4%)
2	B	1.94	37/1937 (1.9%)	1.80	32/2637 (1.2%)
2	O	1.82	30/1908 (1.6%)	1.57	23/2597 (0.9%)
3	C	1.91	41/2272 (1.8%)	1.79	46/3102 (1.5%)
3	P	1.91	37/2272 (1.6%)	1.77	52/3102 (1.7%)
4	D	2.05	31/1277 (2.4%)	1.87	40/1720 (2.3%)
4	Q	1.67	17/1259 (1.4%)	1.88	23/1698 (1.4%)
5	E	2.01	24/871 (2.8%)	2.45	41/1182 (3.5%)
5	R	1.89	18/882 (2.0%)	1.60	14/1196 (1.2%)
6	F	1.96	16/795 (2.0%)	1.65	7/1079 (0.6%)
6	S	1.89	17/780 (2.2%)	1.69	14/1058 (1.3%)
7	G	2.03	15/702 (2.1%)	1.87	19/953 (2.0%)
7	T	1.95	14/702 (2.0%)	1.59	9/953 (0.9%)
8	H	1.77	7/682 (1.0%)	1.52	9/921 (1.0%)
8	U	1.74	10/682 (1.5%)	1.36	5/921 (0.5%)
9	I	1.96	13/605 (2.1%)	1.73	13/802 (1.6%)
9	V	1.70	7/605 (1.2%)	1.85	9/802 (1.1%)
10	J	1.80	6/471 (1.3%)	1.54	4/636 (0.6%)
10	W	1.65	4/480 (0.8%)	1.36	2/648 (0.3%)
11	K	2.09	14/398 (3.5%)	1.91	7/546 (1.3%)
11	X	1.63	7/405 (1.7%)	1.47	5/556 (0.9%)
12	L	2.01	6/393 (1.5%)	1.76	11/526 (2.1%)
12	Y	1.97	11/401 (2.7%)	1.52	3/536 (0.6%)
13	M	1.80	6/345 (1.7%)	1.68	5/470 (1.1%)
13	Z	1.70	4/345 (1.2%)	1.43	3/470 (0.6%)
All	All	1.87	503/30049 (1.7%)	1.77	569/40820 (1.4%)

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	4
1	N	0	3
2	B	0	3
3	C	0	1
4	Q	0	1
5	E	0	2
6	F	0	1
6	S	0	2
7	G	0	1
7	T	0	1
10	J	0	1
11	K	0	1
13	M	0	1
All	All	0	22

The worst 5 of 503 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	O	65	TRP	CB-CG	-17.96	1.18	1.50
2	B	65	TRP	CB-CG	-17.06	1.19	1.50
4	D	58	GLU	CD-OE1	15.90	1.43	1.25
11	K	47	ARG	CZ-NH2	14.60	1.52	1.33
7	T	36	TRP	CB-CG	13.34	1.74	1.50

The worst 5 of 569 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	E	90	ARG	NE-CZ-NH1	35.49	138.05	120.30
4	Q	20	ARG	NE-CZ-NH1	32.93	136.77	120.30
4	Q	20	ARG	NE-CZ-NH2	-31.64	104.48	120.30
5	E	90	ARG	NE-CZ-NH2	-24.93	107.83	120.30
9	V	10	ARG	NE-CZ-NH2	-21.32	109.64	120.30

There are no chirality outliers.

5 of 22 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	296	GLY	Mainchain
1	A	304	TYR	Sidechain
1	A	379	TYR	Sidechain

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Mol	Chain	Res	Type	Group
1	A	38	ARG	Sidechain
2	B	68	LEU	Mainchain

## 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	4168	0	4137	74	0
1	N	4154	0	4129	81	0
2	B	1899	0	1898	65	0
2	O	1870	0	1868	40	0
3	C	2185	0	2097	38	0
3	P	2185	0	2097	51	0
4	D	1242	0	1235	19	0
4	Q	1224	0	1211	25	0
5	E	852	0	845	1	0
5	R	863	0	857	8	2
6	F	778	0	754	27	0
6	S	763	0	742	42	0
7	G	686	0	651	42	0
7	T	686	0	651	45	0
8	H	662	0	623	24	0
8	U	662	0	623	15	0
9	I	601	0	613	17	2
9	V	601	0	613	13	0
10	J	460	0	459	13	0
10	W	469	0	464	9	0
11	K	384	0	366	6	0
11	X	391	0	374	4	0
12	L	380	0	380	19	0
12	Y	388	0	388	29	0
13	M	335	0	352	7	0
13	Z	335	0	352	5	0
14	A	120	0	107	10	0
14	N	120	0	107	11	0
15	A	1	0	0	0	0
15	N	1	0	0	0	0
16	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
16	N	1	0	0	0	0
17	A	1	0	0	0	0
17	N	1	0	0	0	0
18	A	2	0	0	1	0
18	N	2	0	0	1	0
19	A	102	0	152	11	0
19	C	102	0	152	8	0
19	N	51	0	76	1	0
19	P	102	0	152	9	0
19	Q	51	0	76	11	0
20	B	63	0	109	3	0
20	D	63	0	106	21	0
20	L	63	0	110	19	0
20	N	63	0	110	6	0
20	Q	63	0	110	15	0
20	Y	63	0	110	23	0
21	B	2	0	0	0	0
21	O	2	0	0	0	0
22	B	29	0	39	0	0
22	C	58	0	77	3	0
22	J	29	0	37	6	0
22	O	29	0	39	0	0
22	P	58	0	77	7	0
22	W	29	0	38	6	0
23	B	52	0	80	13	0
23	O	52	0	80	19	0
24	C	1	0	0	0	0
24	P	1	0	0	0	0
25	C	100	0	154	26	0
25	G	100	0	156	37	0
25	P	100	0	156	24	0
25	T	100	0	156	28	0
26	C	53	0	77	13	0
26	G	106	0	154	16	0
26	P	53	0	77	16	0
26	T	106	0	154	18	0
27	F	1	0	0	0	0
27	S	1	0	0	0	0
28	J	33	0	42	5	0
28	M	33	0	42	0	0
28	P	33	0	42	10	0
28	Z	33	0	42	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
29	A	297	0	0	17	0
29	B	274	0	0	13	0
29	C	176	0	0	8	0
29	D	266	0	0	5	0
29	E	178	0	0	2	0
29	F	199	0	0	9	0
29	G	100	0	0	7	0
29	H	122	0	0	8	0
29	I	88	0	0	2	0
29	J	63	0	0	2	0
29	K	69	0	0	4	0
29	L	48	0	0	2	0
29	M	47	0	0	0	0
29	N	290	0	0	12	0
29	O	243	0	0	4	0
29	P	173	0	0	8	0
29	Q	164	0	0	8	0
29	R	151	0	0	4	0
29	S	186	0	0	9	0
29	T	94	0	0	3	0
29	U	110	0	0	5	0
29	V	71	0	0	2	0
29	W	58	0	0	1	0
29	X	57	0	0	1	0
29	Y	40	0	0	3	0
29	Z	37	0	0	0	0
All	All	35054	0	31975	816	2

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 13.

The worst 5 of 816 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)
19:C:302:PGV:C21	19:C:302:PGV:C22	1.77	1.62
26:T:101:PEK:C3	26:T:101:PEK:C2	1.76	1.56
2:B:1:FME:CN	2:B:1:FME:N	1.70	1.53
3:P:224:LYS:CE	3:P:224:LYS:NZ	1.77	1.47
20:D:201:TGL:OG2	20:D:201:TGL:CB1	1.63	1.47

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:I:2:THR:CB	5:R:80:GLU:OE1[3_647]	2.00	0.20
9:I:2:THR:CG2	5:R:80:GLU:OE1[3_647]	2.19	0.01

## 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%)	515 (97%)	15 (3%)	0	100	100
1	N	528/514 (103%)	511 (97%)	17 (3%)	0	100	100
2	B	234/227 (103%)	230 (98%)	4 (2%)	0	100	100
2	O	230/227 (101%)	225 (98%)	4 (2%)	1 (0%)	34	13
3	C	266/261 (102%)	261 (98%)	4 (2%)	1 (0%)	34	13
3	P	266/261 (102%)	261 (98%)	5 (2%)	0	100	100
4	D	147/147 (100%)	143 (97%)	4 (3%)	0	100	100
4	Q	145/147 (99%)	139 (96%)	5 (3%)	1 (1%)	22	6
5	E	103/109 (94%)	102 (99%)	1 (1%)	0	100	100
5	R	104/109 (95%)	104 (100%)	0	0	100	100
6	F	100/98 (102%)	95 (95%)	3 (3%)	2 (2%)	7	1
6	S	98/98 (100%)	92 (94%)	2 (2%)	4 (4%)	3	0
7	G	82/85 (96%)	68 (83%)	8 (10%)	6 (7%)	1	0
7	T	82/85 (96%)	71 (87%)	5 (6%)	6 (7%)	1	0
8	H	77/85 (91%)	68 (88%)	6 (8%)	3 (4%)	3	0
8	U	77/85 (91%)	68 (88%)	5 (6%)	4 (5%)	2	0
9	I	71/73 (97%)	71 (100%)	0	0	100	100
9	V	71/73 (97%)	69 (97%)	1 (1%)	1 (1%)	11	1
10	J	56/59 (95%)	56 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
10	W	57/59 (97%)	57 (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%)	2 (5%)	0	100	100
13	Z	41/46 (89%)	39 (95%)	1 (2%)	1 (2%)	6	0
All	All	3590/3614 (99%)	3460 (96%)	100 (3%)	30 (1%)	17	5

5 of 30 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
6	F	94	HIS
6	F	96	LEU
7	G	4	ALA
7	G	5	LYS
7	G	8	HIS

### 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	444/426 (104%)	438 (99%)	6 (1%)	67	42
1	N	442/426 (104%)	432 (98%)	10 (2%)	50	20
2	B	219/210 (104%)	208 (95%)	11 (5%)	24	4
2	O	215/210 (102%)	206 (96%)	9 (4%)	30	6
3	C	233/226 (103%)	230 (99%)	3 (1%)	69	44
3	P	233/226 (103%)	230 (99%)	3 (1%)	69	44
4	D	133/129 (103%)	128 (96%)	5 (4%)	33	7
4	Q	131/129 (102%)	124 (95%)	7 (5%)	22	3
5	E	92/95 (97%)	90 (98%)	2 (2%)	52	22

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
5	R	93/95 (98%)	91 (98%)	2 (2%)	52	22
6	F	85/81 (105%)	79 (93%)	6 (7%)	14	1
6	S	83/81 (102%)	75 (90%)	8 (10%)	8	0
7	G	68/68 (100%)	63 (93%)	5 (7%)	13	1
7	T	68/68 (100%)	61 (90%)	7 (10%)	7	0
8	H	71/75 (95%)	65 (92%)	6 (8%)	10	1
8	U	71/75 (95%)	67 (94%)	4 (6%)	21	3
9	I	57/57 (100%)	55 (96%)	2 (4%)	36	9
9	V	57/57 (100%)	52 (91%)	5 (9%)	10	0
10	J	49/50 (98%)	49 (100%)	0	100	100
10	W	50/50 (100%)	48 (96%)	2 (4%)	31	6
11	K	39/46 (85%)	38 (97%)	1 (3%)	46	16
11	X	40/46 (87%)	40 (100%)	0	100	100
12	L	39/40 (98%)	38 (97%)	1 (3%)	46	16
12	Y	40/40 (100%)	38 (95%)	2 (5%)	24	4
13	M	37/38 (97%)	33 (89%)	4 (11%)	6	0
13	Z	37/38 (97%)	33 (89%)	4 (11%)	6	0
All	All	3126/3082 (101%)	3011 (96%)	115 (4%)	35	8

5 of 115 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	N	180	GLN
12	Y	20	ARG
3	P	159	MET
12	Y	2	HIS
8	U	9	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 39 such sidechains are listed below:

Mol	Chain	Res	Type
5	R	94	ASN
8	U	37	HIS
6	S	54	ASN
7	T	8	HIS

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Mol	Chain	Res	Type
10	W	29	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	N	1	1	8,9,10	1.31	1 (12%)	7,9,11	1.73	2 (28%)
2	FME	O	1	2	8,9,10	1.93	4 (50%)	7,9,11	2.34	3 (42%)
2	FME	B	1	2	8,9,10	5.84	6 (75%)	7,9,11	8.82	4 (57%)
1	FME	A	1	1	8,9,10	1.69	2 (25%)	7,9,11	1.64	3 (42%)
7	TPO	T	11	7	8,10,11	1.69	2 (25%)	10,14,16	1.24	1 (10%)
9	SAC	V	1	9	7,8,9	2.32	2 (28%)	8,9,11	3.04	4 (50%)
7	TPO	G	11	7	8,10,11	2.10	4 (50%)	10,14,16	1.99	2 (20%)
9	SAC	I	1	9	7,8,9	2.70	3 (42%)	8,9,11	1.44	1 (12%)

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	N	1	1	-	4/7/9/11	-
2	FME	O	1	2	-	0/7/9/11	-
2	FME	B	1	2	-	1/7/9/11	-
1	FME	A	1	1	-	3/7/9/11	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	TPO	T	11	7	-	4/9/11/13	-
9	SAC	V	1	9	-	4/7/8/10	-
7	TPO	G	11	7	-	5/9/11/13	-
9	SAC	I	1	9	-	1/7/8/10	-

The worst 5 of 24 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	FME	CN-N	11.15	1.71	1.33
2	B	1	FME	O1-CN	-8.94	0.95	1.22
9	I	1	SAC	OAC-C1A	5.36	1.35	1.23
9	V	1	SAC	OAC-C1A	4.57	1.33	1.23
2	B	1	FME	CB-CA	4.52	1.61	1.53

The worst 5 of 20 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	FME	O1-CN-N	-17.24	79.86	125.27
2	B	1	FME	CA-N-CN	-13.95	101.37	122.82
2	B	1	FME	CG-CB-CA	-5.63	97.31	112.95
7	G	11	TPO	CG2-CB-CA	5.16	123.35	113.16
9	V	1	SAC	C-CA-N	-5.02	100.67	109.73

There are no chirality outliers.

5 of 22 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	1	FME	N-CA-CB-CG
2	B	1	FME	O1-CN-N-CA
7	G	11	TPO	N-CA-CB-CG2
7	G	11	TPO	N-CA-CB-OG1
7	G	11	TPO	C-CA-CB-CG2

There are no ring outliers.

3 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1	FME	7	0
7	T	11	TPO	3	0
7	G	11	TPO	4	0

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 56 ligands modelled in this entry, 8 are monoatomic and 2 are unknown - leaving 46 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	$\# Z  > 2$
18	PER	A	606	15,14	0,1,1	-	-	-		
25	CDL	G	102	-	99,99,99	1.49	13 (13%)	105,111,111	1.82	22 (20%)
26	PEK	P	308	-	52,52,52	1.68	6 (11%)	55,57,57	1.92	14 (25%)
22	CHD	C	305	-	32,32,32	2.46	15 (46%)	51,51,51	2.68	20 (39%)
20	TGL	B	301	-	62,62,62	1.71	8 (12%)	65,65,65	2.82	21 (32%)
22	CHD	P	307	-	32,32,32	1.98	13 (40%)	51,51,51	2.47	20 (39%)
21	CUA	B	302	2	0,1,1	-	-	-		
20	TGL	D	201	-	62,62,62	2.77	11 (17%)	65,65,65	3.36	24 (36%)
19	PGV	Q	201	-	50,50,50	1.31	3 (6%)	53,56,56	1.74	10 (18%)
26	PEK	G	103	-	52,52,52	1.29	3 (5%)	55,57,57	1.64	11 (20%)
19	PGV	C	307	-	50,50,50	1.29	4 (8%)	53,56,56	1.60	8 (15%)
25	CDL	C	303	-	99,99,99	1.82	20 (20%)	105,111,111	2.30	34 (32%)
22	CHD	O	302	-	32,32,32	2.42	11 (34%)	51,51,51	2.45	21 (41%)
19	PGV	A	608	-	50,50,50	1.97	6 (12%)	53,56,56	2.43	14 (26%)
19	PGV	A	607	-	50,50,50	1.49	9 (18%)	53,56,56	1.54	11 (20%)
20	TGL	Q	202	-	62,62,62	2.36	11 (17%)	65,65,65	2.77	14 (21%)
26	PEK	T	102	-	52,52,52	1.30	2 (3%)	55,57,57	1.55	8 (14%)
21	CUA	O	301	2	0,1,1	-	-	-		
19	PGV	C	302	-	50,50,50	1.41	5 (10%)	53,56,56	2.06	5 (9%)
22	CHD	P	305	-	32,32,32	1.53	6 (18%)	51,51,51	3.52	29 (56%)
26	PEK	C	306	-	52,52,52	1.61	5 (9%)	55,57,57	2.21	18 (32%)
14	HEA	A	602	18,1	57,67,67	1.74	12 (21%)	61,103,103	2.14	17 (27%)



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
19	PGV	P	301	-	50,50,50	1.11	2 (4%)	53,56,56	1.68	10 (18%)
22	CHD	C	304	-	32,32,32	1.48	5 (15%)	51,51,51	3.85	30 (58%)
14	HEA	N	601	1	57,67,67	2.31	20 (35%)	61,103,103	2.62	20 (32%)
28	DMU	J	101	-	34,34,34	1.00	1 (2%)	45,45,45	1.48	6 (13%)
18	PER	N	606	15,14	0,1,1	-	-	-	-	-
20	TGL	Y	101	-	62,62,62	2.05	11 (17%)	65,65,65	3.06	27 (41%)
22	CHD	B	303	-	32,32,32	2.75	18 (56%)	51,51,51	2.65	23 (45%)
28	DMU	M	101	-	34,34,34	1.49	5 (14%)	45,45,45	2.24	16 (35%)
19	PGV	P	303	-	50,50,50	1.06	3 (6%)	53,56,56	1.51	13 (24%)
28	DMU	P	306	-	34,34,34	0.87	1 (2%)	45,45,45	1.80	9 (20%)
14	HEA	A	601	1	57,67,67	2.33	18 (31%)	61,103,103	2.52	22 (36%)
23	PSC	O	303	-	51,51,51	1.45	3 (5%)	57,59,59	1.73	11 (19%)
25	CDL	T	103	-	99,99,99	1.53	12 (12%)	105,111,111	1.67	18 (17%)
25	CDL	P	304	-	99,99,99	2.05	22 (22%)	105,111,111	2.17	35 (33%)
19	PGV	N	607	-	50,50,50	1.47	7 (14%)	53,56,56	1.50	11 (20%)
26	PEK	G	101	-	52,52,52	1.20	5 (9%)	55,57,57	1.66	12 (21%)
23	PSC	B	304	-	51,51,51	1.32	3 (5%)	57,59,59	1.77	12 (21%)
14	HEA	N	602	18,1	57,67,67	1.72	15 (26%)	61,103,103	2.70	29 (47%)
26	PEK	T	101	-	52,52,52	1.60	4 (7%)	55,57,57	2.28	10 (18%)
20	TGL	L	101	-	62,62,62	2.10	13 (20%)	65,65,65	2.84	27 (41%)
22	CHD	W	101	-	32,32,32	2.03	13 (40%)	51,51,51	4.55	31 (60%)
22	CHD	J	102	-	32,32,32	1.86	9 (28%)	51,51,51	4.61	35 (68%)
28	DMU	Z	101	-	34,34,34	1.26	5 (14%)	45,45,45	1.84	12 (26%)
20	TGL	N	608	-	62,62,62	1.44	8 (12%)	65,65,65	2.18	14 (21%)

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
25	CDL	G	102	-	-	59/110/110/110	-
26	PEK	P	308	-	-	22/56/56/56	-
22	CHD	C	305	-	-	2/9/74/74	0/4/4/4
20	TGL	B	301	-	-	31/65/65/65	-
22	CHD	P	307	-	-	3/9/74/74	0/4/4/4

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
20	TGL	D	201	-	-	33/65/65/65	-
19	PGV	Q	201	-	-	31/55/55/55	-
26	PEK	G	103	-	-	22/56/56/56	-
19	PGV	C	307	-	-	33/55/55/55	-
25	CDL	C	303	-	-	57/110/110/110	-
22	CHD	O	302	-	-	2/9/74/74	0/4/4/4
19	PGV	A	608	-	-	34/55/55/55	-
19	PGV	A	607	-	-	8/55/55/55	-
20	TGL	Q	202	-	-	36/65/65/65	-
26	PEK	T	102	-	-	28/56/56/56	-
19	PGV	C	302	-	-	14/55/55/55	-
22	CHD	P	305	-	-	7/9/74/74	0/4/4/4
26	PEK	C	306	-	-	34/56/56/56	-
14	HEA	A	602	18,1	-	4/32/76/76	-
19	PGV	P	301	-	-	30/55/55/55	-
22	CHD	C	304	-	-	8/9/74/74	0/4/4/4
14	HEA	N	601	1	-	6/32/76/76	-
28	DMU	J	101	-	-	6/19/59/59	0/2/2/2
20	TGL	Y	101	-	-	41/65/65/65	-
22	CHD	B	303	-	-	2/9/74/74	0/4/4/4
28	DMU	M	101	-	-	5/19/59/59	0/2/2/2
19	PGV	P	303	-	-	10/55/55/55	-
28	DMU	P	306	-	-	6/19/59/59	0/2/2/2
14	HEA	A	601	1	-	5/32/76/76	-
23	PSC	O	303	-	-	27/55/55/55	-
25	CDL	T	103	-	-	62/110/110/110	-
25	CDL	P	304	-	-	62/110/110/110	-
19	PGV	N	607	-	-	9/55/55/55	-
26	PEK	G	101	-	-	16/56/56/56	-
23	PSC	B	304	-	-	31/55/55/55	-
14	HEA	N	602	18,1	-	4/32/76/76	-
26	PEK	T	101	-	-	21/56/56/56	-
20	TGL	L	101	-	-	34/65/65/65	-
22	CHD	W	101	-	-	9/9/74/74	0/4/4/4

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
22	CHD	J	102	-	-	7/9/74/74	0/4/4/4
28	DMU	Z	101	-	-	5/19/59/59	0/2/2/2
20	TGL	N	608	-	-	43/65/65/65	-

The worst 5 of 366 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
20	D	201	TGL	OB1-CB1	12.98	1.61	1.22
20	Q	202	TGL	OB1-CB1	11.24	1.55	1.22
20	D	201	TGL	OG2-CB1	10.44	1.63	1.34
20	Q	202	TGL	OG2-CB1	9.46	1.60	1.34
14	A	601	HEA	C18-C19	-8.37	1.12	1.33

The worst 5 of 754 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
22	W	101	CHD	C17-C13-C12	17.08	133.26	117.67
20	D	201	TGL	OG2-CB1-CB2	-15.65	77.78	111.50
22	W	101	CHD	C18-C13-C12	-13.78	95.04	109.07
22	J	102	CHD	C17-C13-C12	13.68	130.16	117.67
20	D	201	TGL	OG2-CB1-OB1	13.55	156.45	123.70

There are no chirality outliers.

5 of 909 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
19	A	608	PGV	C03-O11-P-O13
19	A	608	PGV	C03-O11-P-O14
19	A	608	PGV	C04-O12-P-O13
19	A	608	PGV	C04-O12-P-O14
19	A	608	PGV	O02-C1-O01-C02

There are no ring outliers.

40 monomers are involved in 372 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
18	A	606	PER	1	0
25	G	102	CDL	37	0
26	P	308	PEK	16	0
22	C	305	CHD	1	0

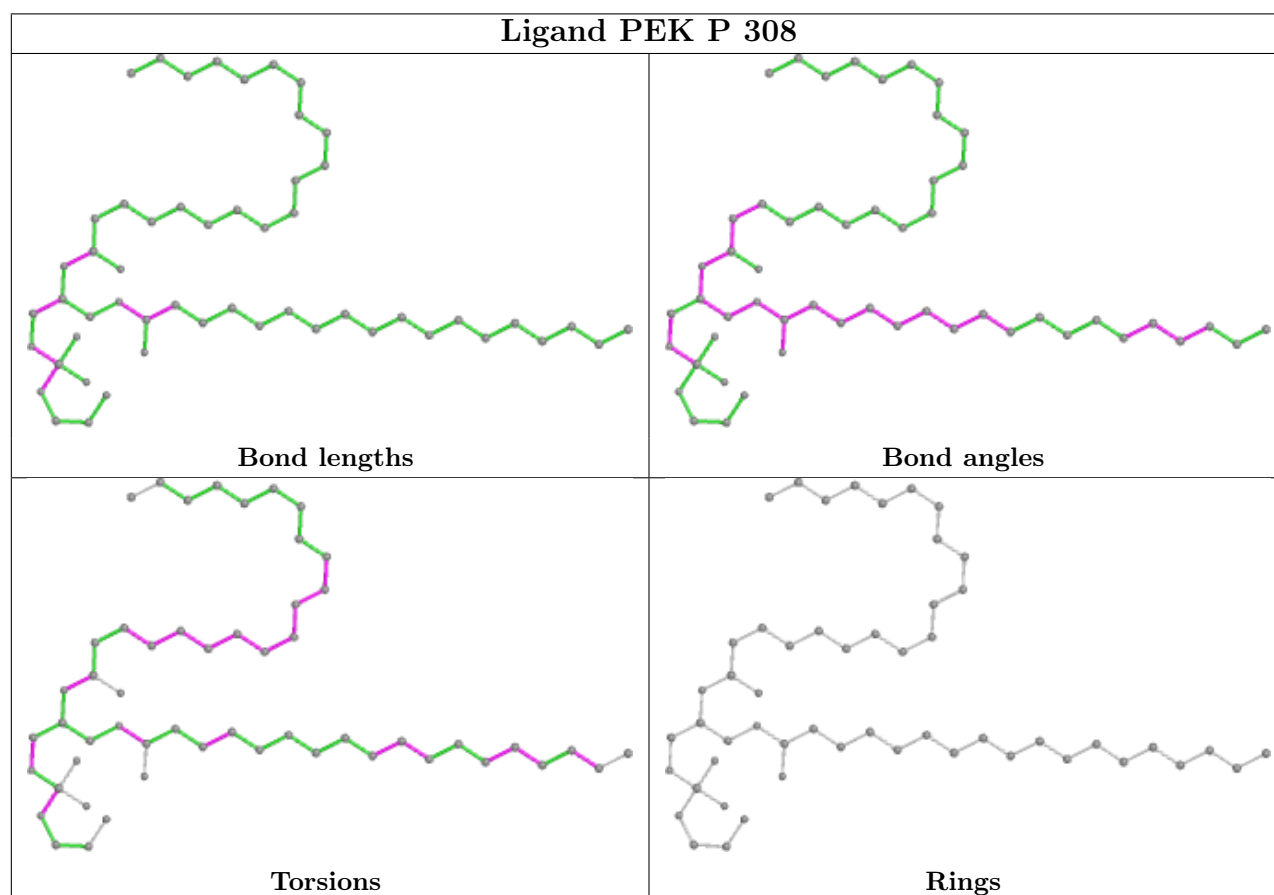
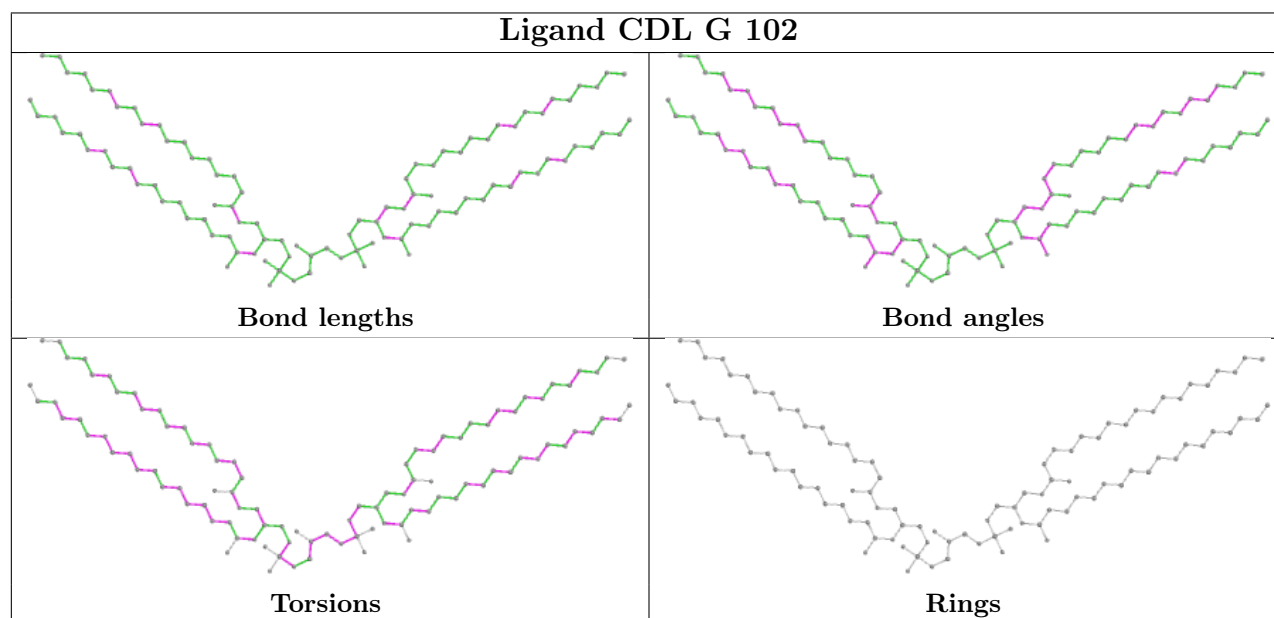
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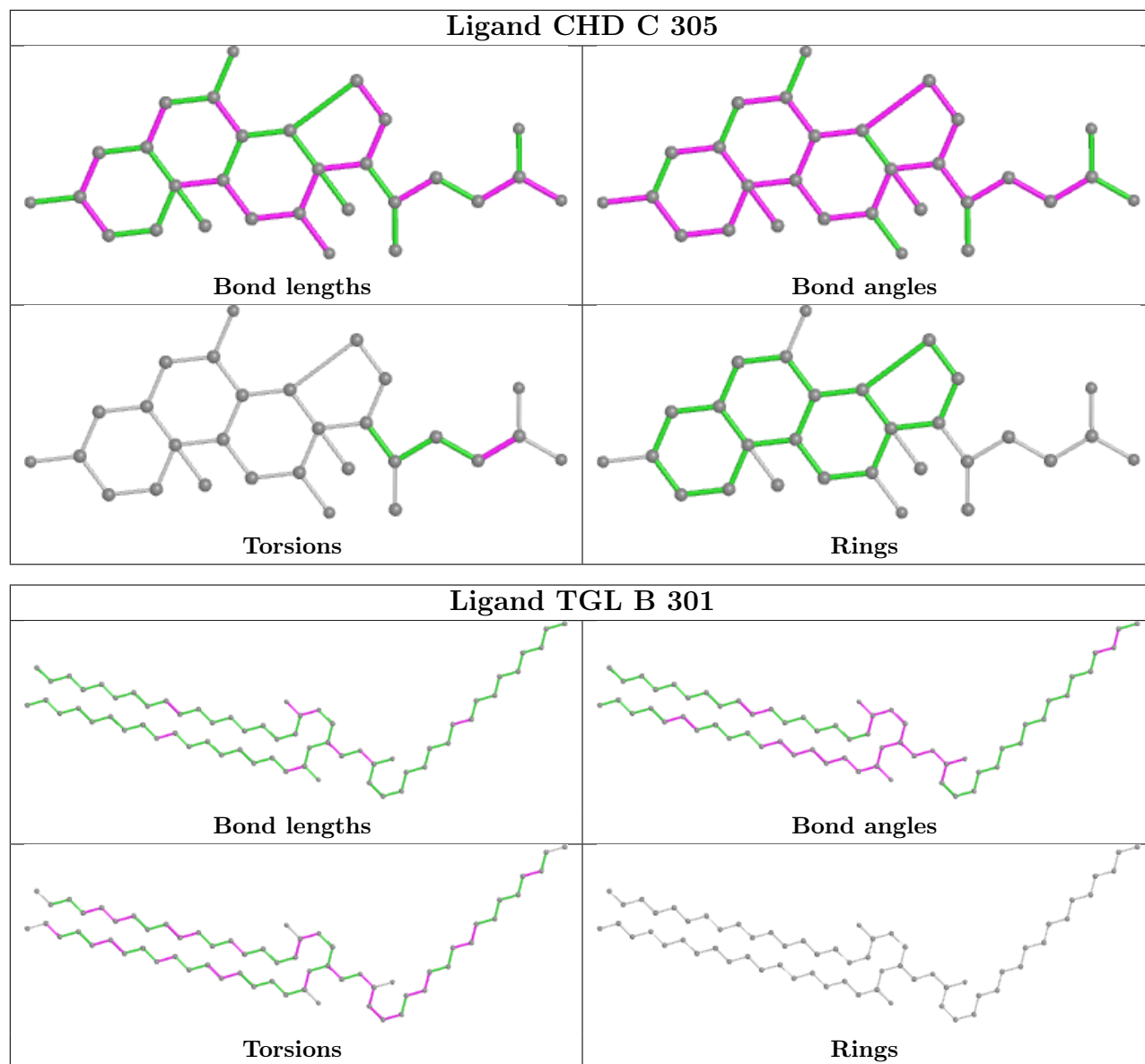
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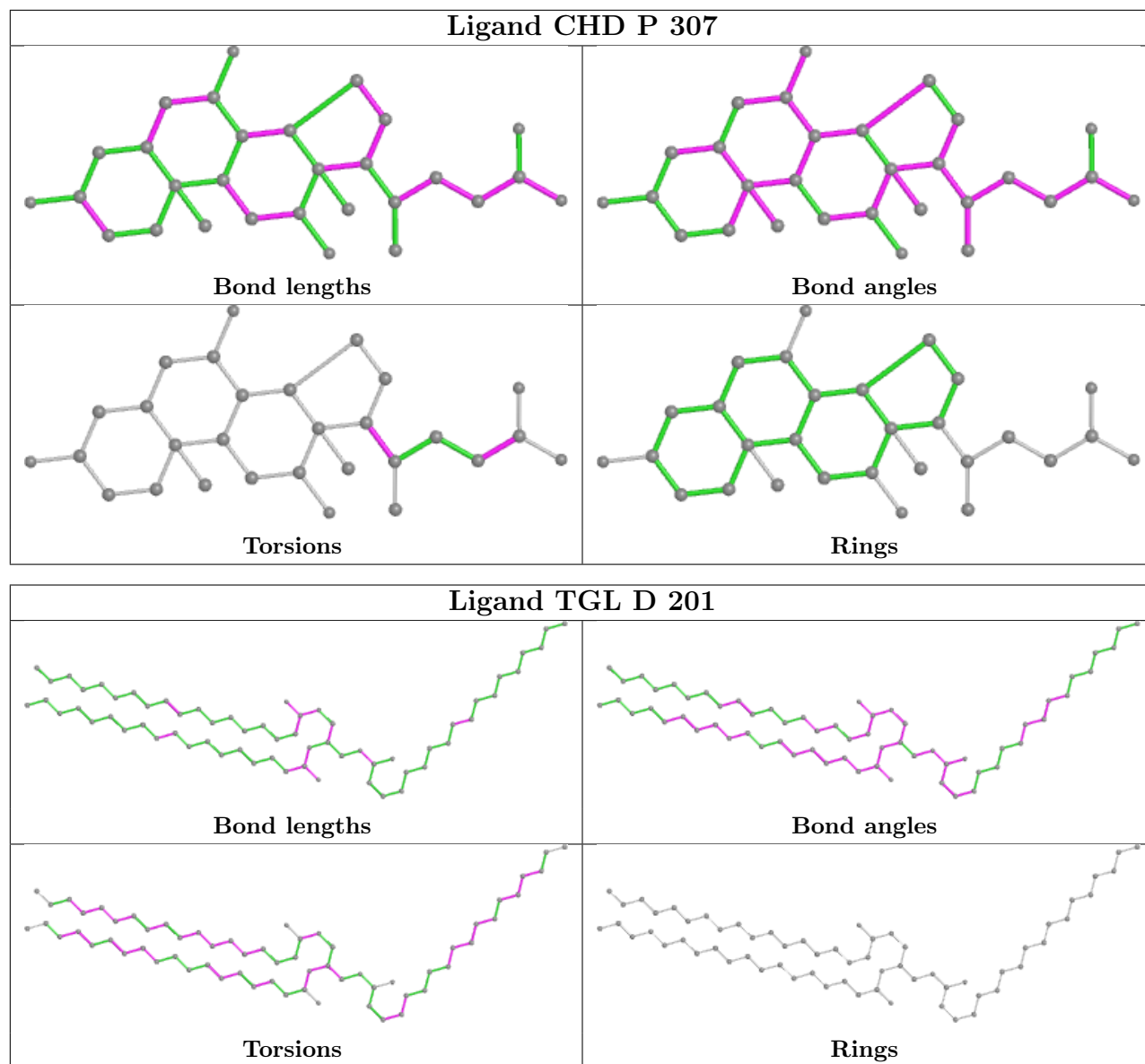
Mol	Chain	Res	Type	Clashes	Symm-Clashes
20	B	301	TGL	3	0
22	P	307	CHD	1	0
20	D	201	TGL	21	0
19	Q	201	PGV	11	0
26	G	103	PEK	9	0
19	C	307	PGV	1	0
25	C	303	CDL	26	0
19	A	608	PGV	7	0
19	A	607	PGV	4	0
20	Q	202	TGL	15	0
26	T	102	PEK	10	0
19	C	302	PGV	7	0
22	P	305	CHD	6	0
26	C	306	PEK	13	0
14	A	602	HEA	1	0
19	P	301	PGV	3	0
22	C	304	CHD	2	0
14	N	601	HEA	10	0
28	J	101	DMU	5	0
18	N	606	PER	1	0
20	Y	101	TGL	23	0
19	P	303	PGV	6	0
28	P	306	DMU	10	0
14	A	601	HEA	9	0
23	O	303	PSC	19	0
25	T	103	CDL	28	0
25	P	304	CDL	24	0
19	N	607	PGV	1	0
26	G	101	PEK	7	0
23	B	304	PSC	13	0
14	N	602	HEA	1	0
26	T	101	PEK	8	0
20	L	101	TGL	19	0
22	W	101	CHD	6	0
22	J	102	CHD	6	0
20	N	608	TGL	6	0

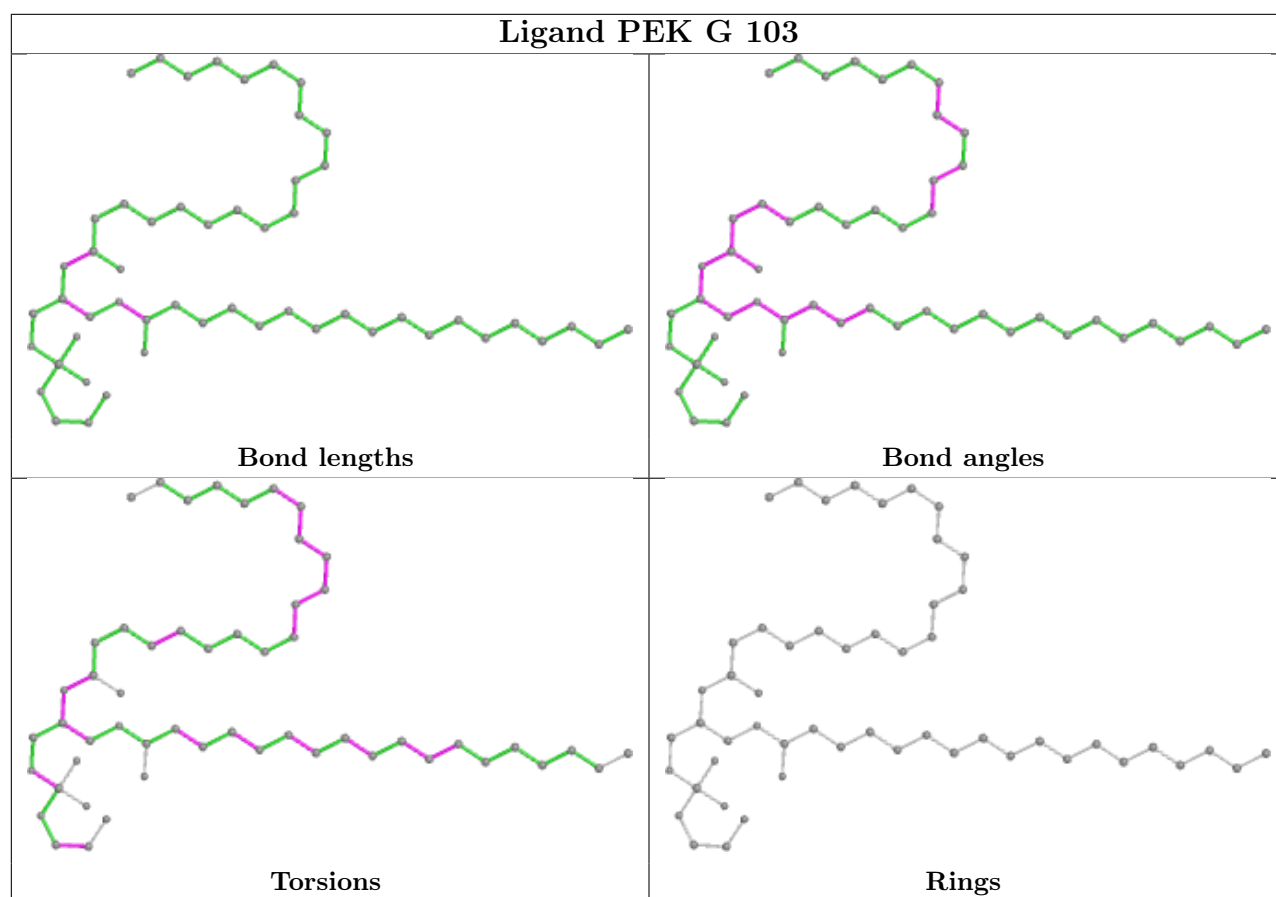
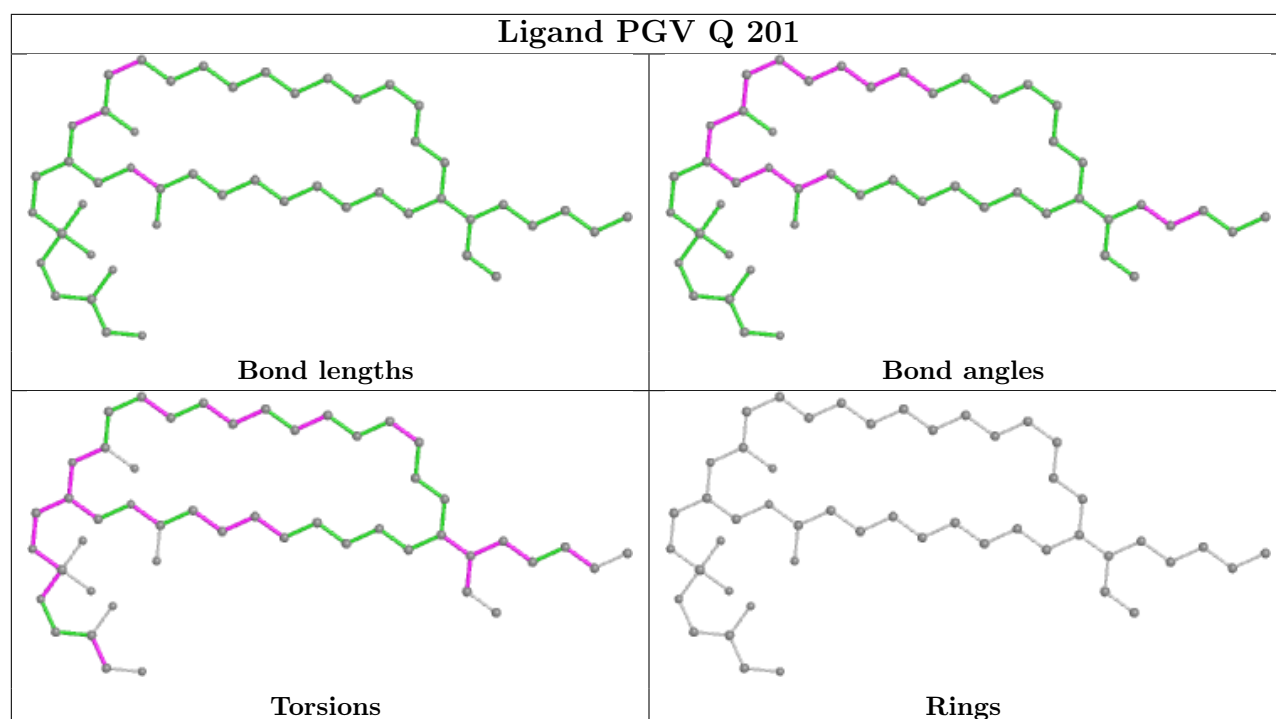
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be

highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

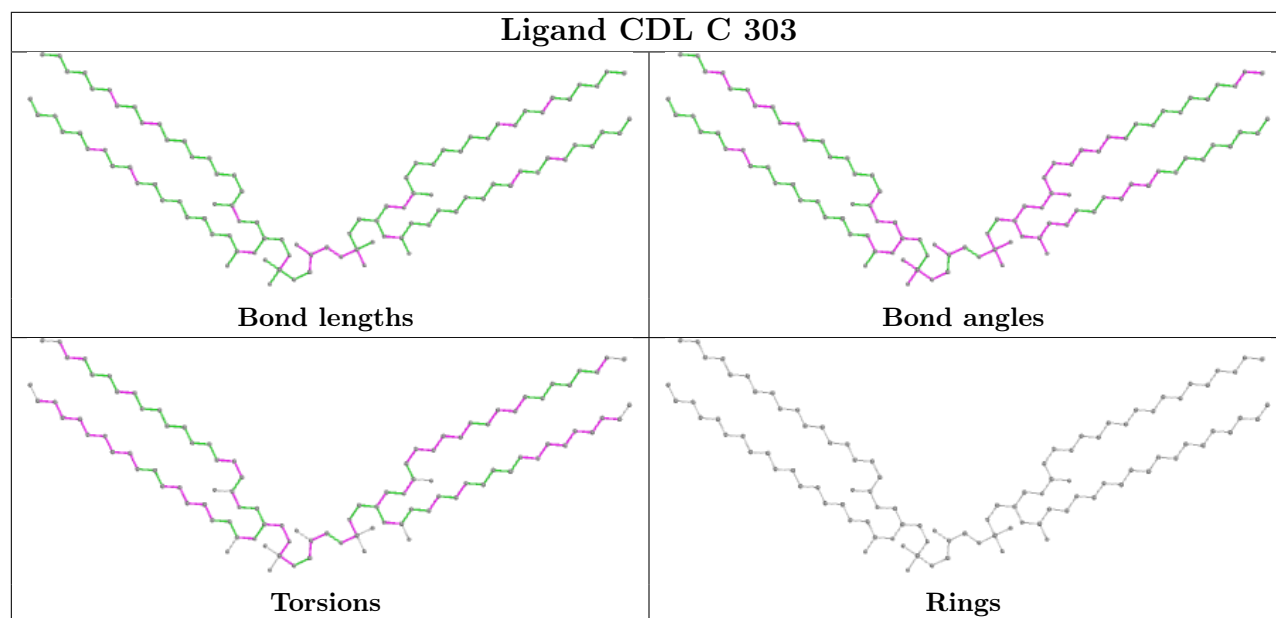
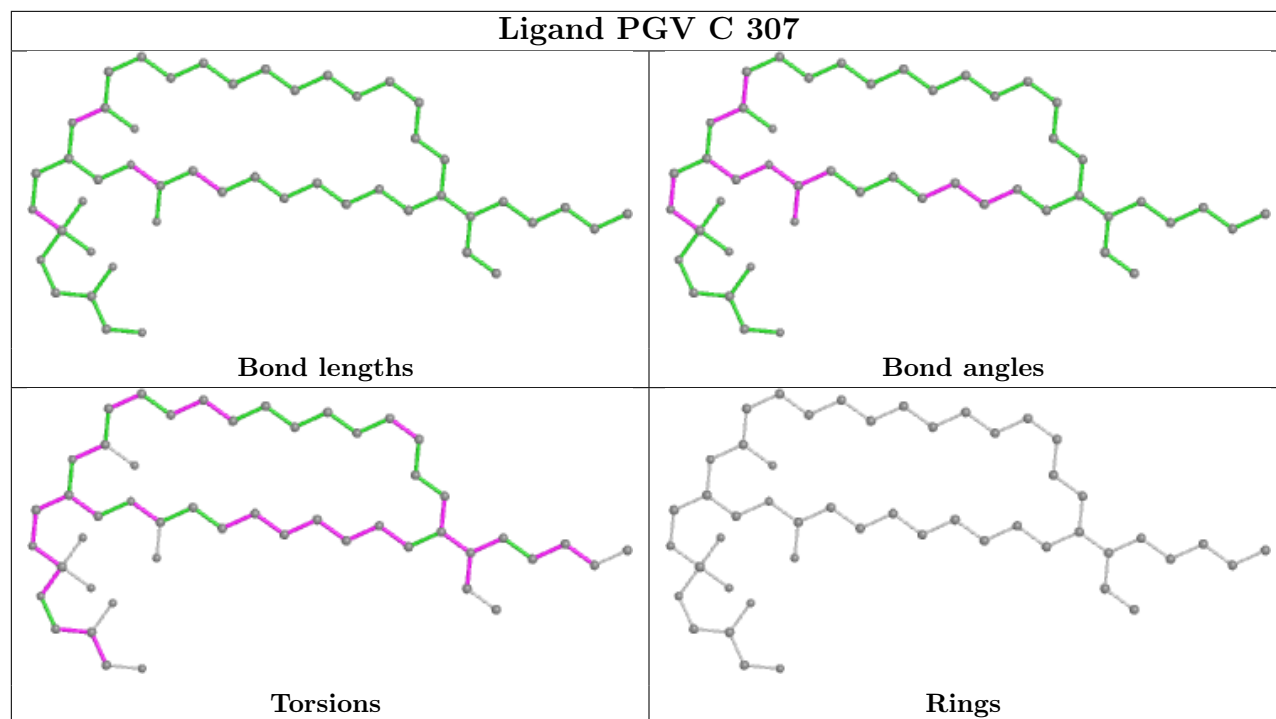


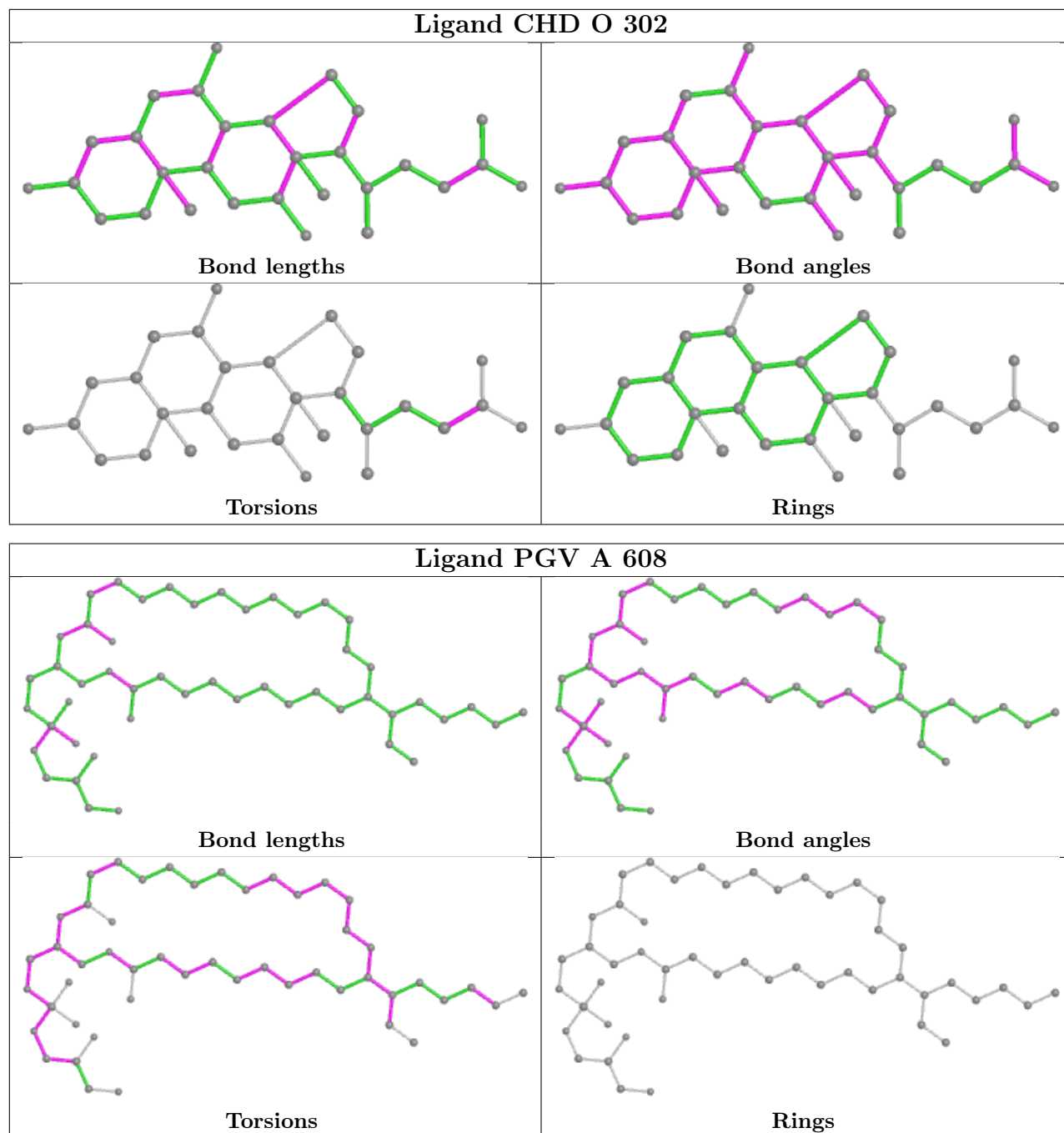


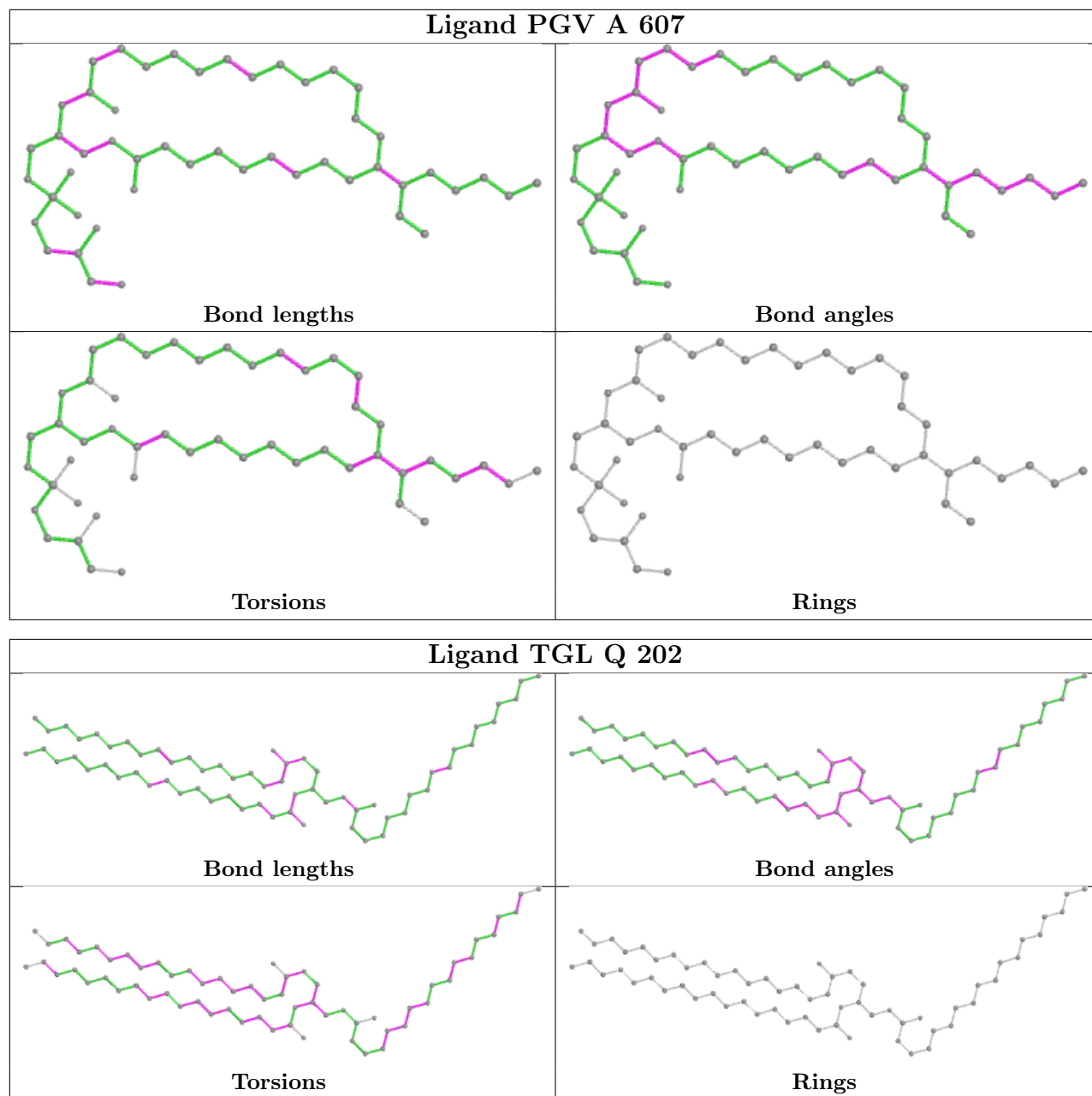


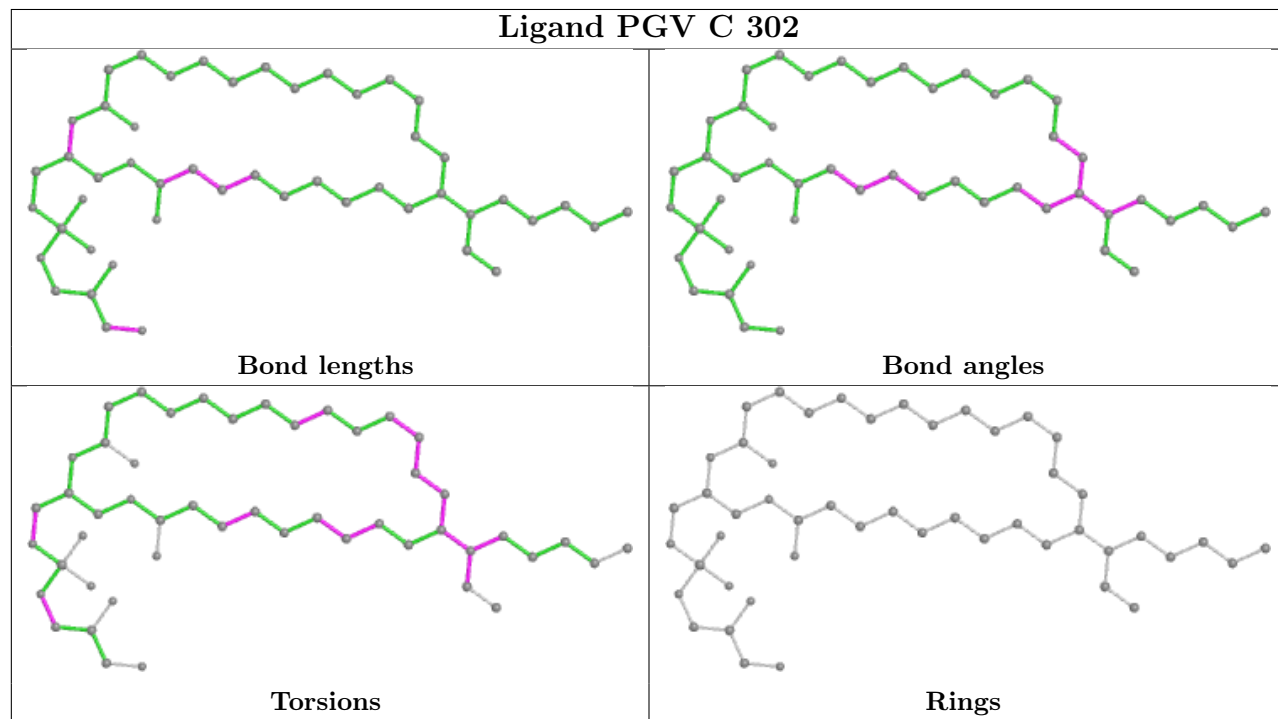
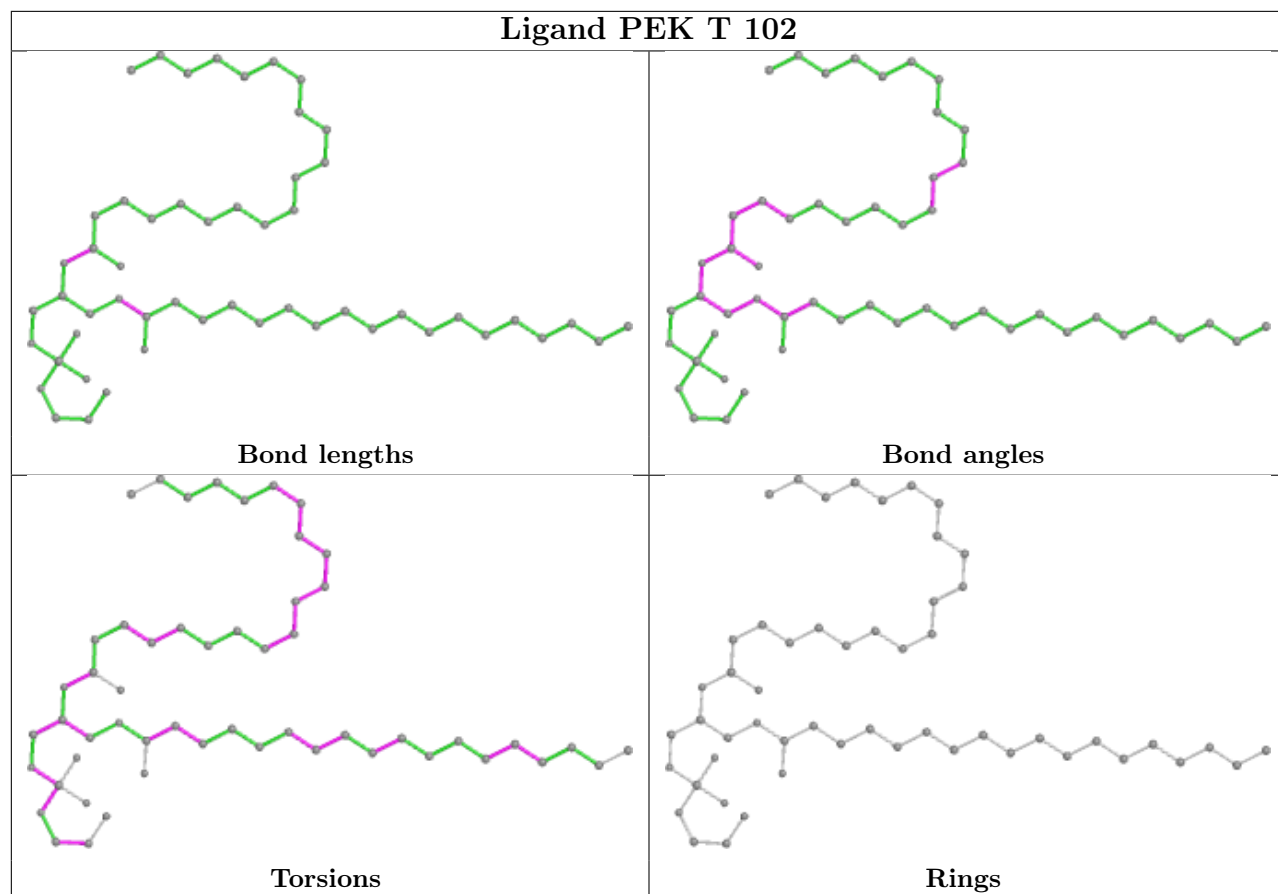


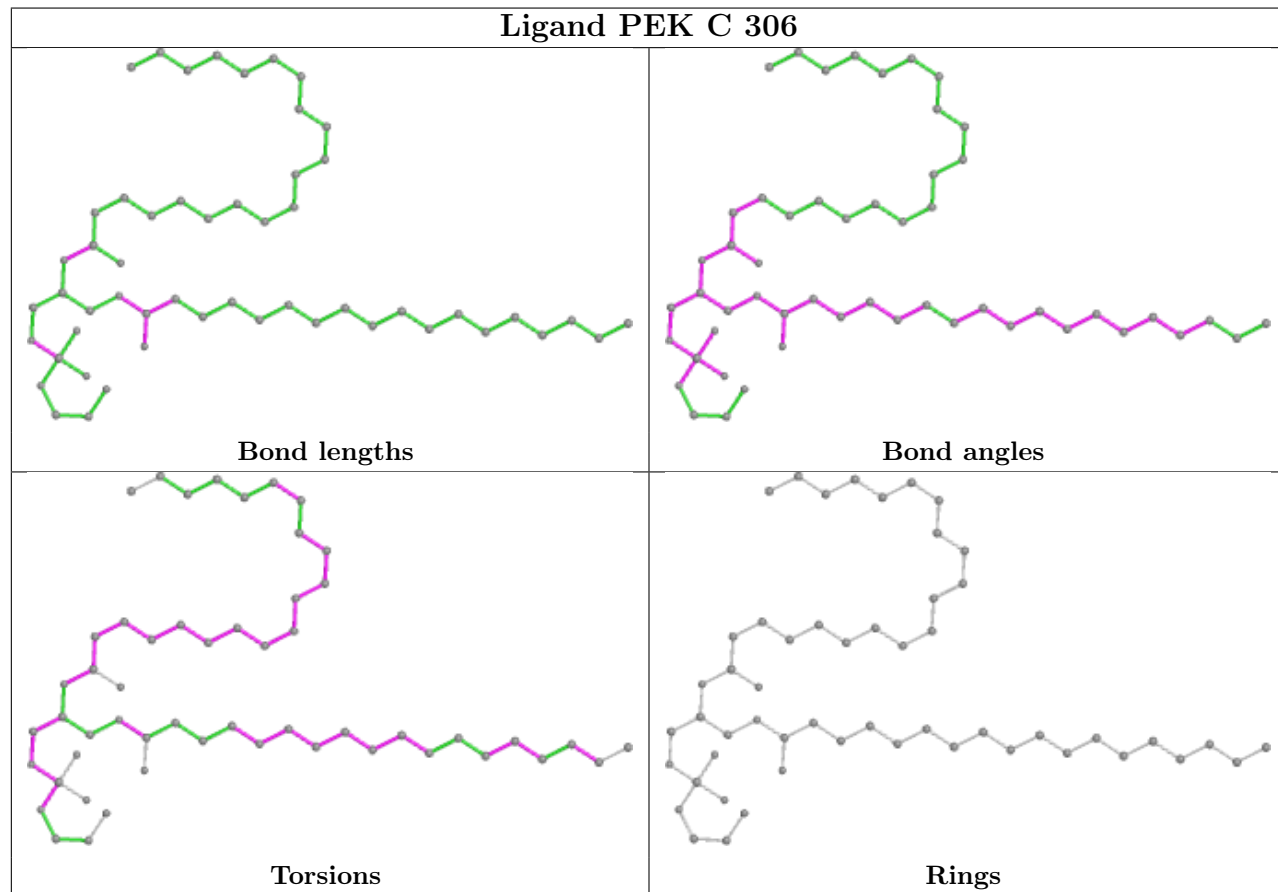
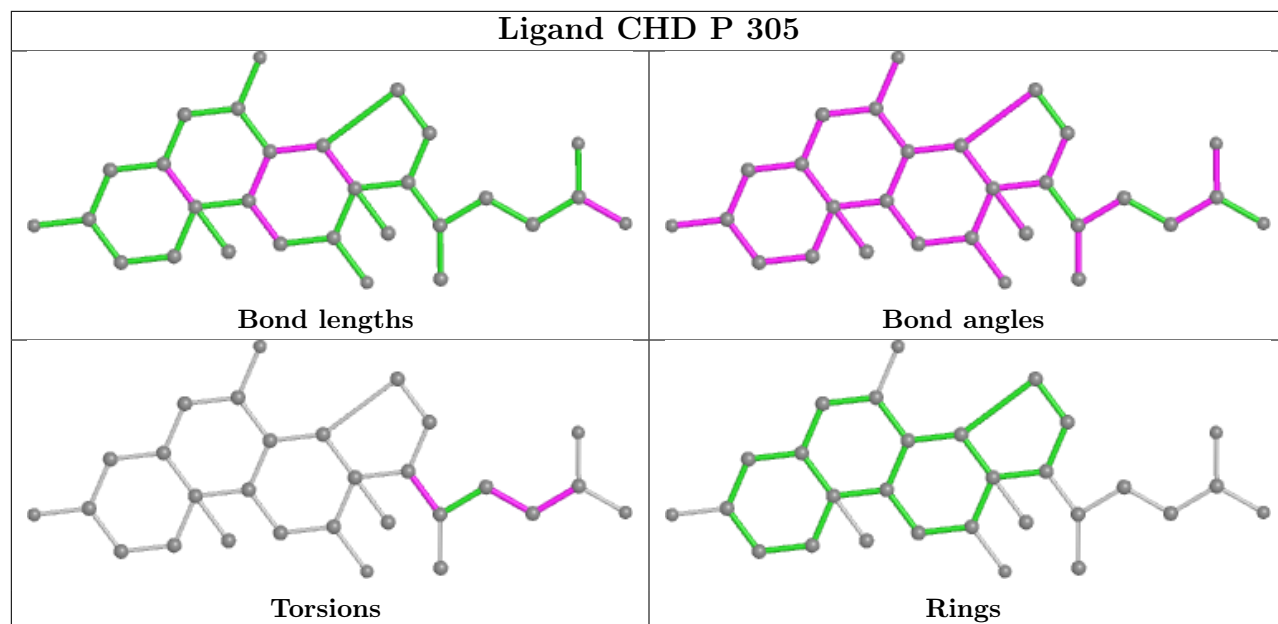


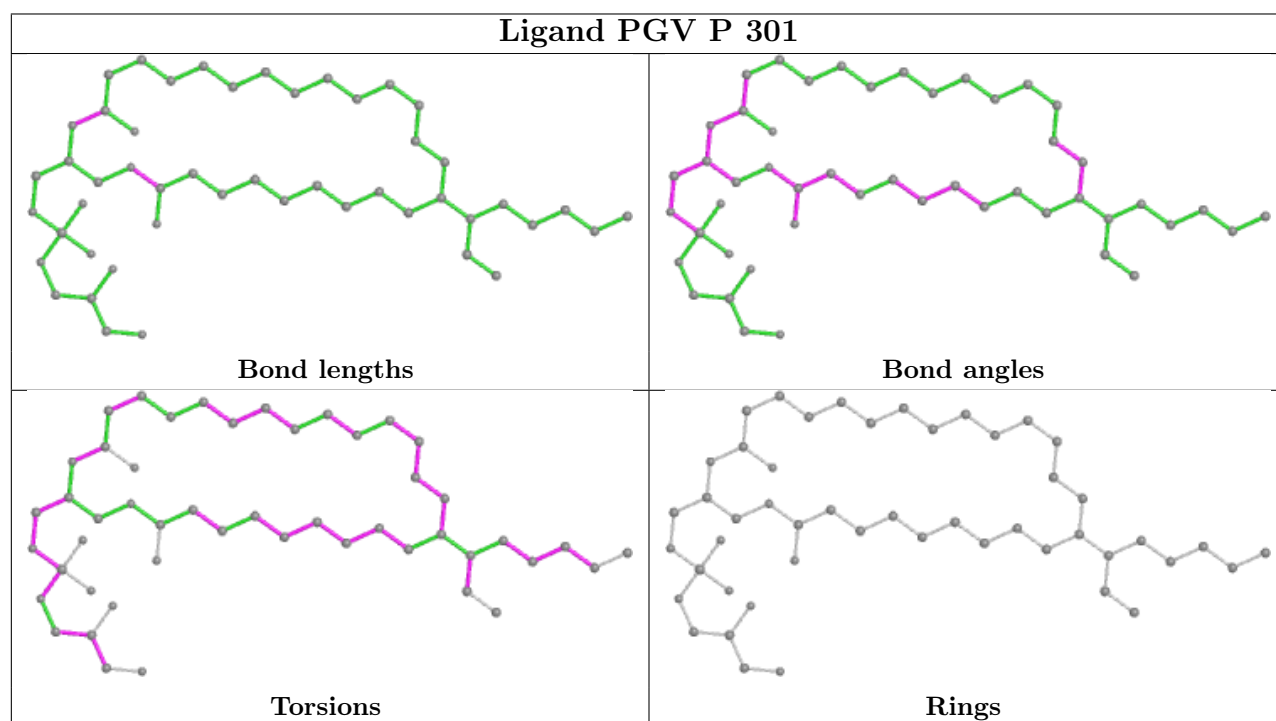
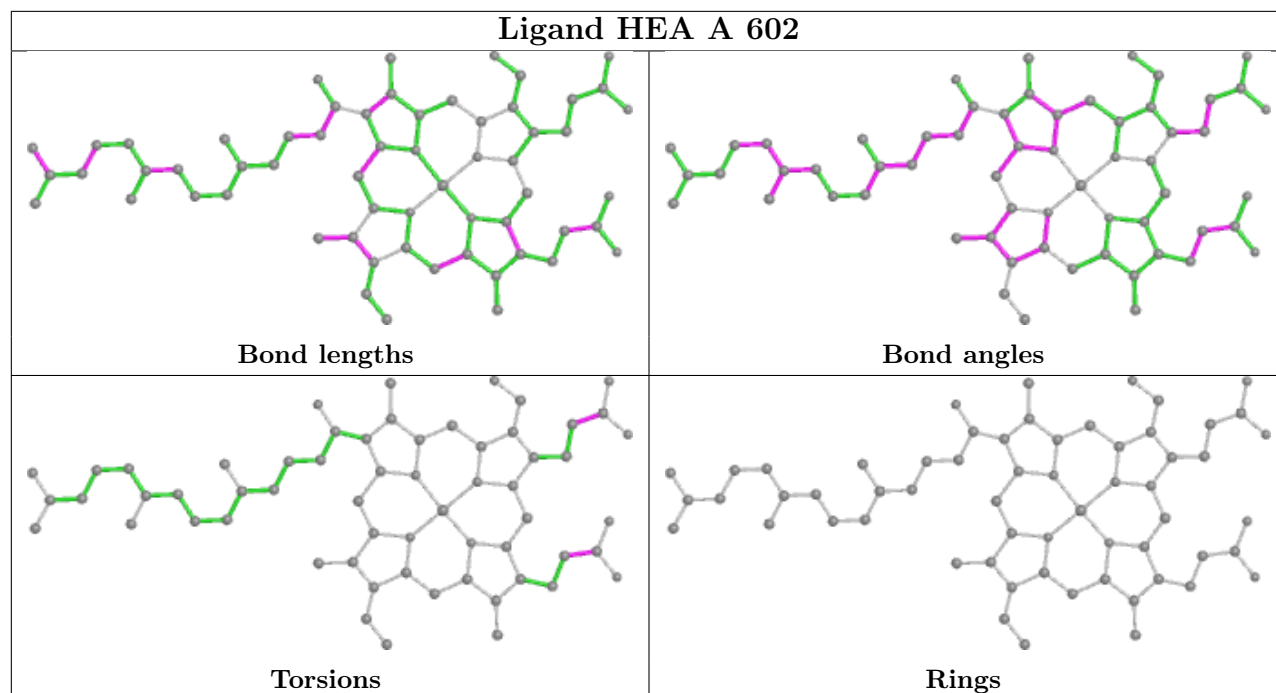


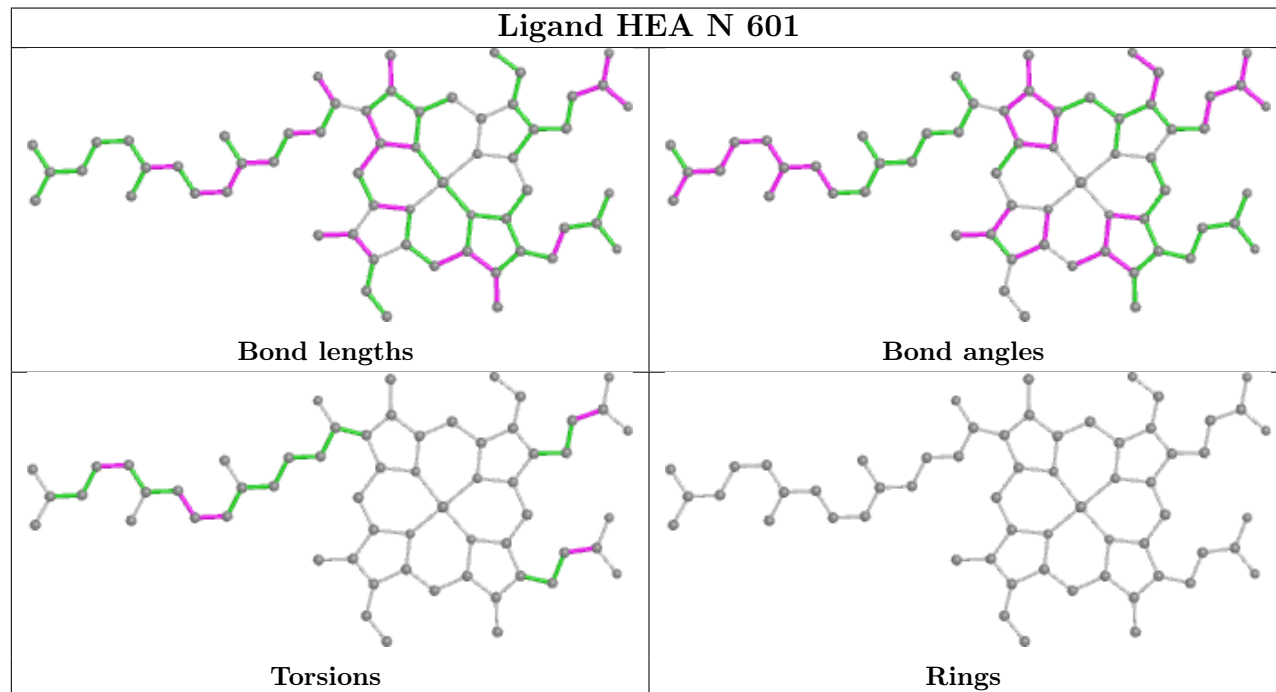
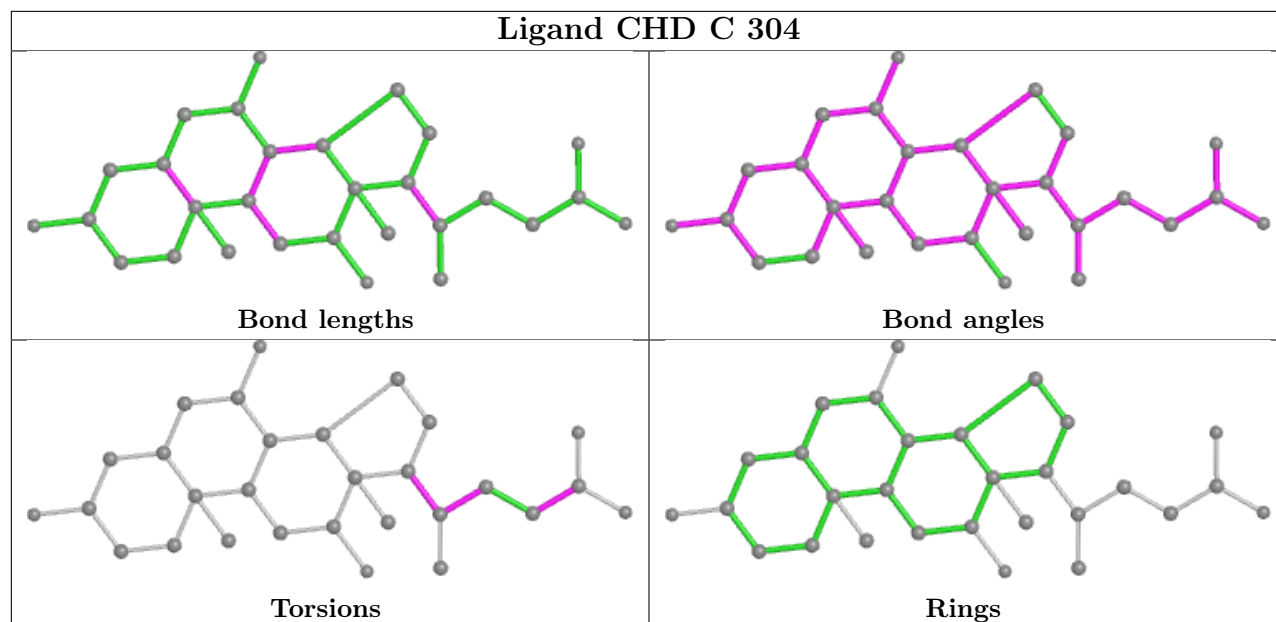


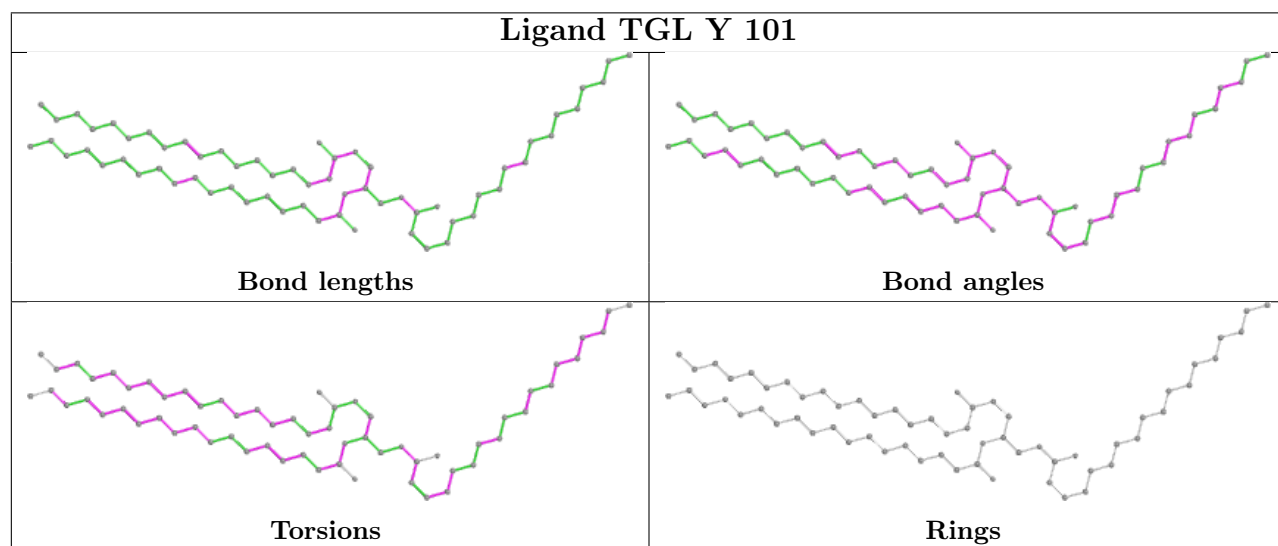
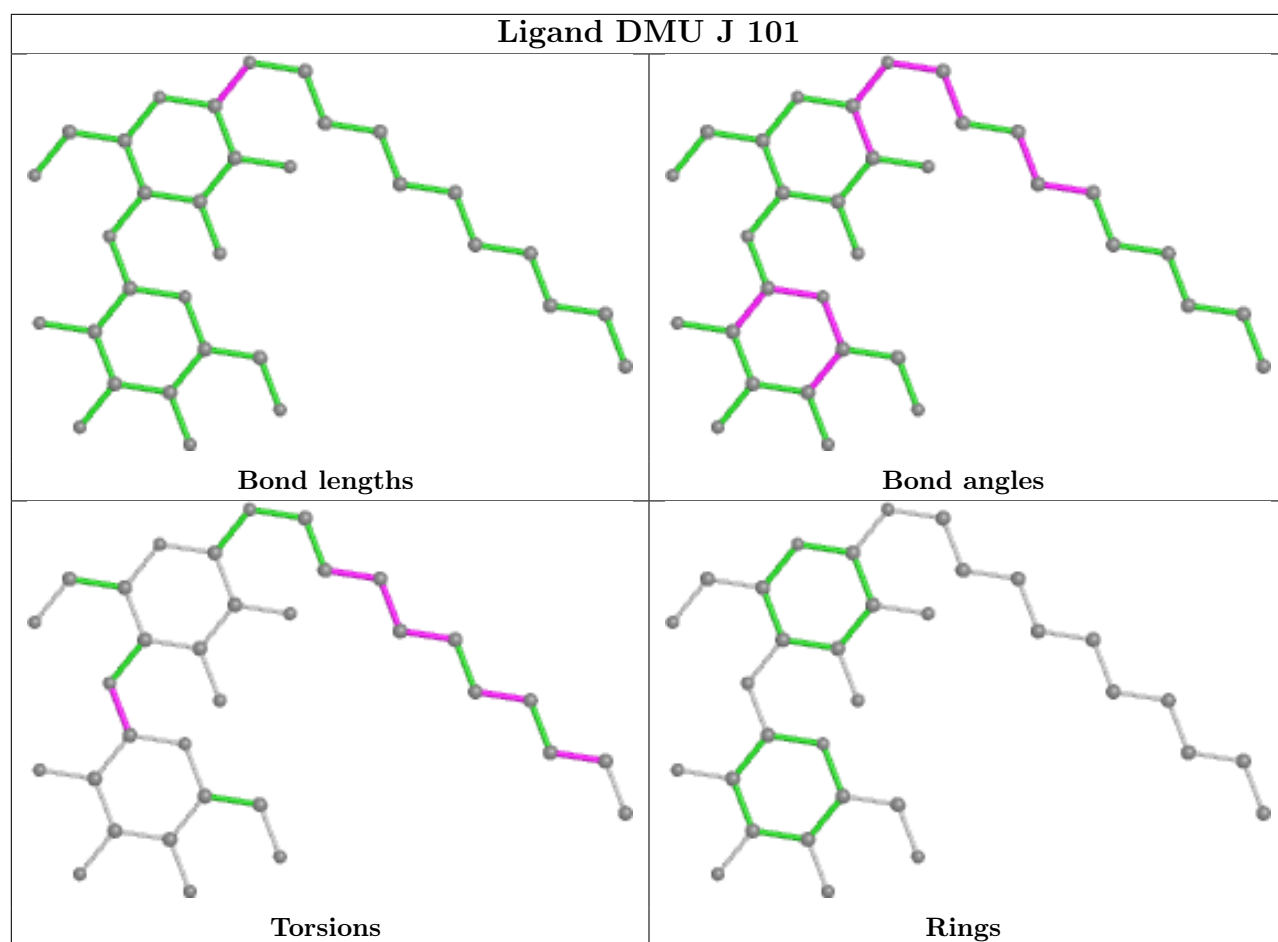




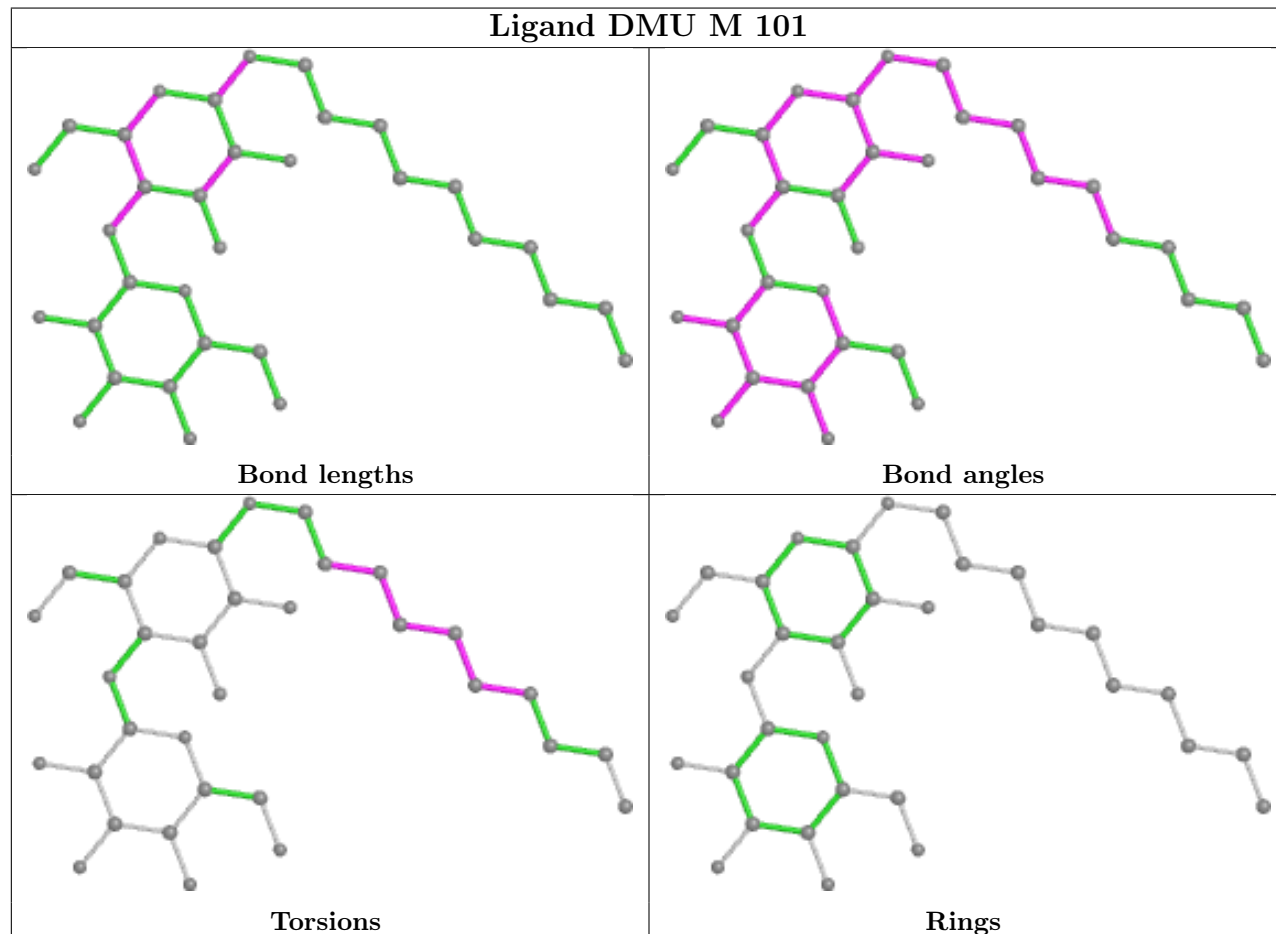
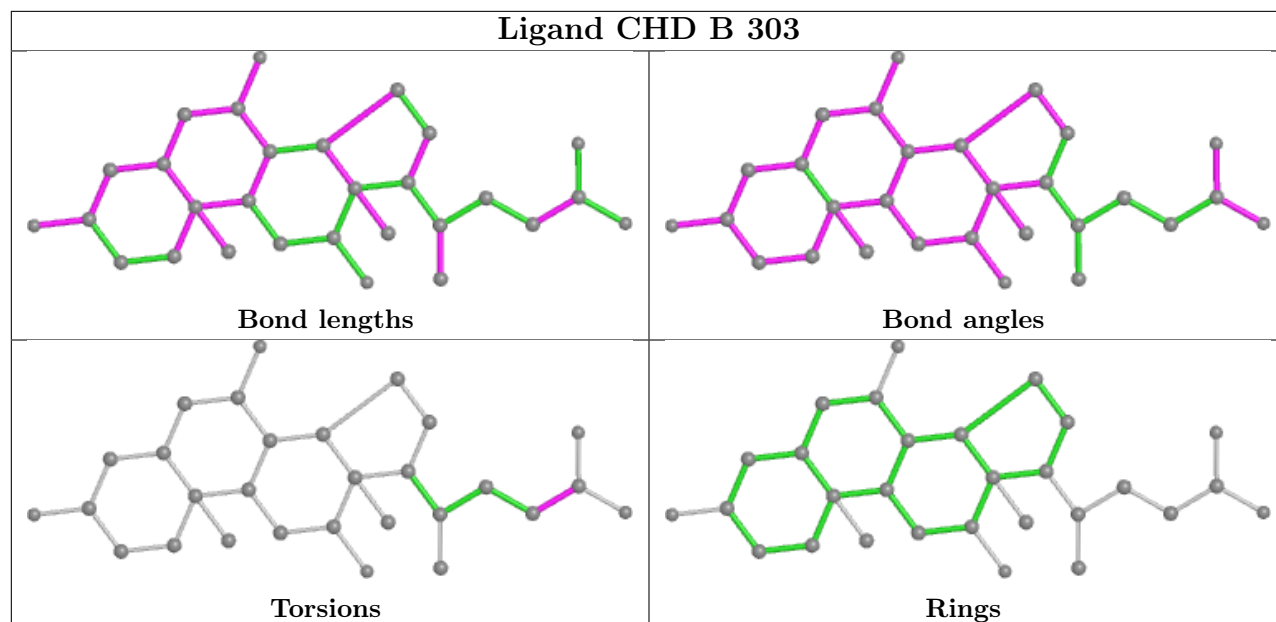


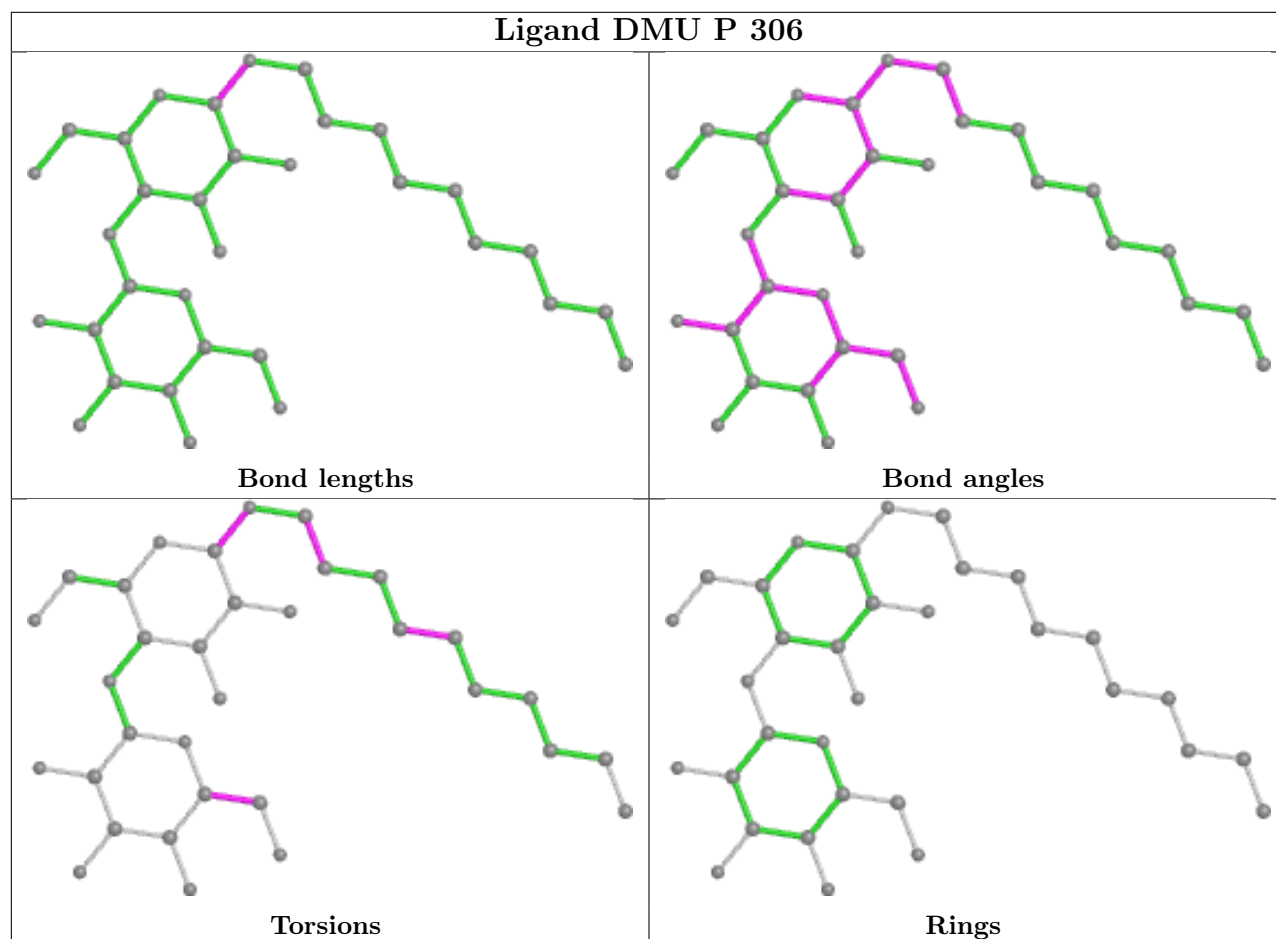
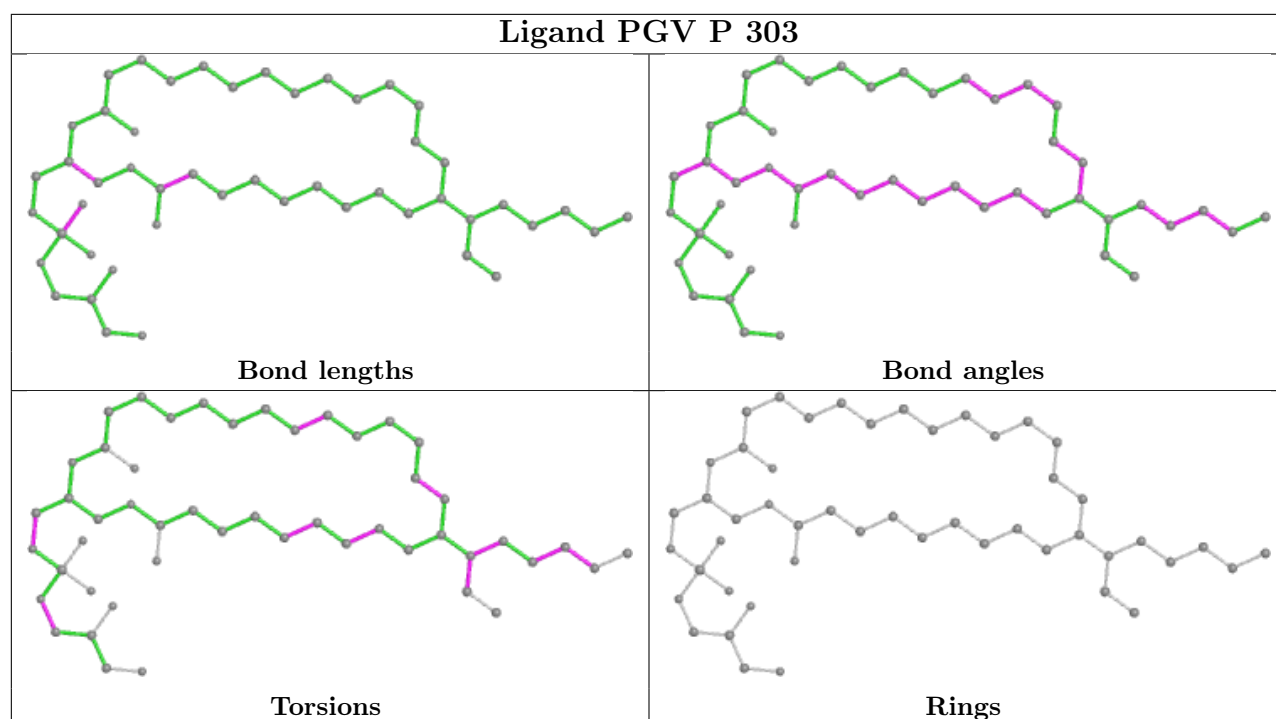


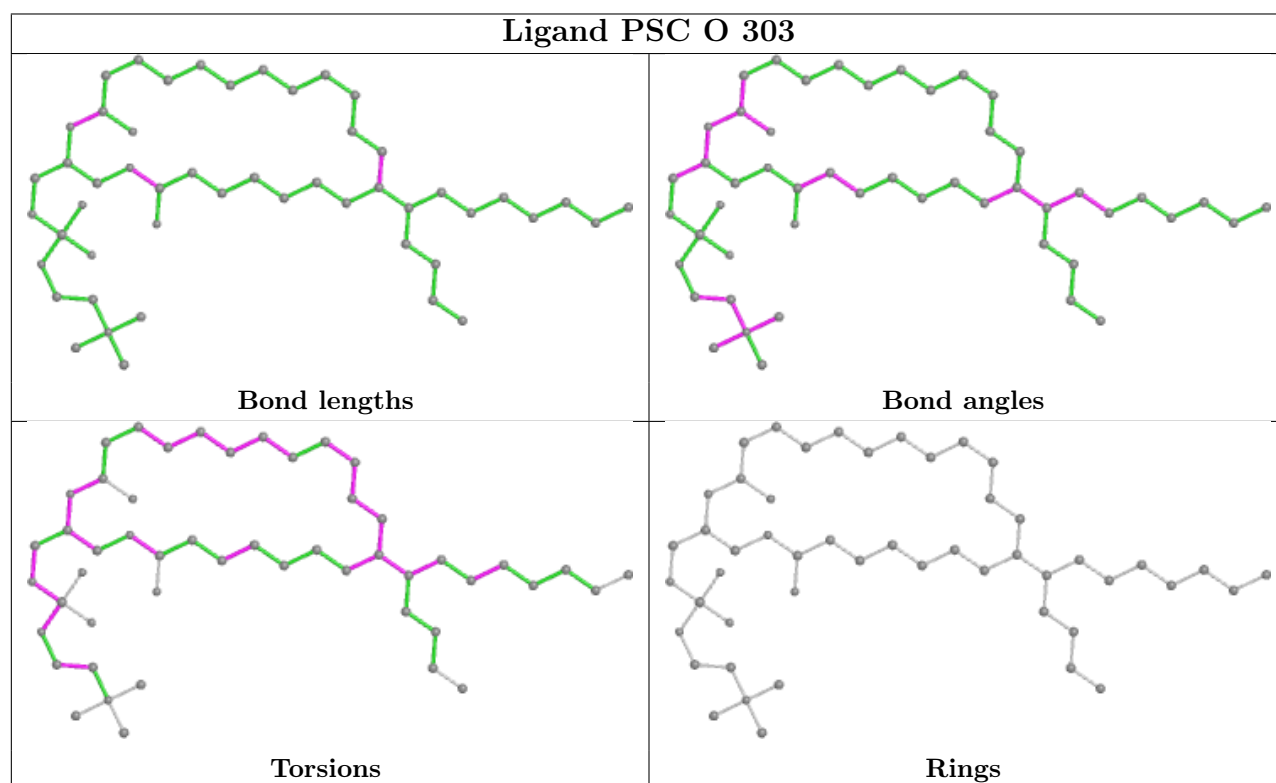
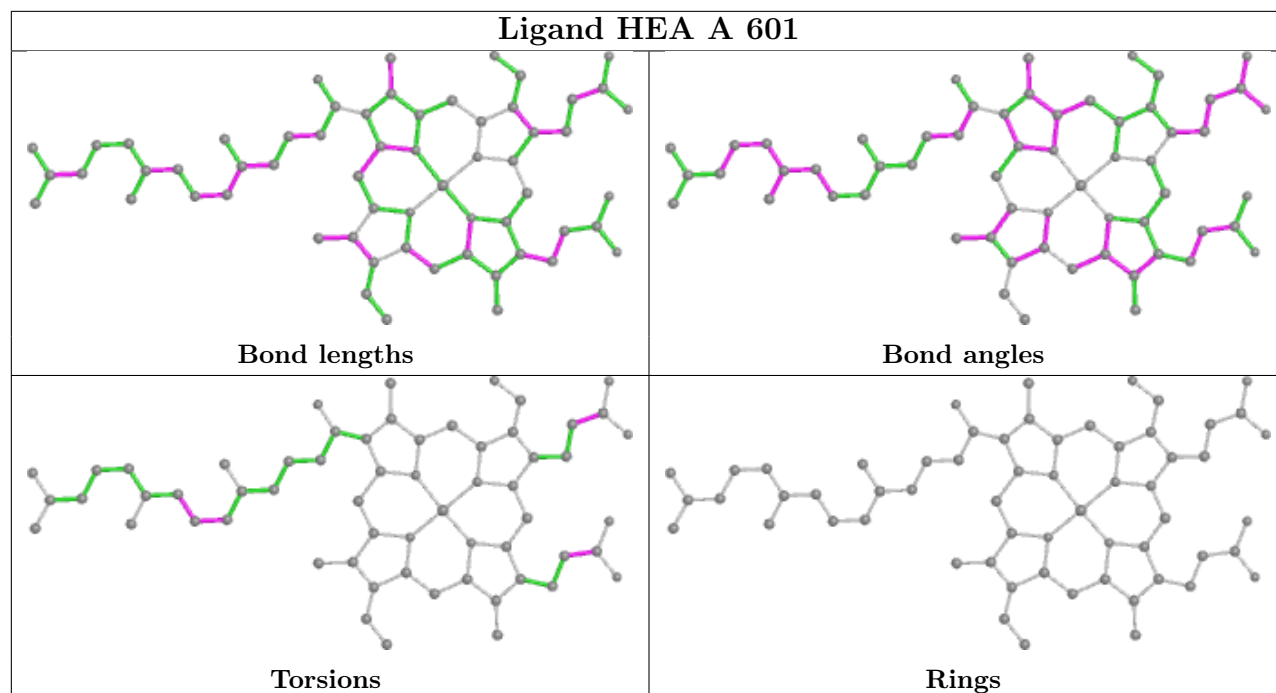


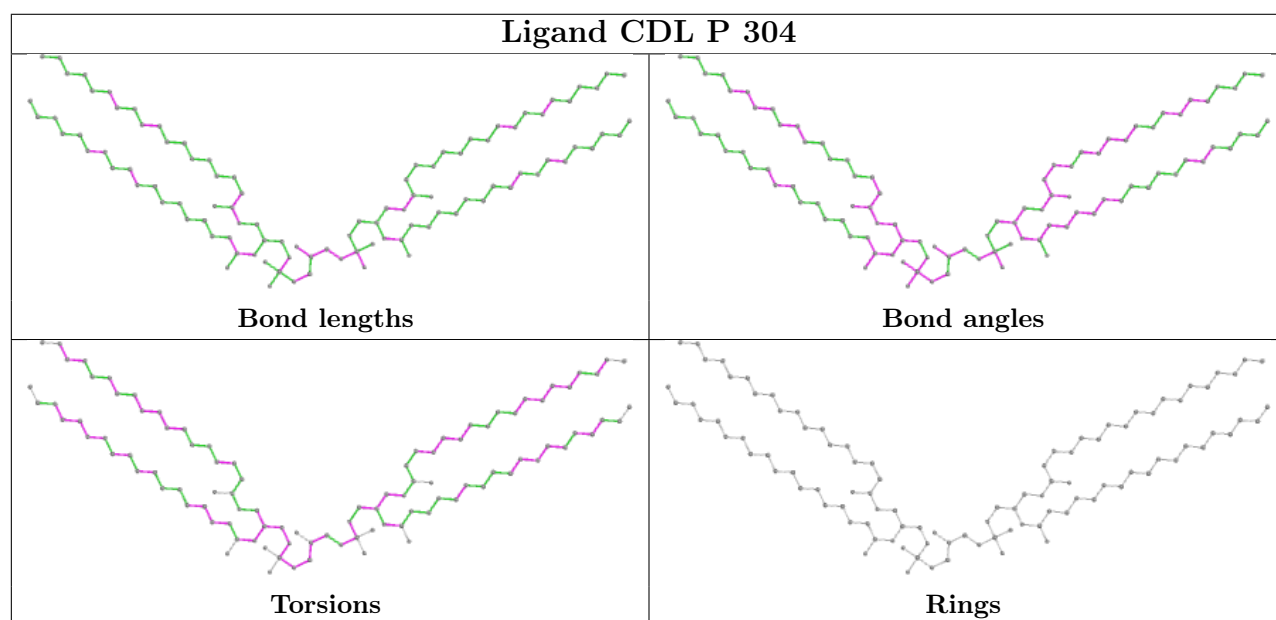
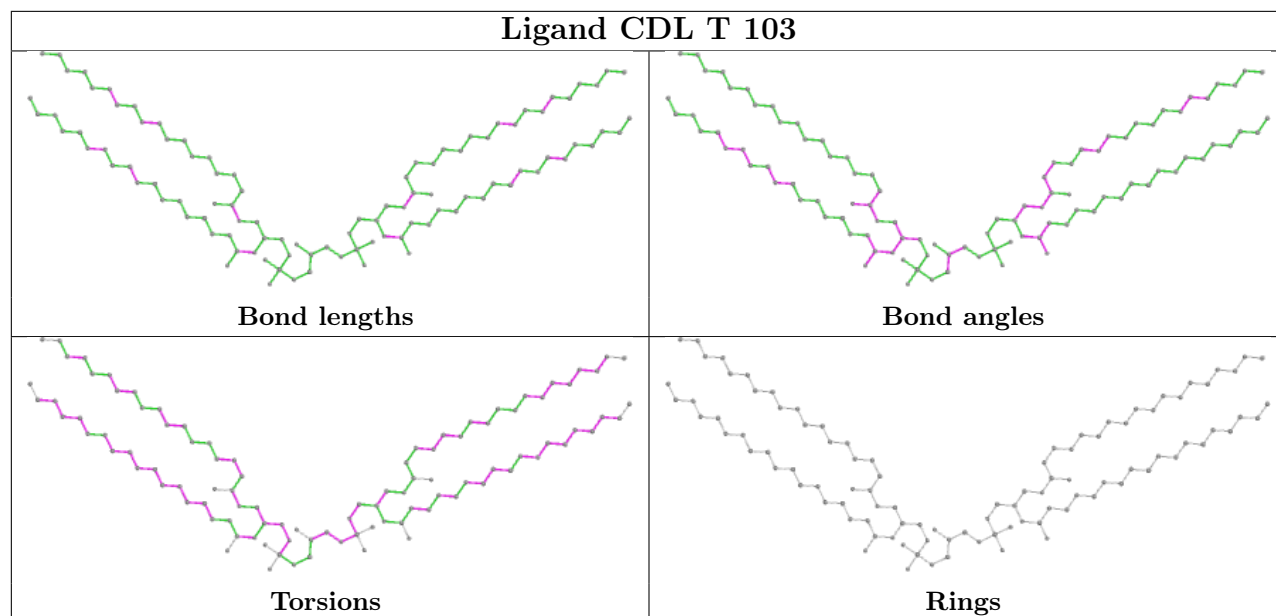


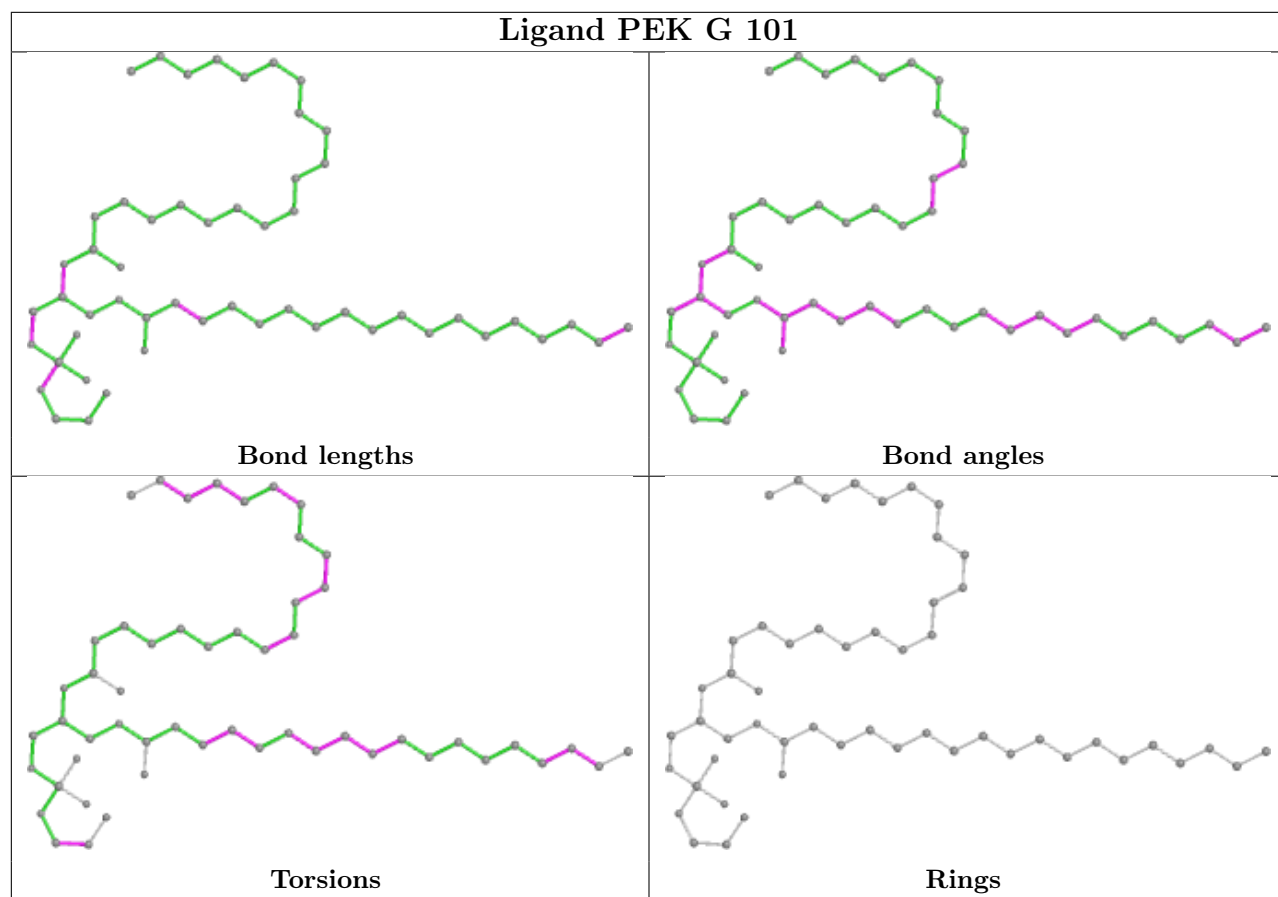
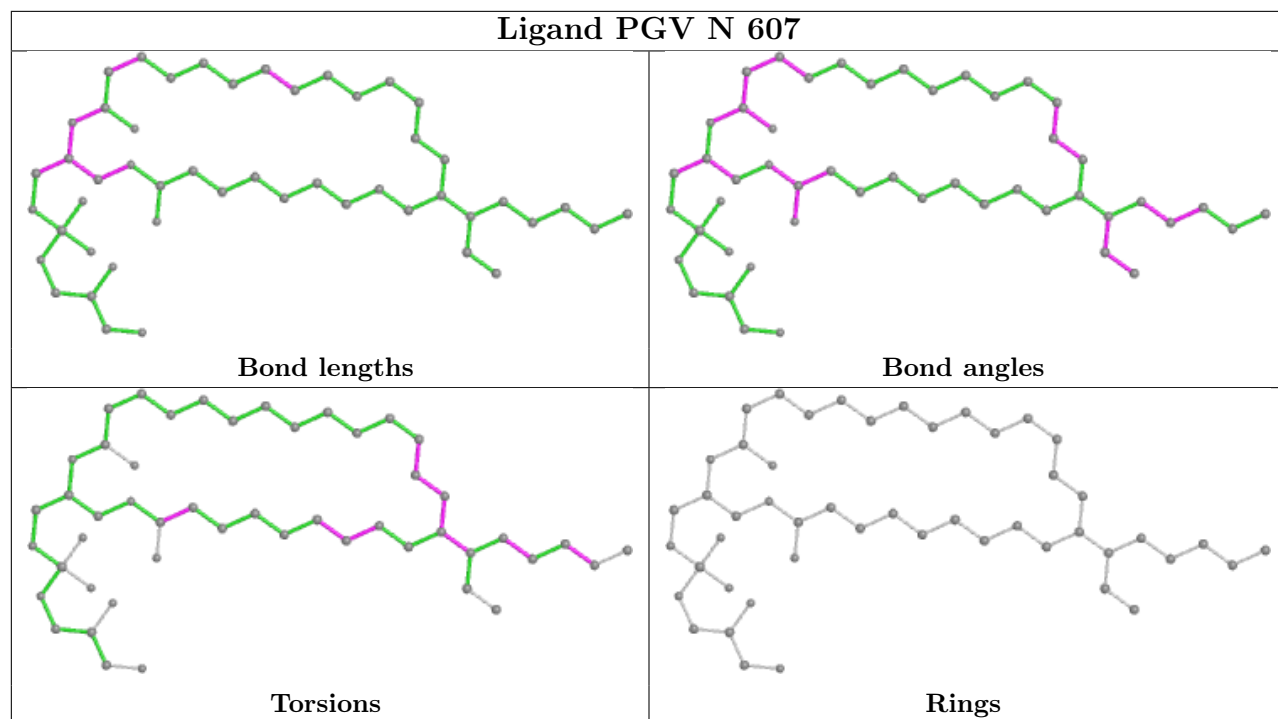


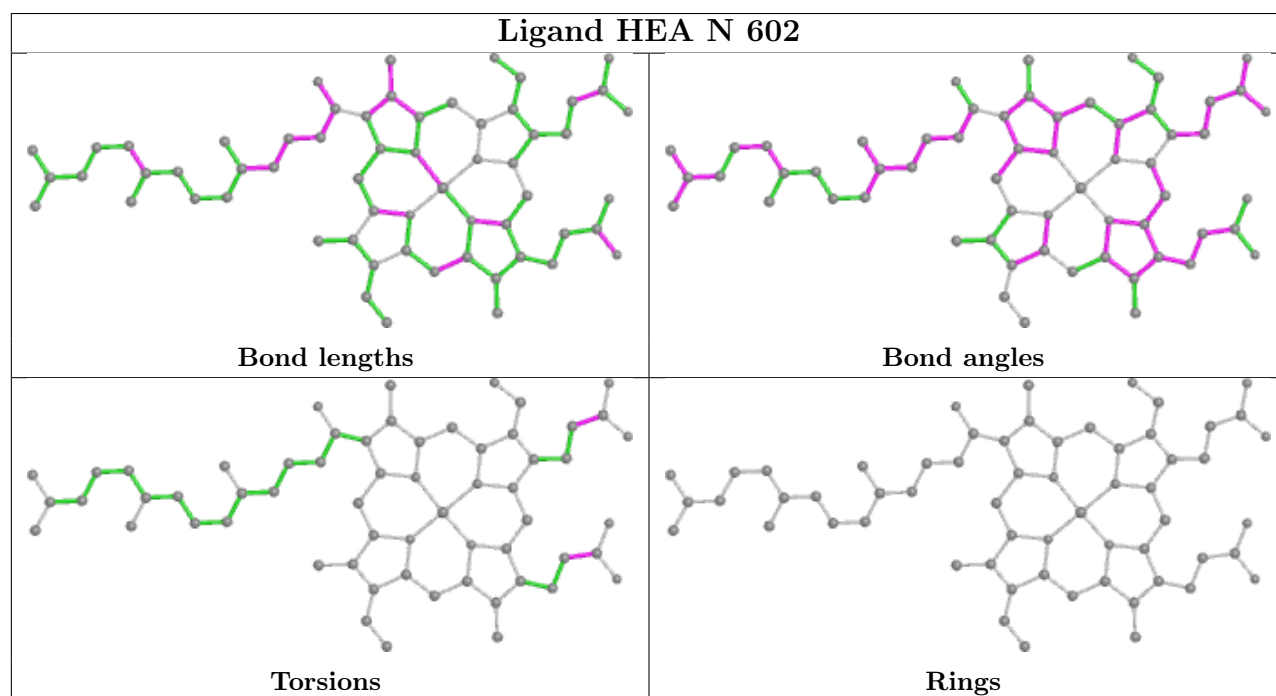
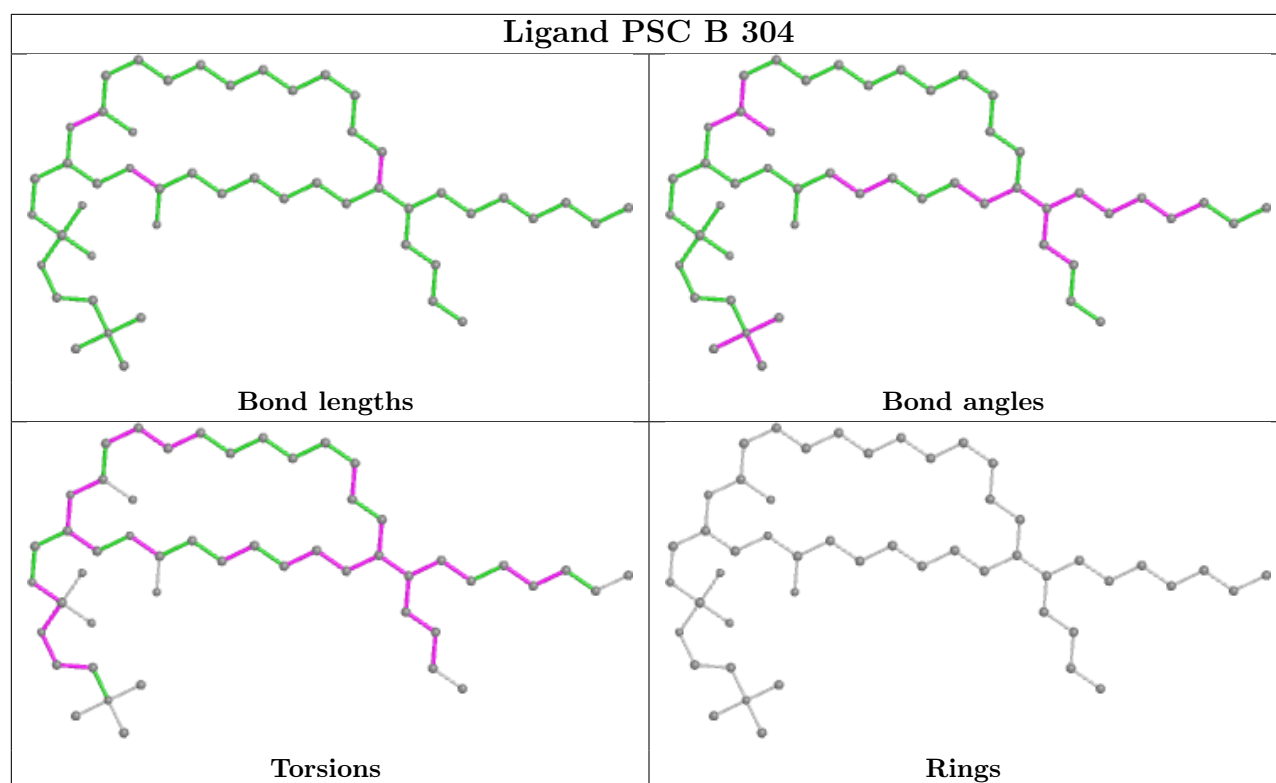


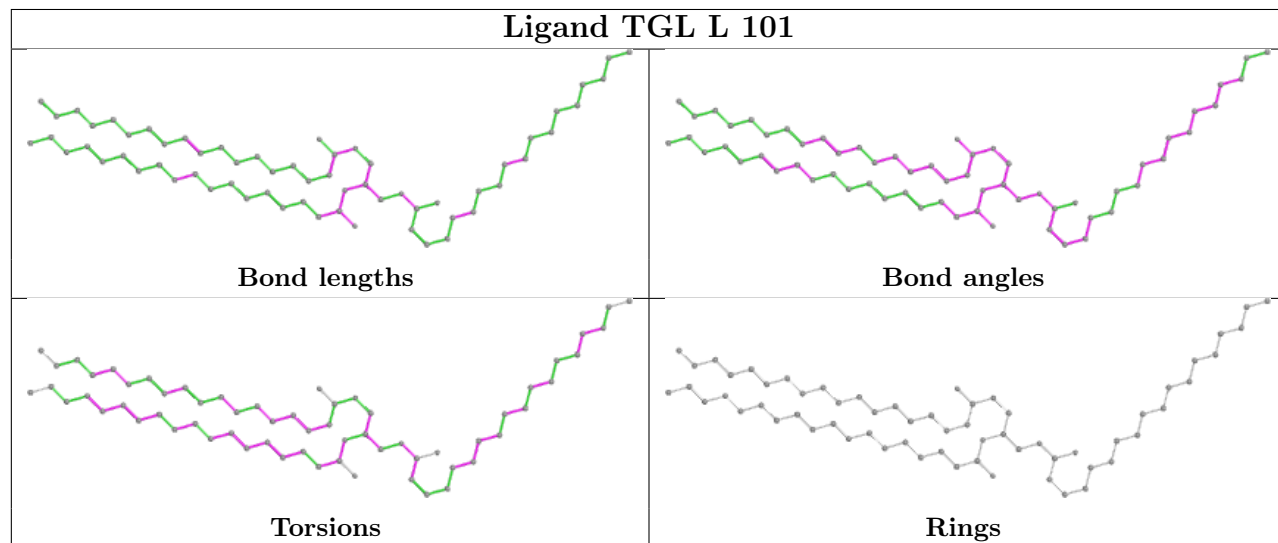
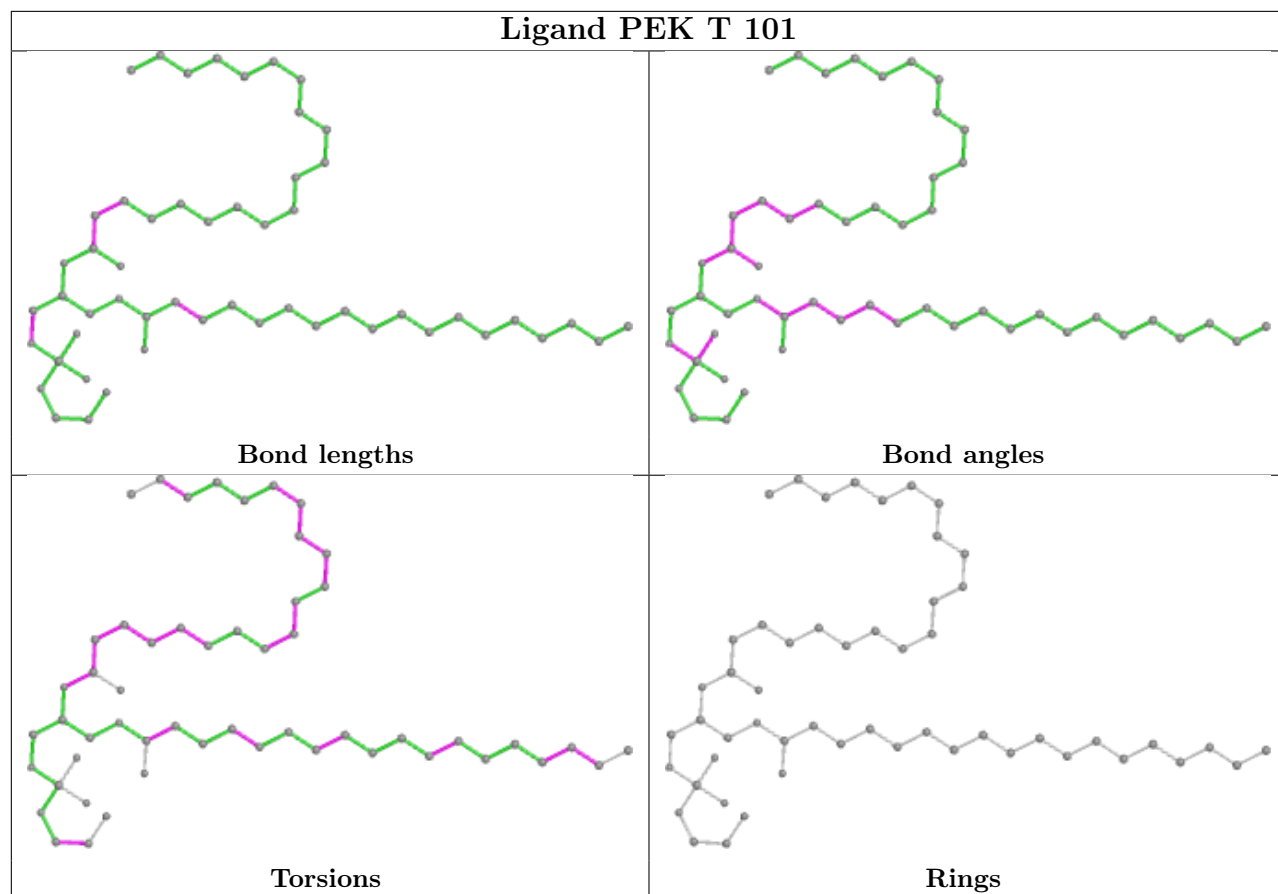




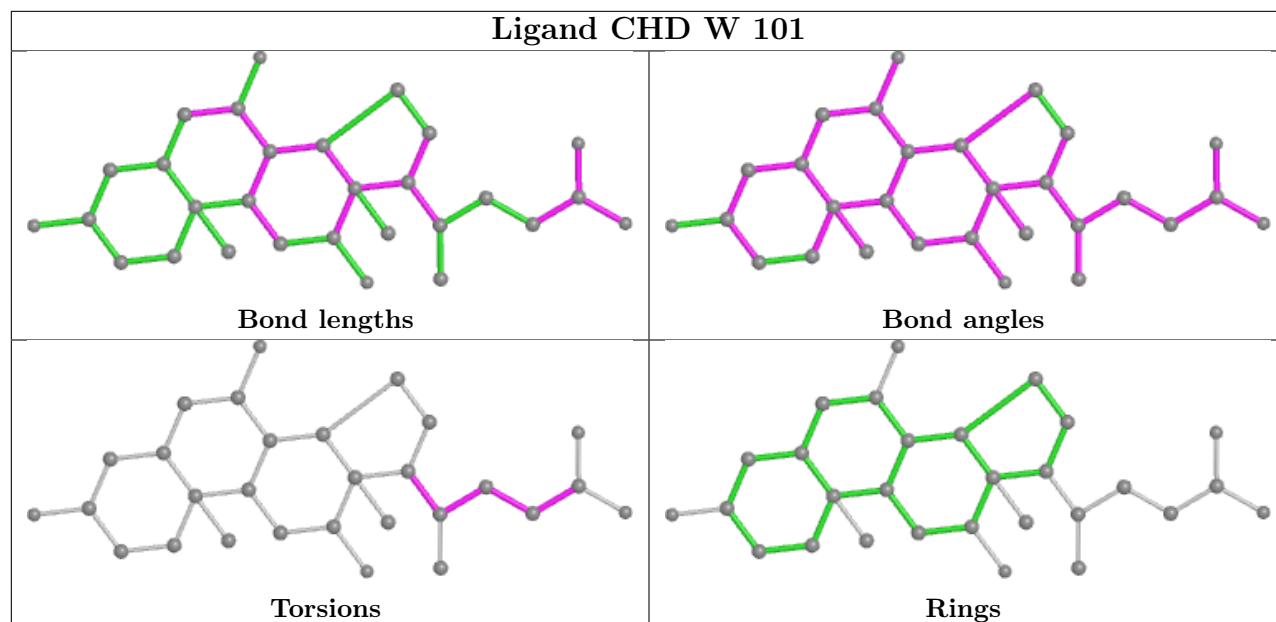




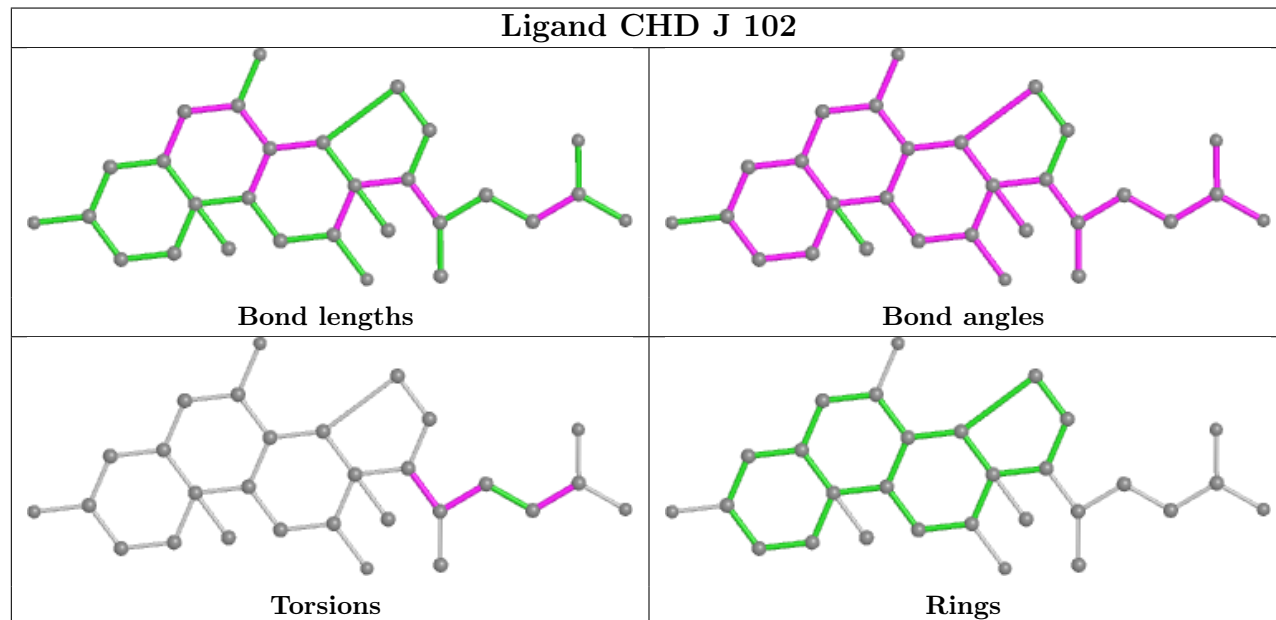




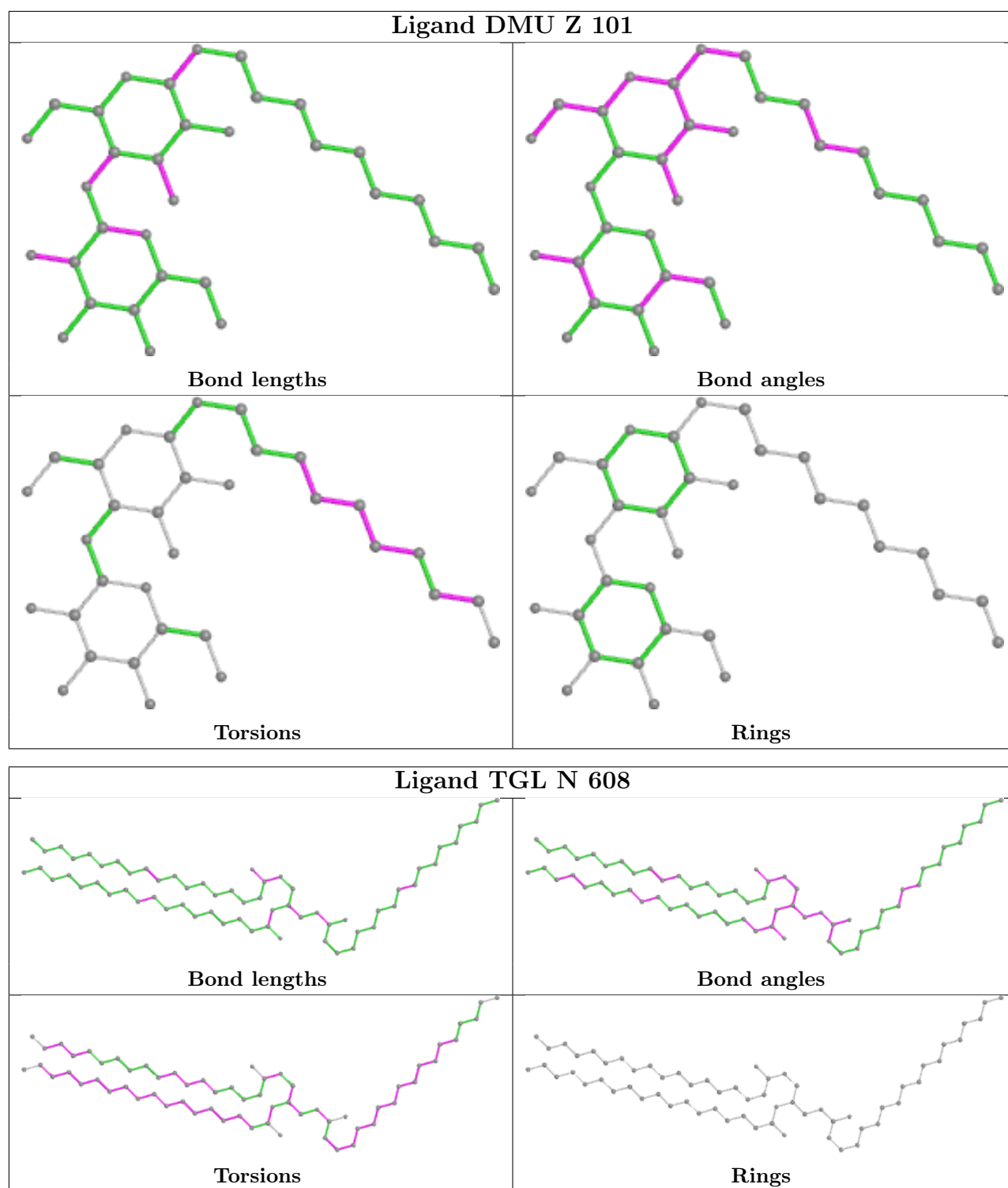
## Ligand CHD W 101



## Ligand CHD J 102







## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	513/514 (99%)	-0.46	0 100 100	17, 21, 29, 68	0
1	N	513/514 (99%)	-0.37	0 100 100	18, 25, 33, 63	0
2	B	226/227 (99%)	-0.50	4 (1%) 68 73	20, 28, 49, 71	0
2	O	226/227 (99%)	-0.50	4 (1%) 68 73	25, 34, 60, 90	0
3	C	259/261 (99%)	-0.78	1 (0%) 92 94	19, 25, 37, 77	0
3	P	259/261 (99%)	-0.69	3 (1%) 79 82	20, 26, 39, 70	0
4	D	144/147 (97%)	-0.79	1 (0%) 87 90	23, 29, 44, 76	0
4	Q	144/147 (97%)	0.57	15 (10%) 6 6	29, 43, 76, 153	0
5	E	105/109 (96%)	-0.62	2 (1%) 66 71	23, 30, 53, 136	0
5	R	105/109 (96%)	-0.12	3 (2%) 51 56	26, 38, 61, 149	0
6	F	98/98 (100%)	0.06	8 (8%) 11 12	21, 31, 94, 141	0
6	S	98/98 (100%)	0.07	9 (9%) 9 9	22, 30, 80, 126	0
7	G	83/85 (97%)	0.59	16 (19%) 1 1	23, 32, 110, 158	0
7	T	83/85 (97%)	0.56	18 (21%) 0 0	23, 36, 112, 154	0
8	H	79/85 (92%)	-0.18	7 (8%) 9 10	25, 36, 92, 133	0
8	U	79/85 (92%)	-0.21	9 (11%) 5 4	31, 40, 103, 127	0
9	I	72/73 (98%)	0.27	7 (9%) 7 8	27, 41, 63, 82	0
9	V	72/73 (98%)	0.55	7 (9%) 7 8	27, 47, 69, 143	0
10	J	58/59 (98%)	0.33	7 (12%) 4 4	25, 35, 65, 134	0
10	W	58/59 (98%)	0.04	6 (10%) 6 6	27, 38, 69, 158	0
11	K	49/56 (87%)	-0.54	0 100 100	28, 35, 49, 58	0
11	X	49/56 (87%)	1.03	7 (14%) 2 2	36, 47, 68, 81	0
12	L	46/47 (97%)	-0.73	1 (2%) 62 67	22, 27, 53, 95	0
12	Y	46/47 (97%)	-0.60	1 (2%) 62 67	27, 34, 58, 125	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
13	M	43/46 (93%)	-0.24	4 (9%) 8 9	24, 28, 64, 118	0
13	Z	43/46 (93%)	0.04	5 (11%) 4 4	31, 38, 79, 145	0
All	All	3550/3614 (98%)	-0.29	145 (4%) 37 41	17, 29, 61, 158	0

The worst 5 of 145 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
6	S	97	ALA	37.3
4	Q	5	VAL	33.6
4	Q	6	VAL	14.6
6	F	97	ALA	13.2
6	F	98	HIS	11.6

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
7	TPO	G	11	11/12	0.49	0.31	61,102,165,180	0
9	SAC	V	1	9/10	0.52	0.37	114,126,154,162	0
7	TPO	T	11	11/12	0.62	0.24	75,96,179,187	0
9	SAC	I	1	9/10	0.91	0.31	54,70,75,82	0
2	FME	O	1	10/11	0.97	0.06	29,34,48,59	0
1	FME	A	1	10/11	0.97	0.10	31,39,64,89	0
2	FME	B	1	10/11	0.97	0.08	22,27,48,55	0
1	FME	N	1	10/11	0.98	0.09	33,40,72,74	0

## 6.3 Carbohydrates [i](#)

There are no monosaccharides 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, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column

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

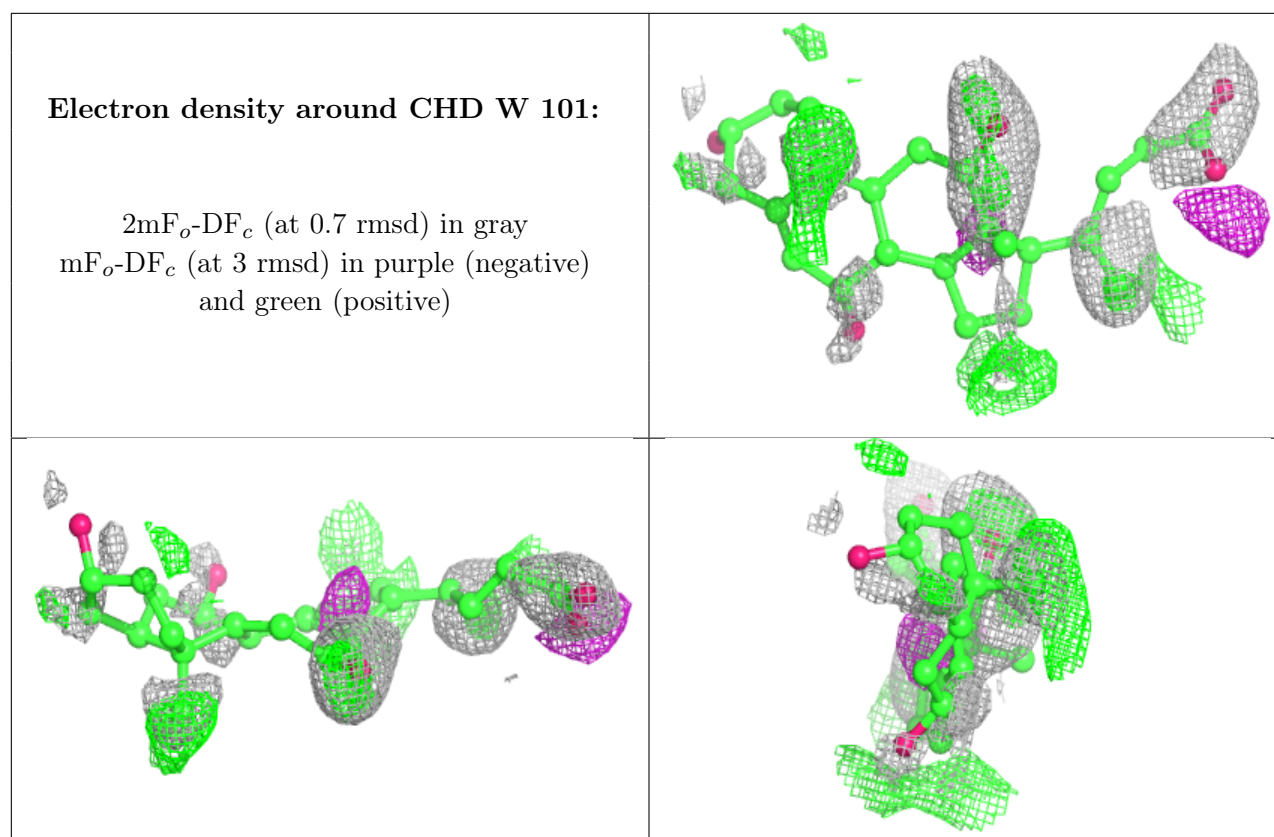
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
22	CHD	W	101	29/29	0.45	0.43	59,136,161,164	0
22	CHD	J	102	29/29	0.57	0.44	57,133,165,172	0
25	CDL	T	103	100/100	0.69	0.28	50,91,168,202	0
25	CDL	G	102	100/100	0.70	0.30	51,96,162,187	0
26	PEK	T	102	53/53	0.73	0.28	44,95,159,176	0
26	PEK	P	308	53/53	0.77	0.21	38,79,120,144	0
26	PEK	C	306	53/53	0.77	0.17	34,78,128,158	0
28	DMU	J	101	33/33	0.77	0.25	44,65,156,176	0
22	CHD	P	305	29/29	0.78	0.35	47,91,109,111	0
22	CHD	C	304	29/29	0.79	0.40	46,96,116,118	0
20	TGL	Q	202	63/63	0.79	0.17	40,62,99,110	0
28	DMU	P	306	33/33	0.80	0.21	36,105,192,198	0
24	UNX	P	302	1/1	0.81	0.39	38,38,38,38	0
19	PGV	P	301	51/51	0.81	0.21	45,85,147,177	0
19	PGV	C	307	51/51	0.81	0.25	47,85,169,181	0
23	PSC	B	304	52/52	0.81	0.24	39,107,190,207	0
20	TGL	Y	101	63/63	0.83	0.18	37,64,106,120	0
26	PEK	G	103	53/53	0.83	0.25	46,93,150,158	0
23	PSC	O	303	52/52	0.84	0.23	37,75,191,211	0
25	CDL	C	303	100/100	0.85	0.32	29,72,127,137	0
20	TGL	D	201	63/63	0.85	0.14	30,55,90,101	0
19	PGV	Q	201	51/51	0.85	0.26	42,79,161,191	0
28	DMU	Z	101	33/33	0.85	0.20	41,47,61,63	0
20	TGL	N	608	63/63	0.86	0.20	47,76,120,136	0
25	CDL	P	304	100/100	0.86	0.27	32,78,140,149	0
20	TGL	L	101	63/63	0.87	0.17	30,58,96,101	0
19	PGV	A	608	51/51	0.87	0.17	31,68,107,120	0
24	UNX	C	301	1/1	0.89	0.23	39,39,39,39	0
20	TGL	B	301	63/63	0.90	0.12	40,65,90,103	0
28	DMU	M	101	33/33	0.90	0.14	33,38,51,53	0
22	CHD	C	305	29/29	0.93	0.07	26,28,31,34	0
22	CHD	P	307	29/29	0.94	0.06	25,29,32,35	0
22	CHD	O	302	29/29	0.95	0.07	23,26,29,34	0
22	CHD	B	303	29/29	0.95	0.06	24,26,29,37	0
18	PER	N	606	2/2	0.96	0.18	20,20,20,26	0
26	PEK	T	101	53/53	0.97	0.08	26,42,83,92	0
18	PER	A	606	2/2	0.97	0.18	17,17,17,22	0
26	PEK	G	101	53/53	0.97	0.08	23,40,82,96	0
19	PGV	N	607	51/51	0.98	0.07	22,32,67,68	0
19	PGV	C	302	51/51	0.98	0.07	22,29,81,98	0
19	PGV	P	303	51/51	0.98	0.06	21,30,71,79	0
19	PGV	A	607	51/51	0.98	0.07	20,29,56,65	0

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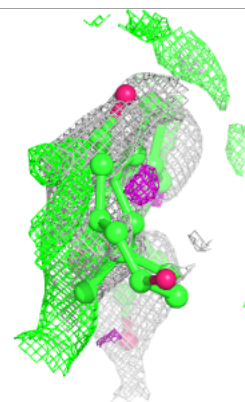
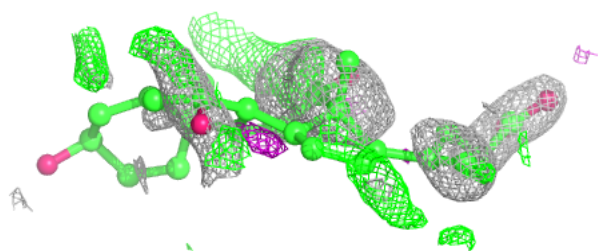
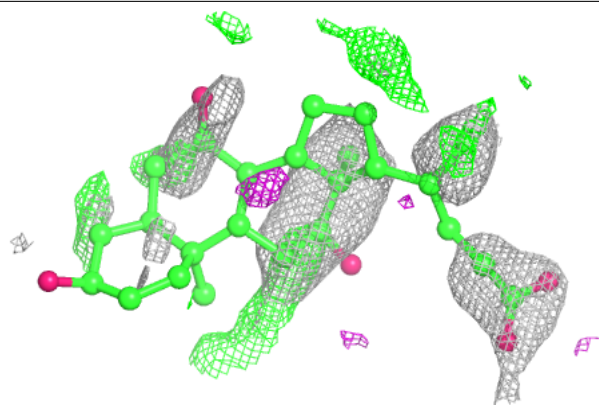
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
14	HEA	A	601	60/60	0.99	0.07	17,19,37,48	0
14	HEA	A	602	60/60	0.99	0.07	17,18,25,30	0
14	HEA	N	601	60/60	0.99	0.08	21,23,43,51	0
14	HEA	N	602	60/60	0.99	0.09	19,21,28,31	0
16	MG	A	604	1/1	0.99	0.04	19,19,19,19	0
16	MG	N	604	1/1	0.99	0.08	25,25,25,25	0
17	NA	A	605	1/1	1.00	0.02	23,23,23,23	0
17	NA	N	605	1/1	1.00	0.03	30,30,30,30	0
27	ZN	F	101	1/1	1.00	0.04	26,26,26,26	0
27	ZN	S	101	1/1	1.00	0.04	26,26,26,26	0
21	CUA	B	302	2/2	1.00	0.08	21,21,21,21	0
21	CUA	O	301	2/2	1.00	0.07	27,27,27,27	0
15	CU	A	603	1/1	1.00	0.07	19,19,19,19	0
15	CU	N	603	1/1	1.00	0.08	21,21,21,21	0

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

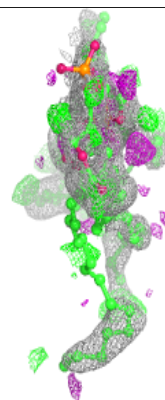
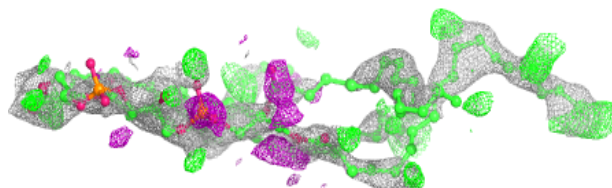
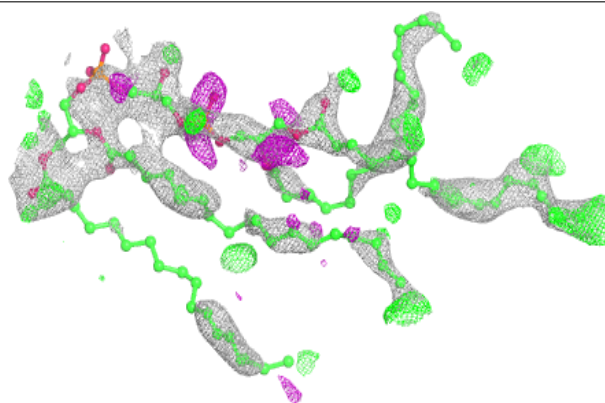


**Electron density around CHD J 102:**

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

**Electron density around CDL T 103:**

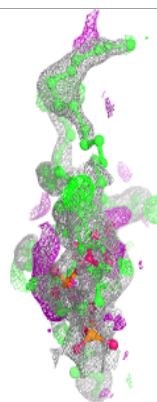
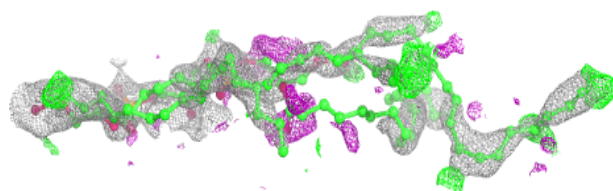
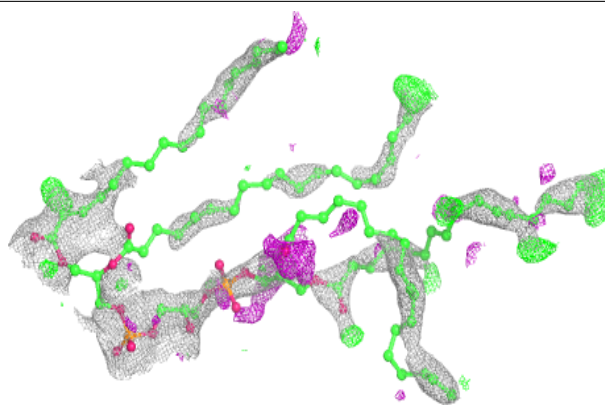
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



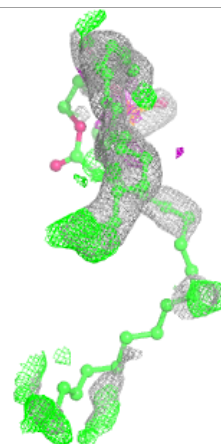
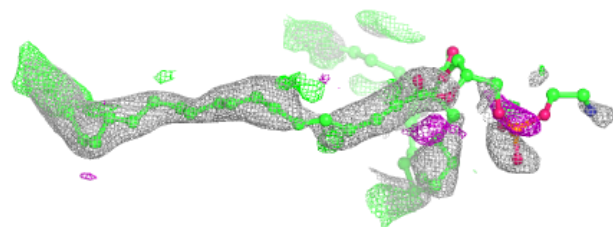
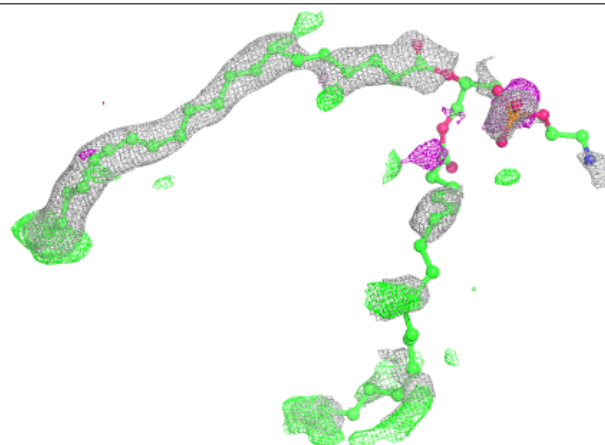


**Electron density around CDL G 102:**

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

**Electron density around PEK T 102:**

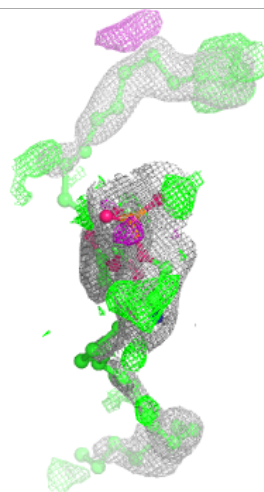
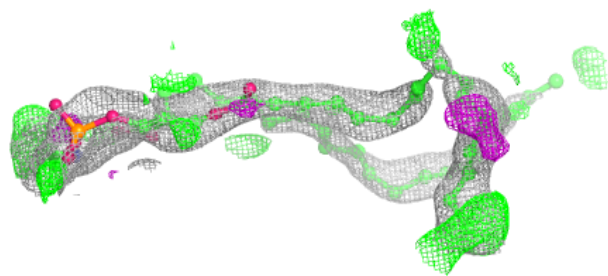
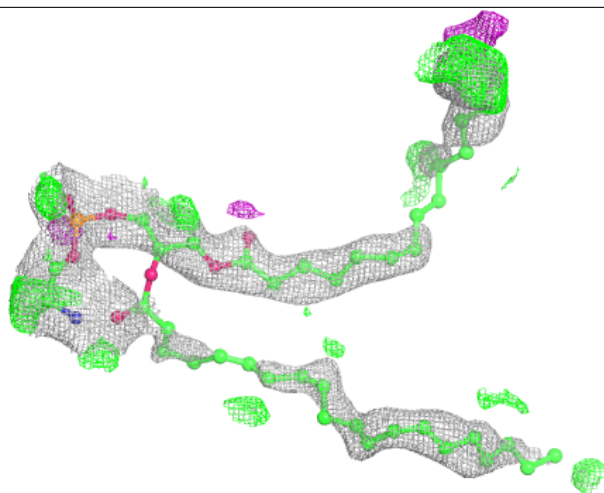
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





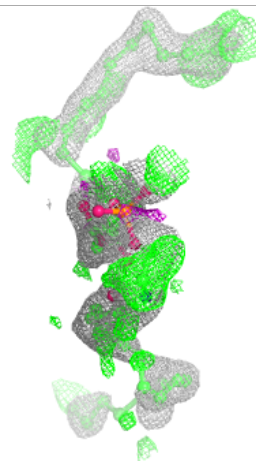
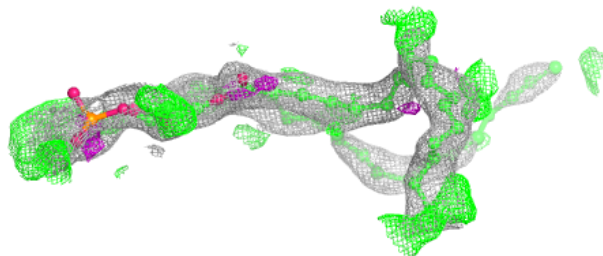
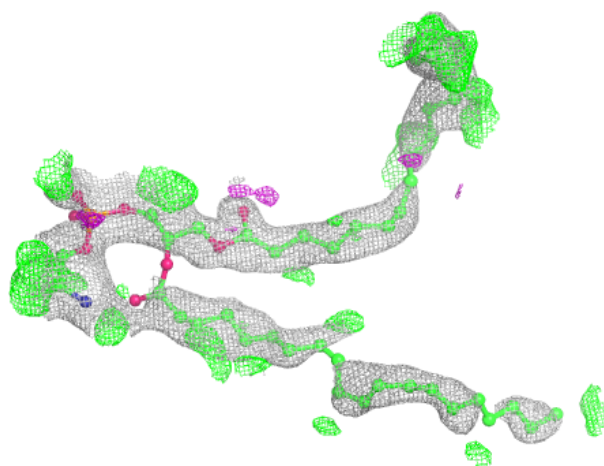
**Electron density around PEK P 308:**

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



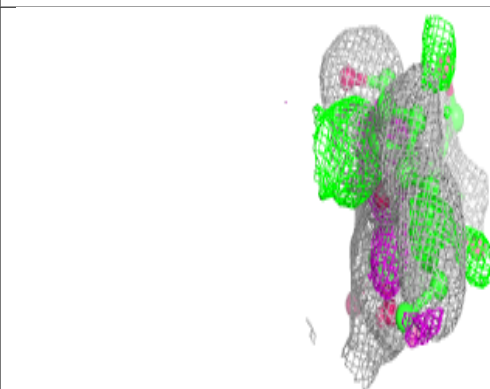
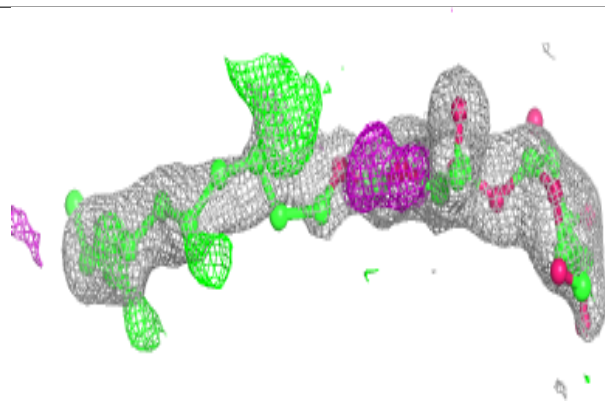
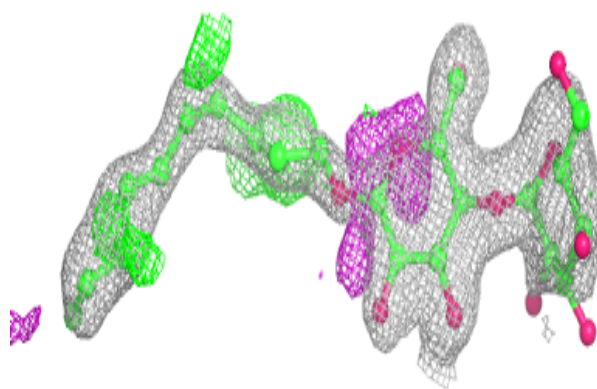
**Electron density around PEK C 306:**

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

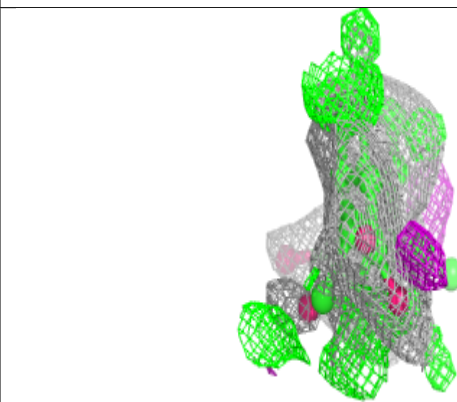
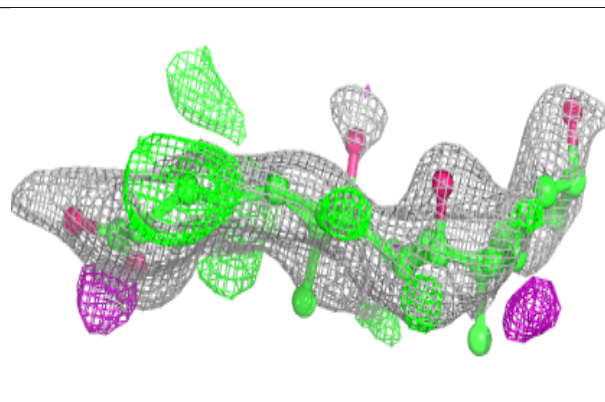
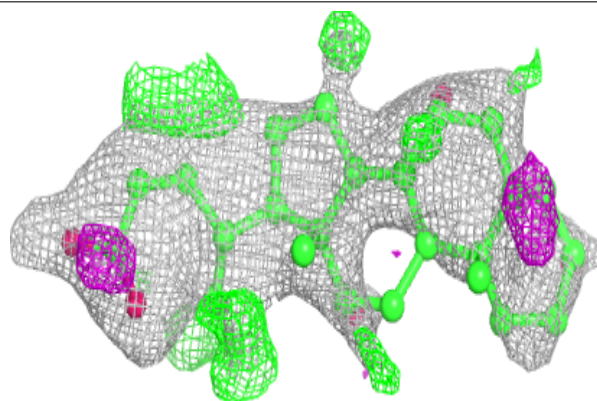


**Electron density around DMU J 101:**

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

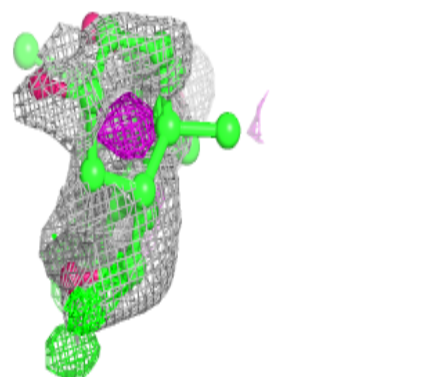
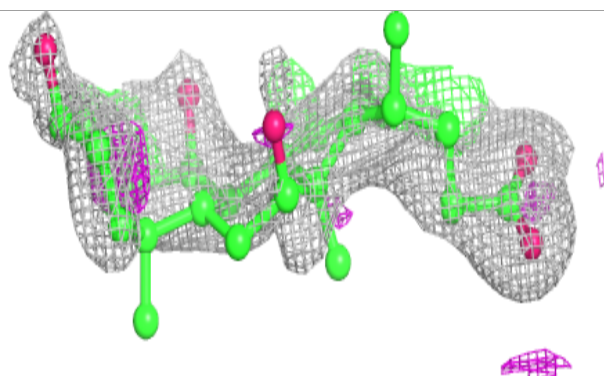
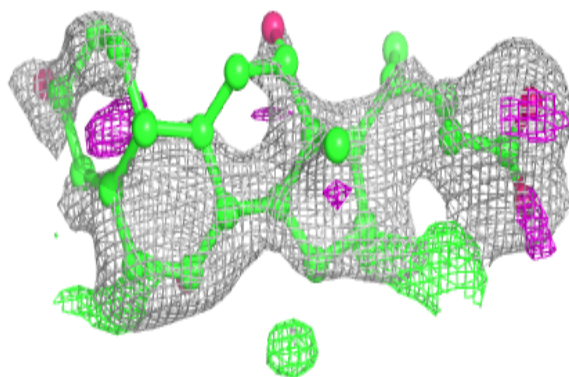
**Electron density around CHD P 305:**

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

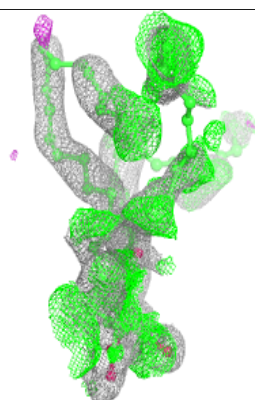
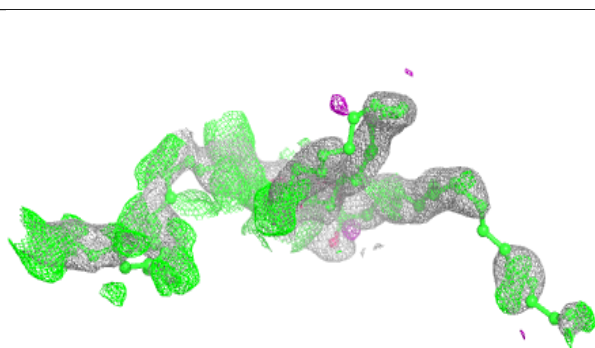
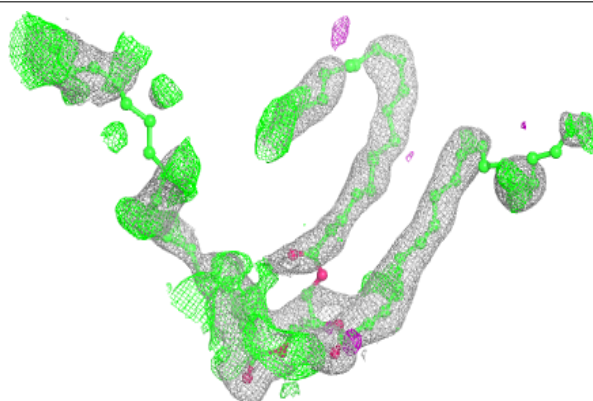


**Electron density around CHD C 304:**

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

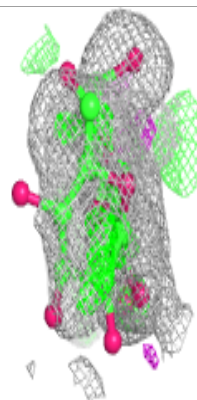
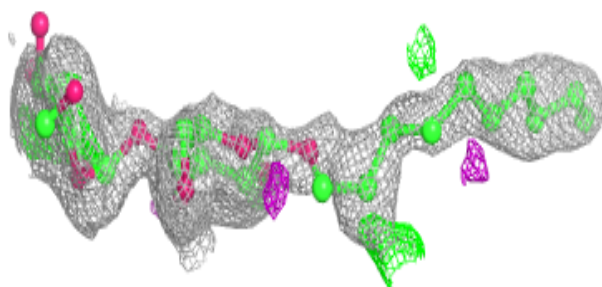
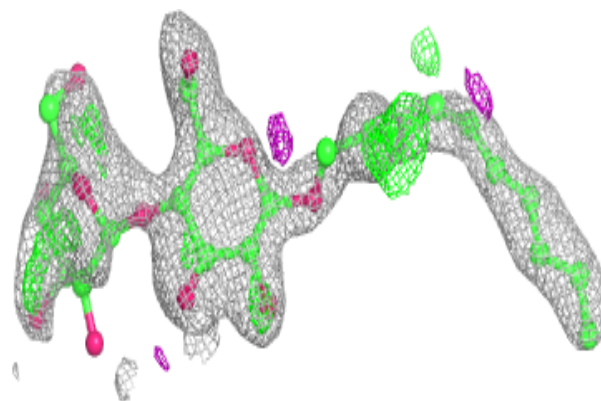
**Electron density around TGL Q 202:**

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

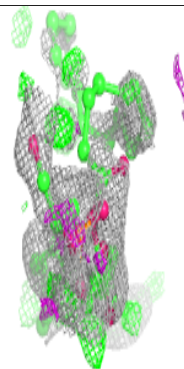
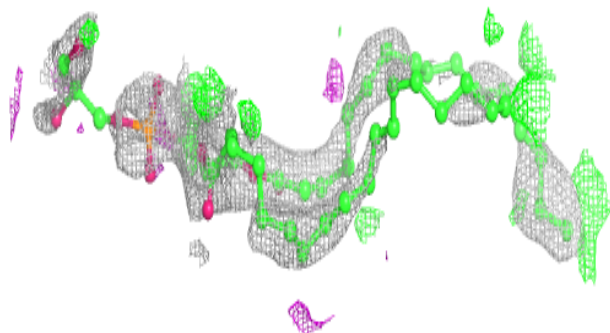
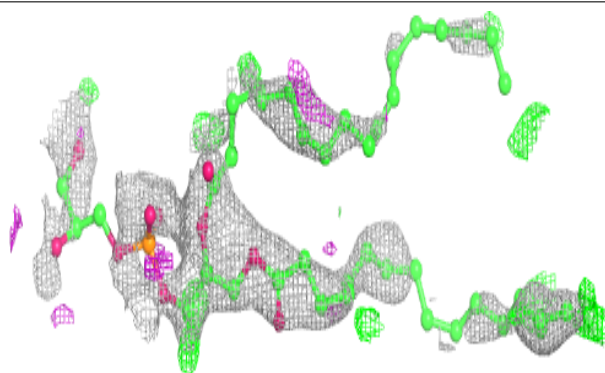


**Electron density around DMU P 306:**

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

**Electron density around PGV P 301:**

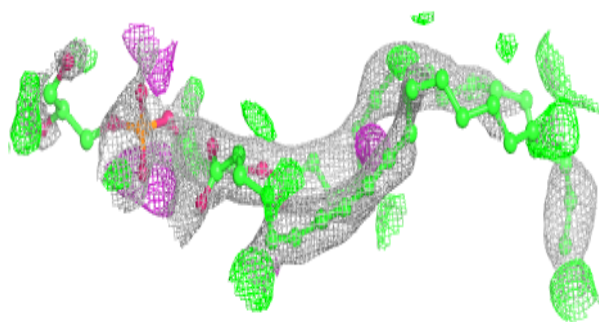
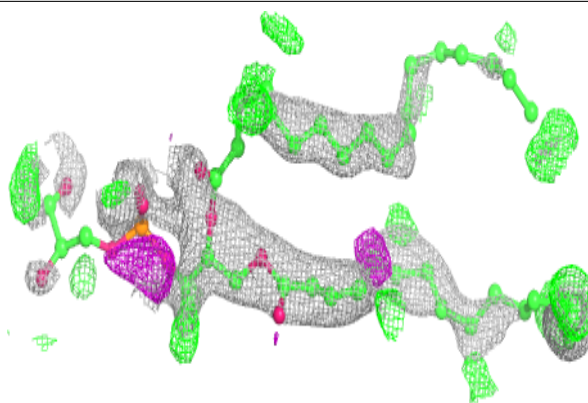
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



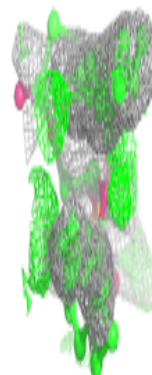
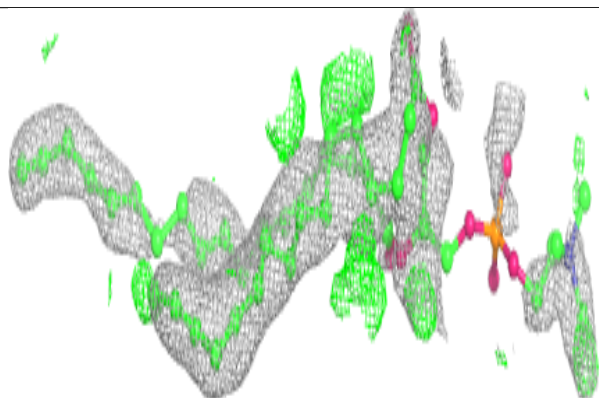
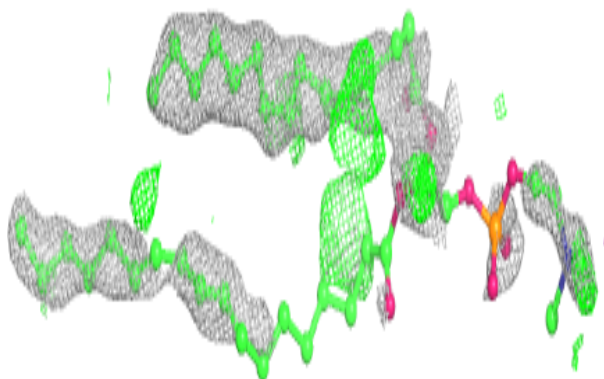


**Electron density around PGV C 307:**

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

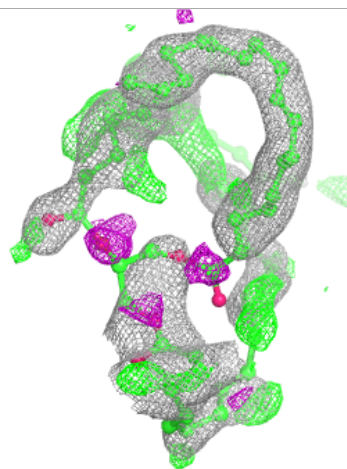
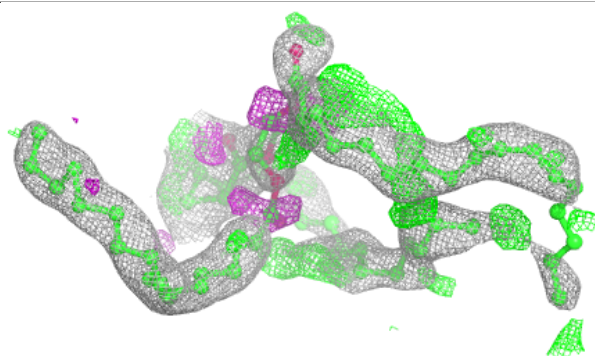
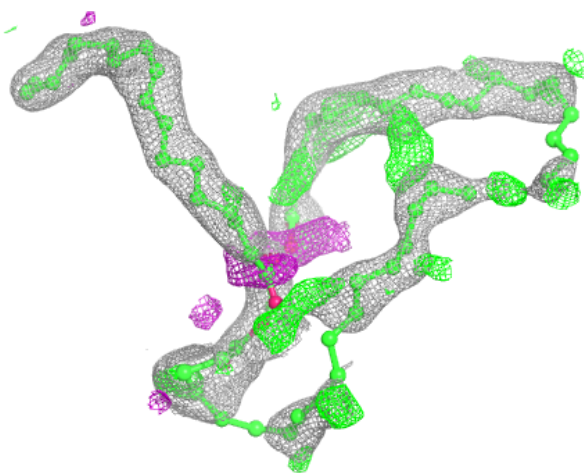
**Electron density around PSC B 304:**

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



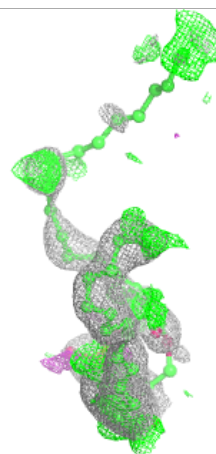
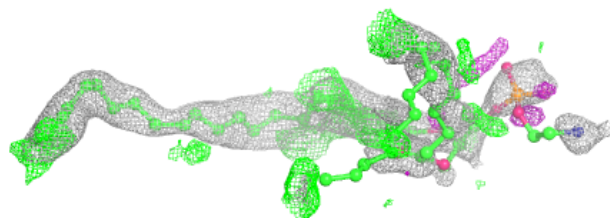
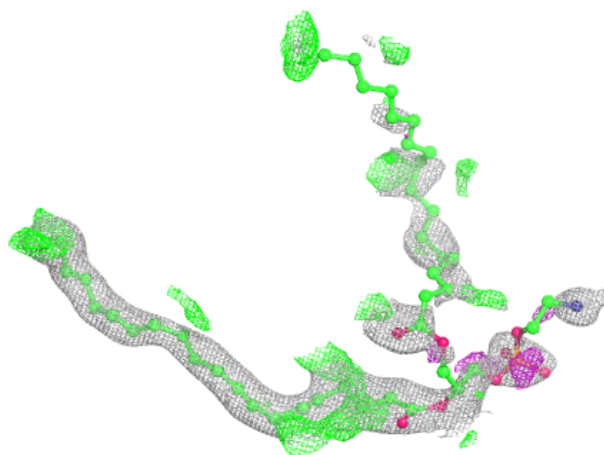
**Electron density around TGL Y 101:**

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



**Electron density around PEK G 103:**

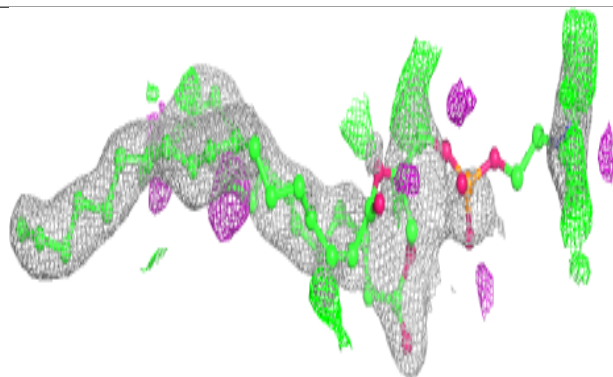
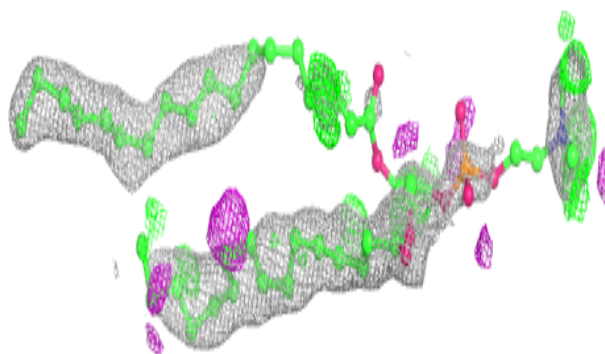
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



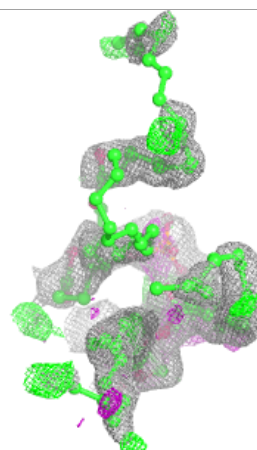
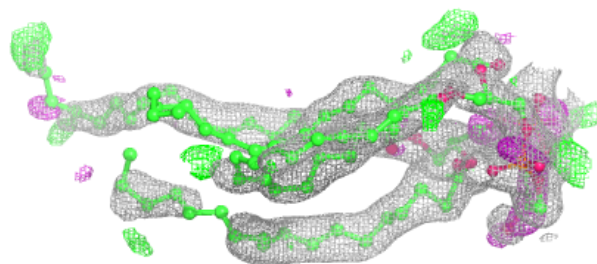
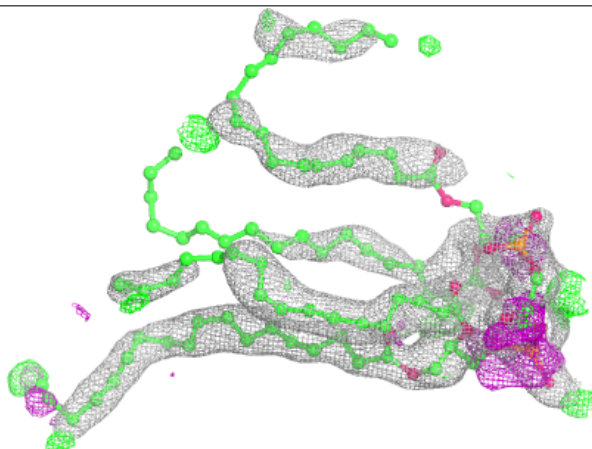


**Electron density around PSC O 303:**

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

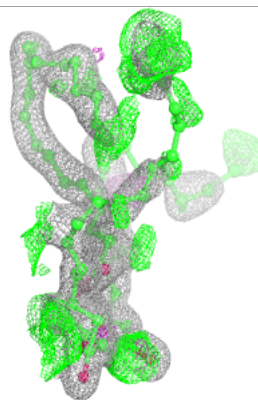
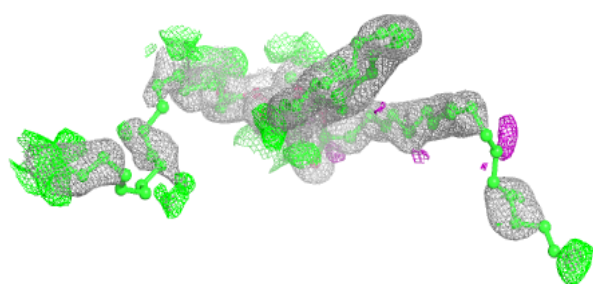
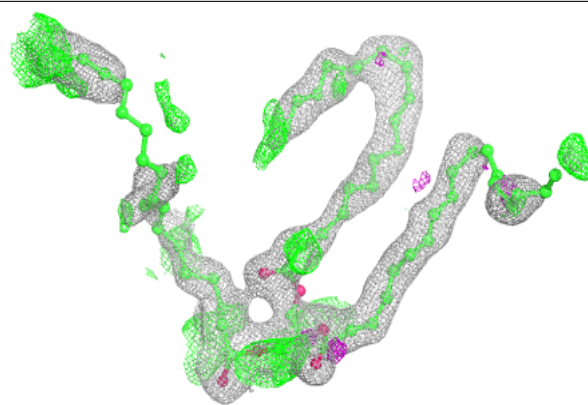
**Electron density around CDL C 303:**

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

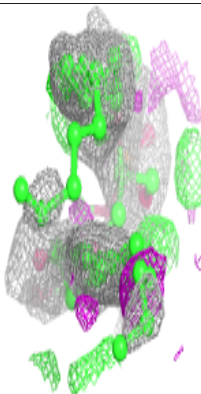
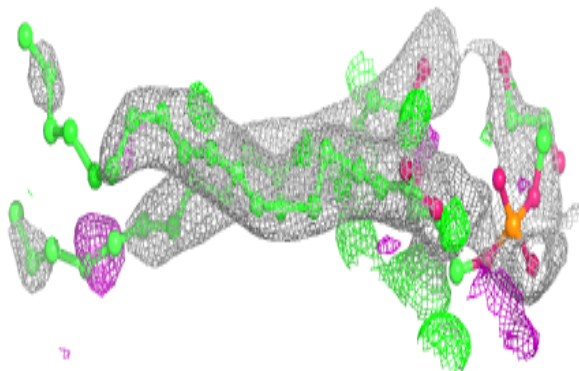
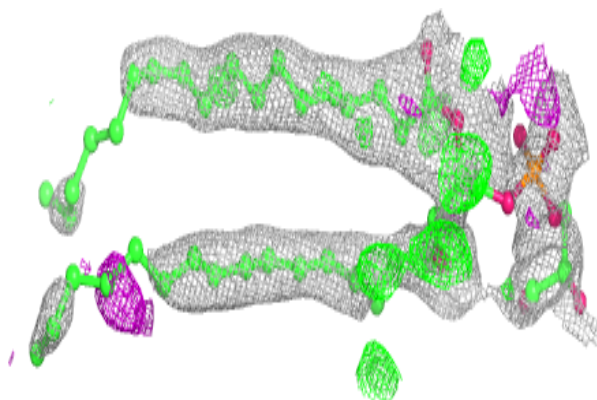


**Electron density around TGL D 201:**

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

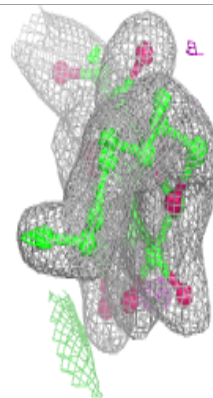
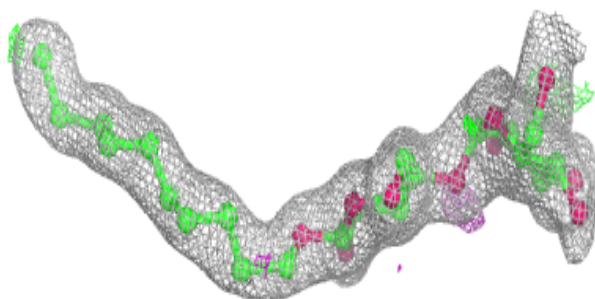
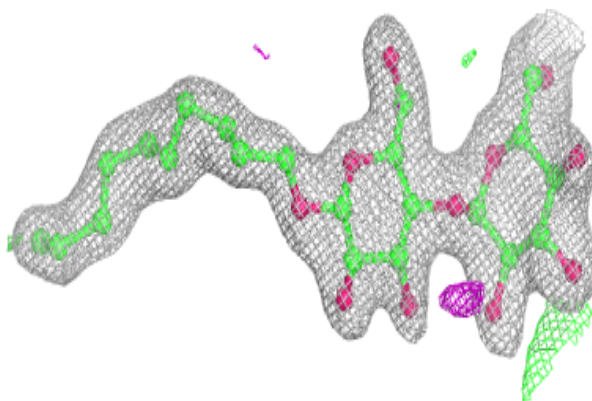
**Electron density around PGV Q 201:**

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

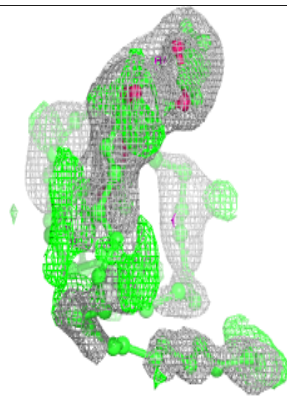
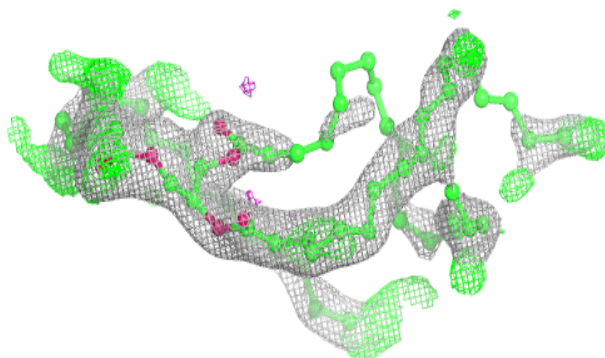
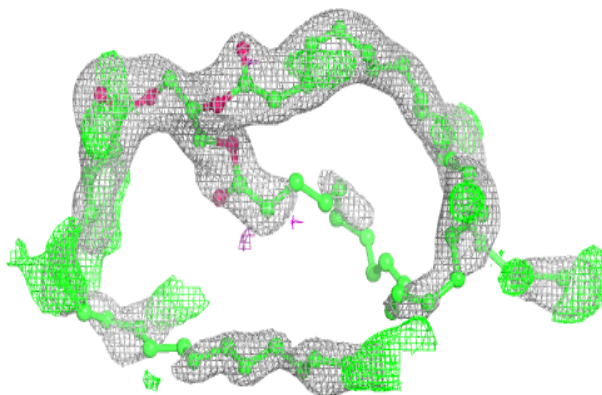


**Electron density around DMU Z 101:**

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

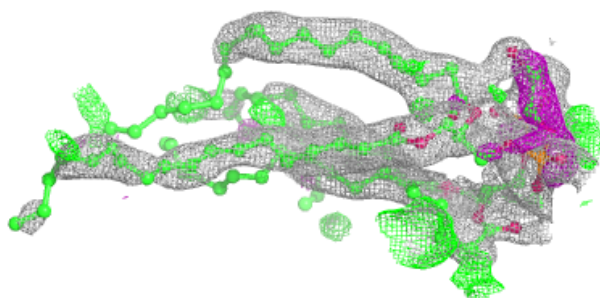
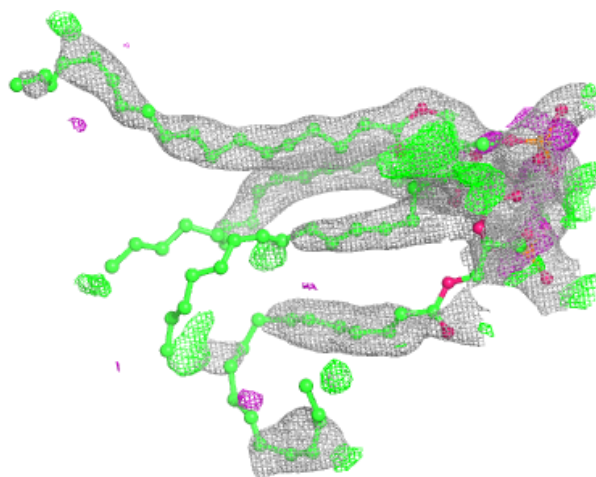
**Electron density around TGL N 608:**

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



**Electron density around CDL P 304:**

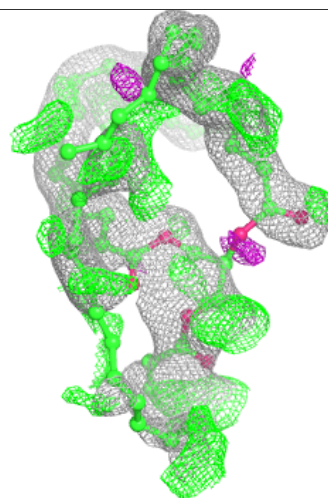
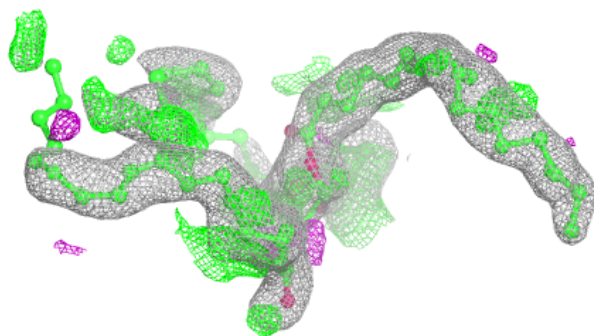
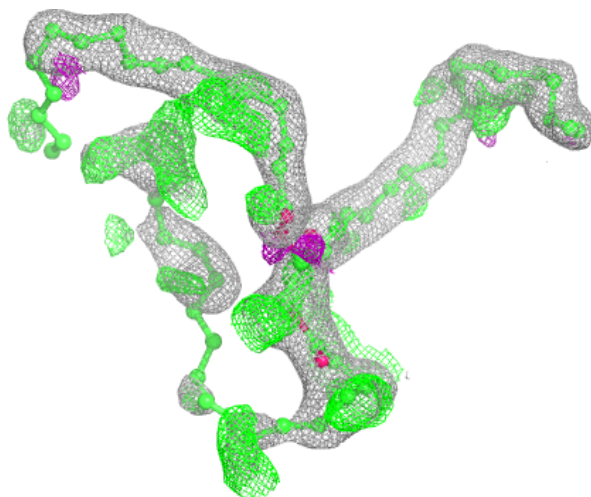
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





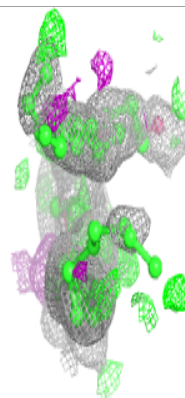
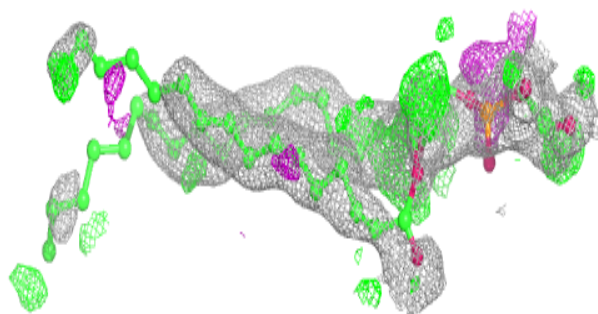
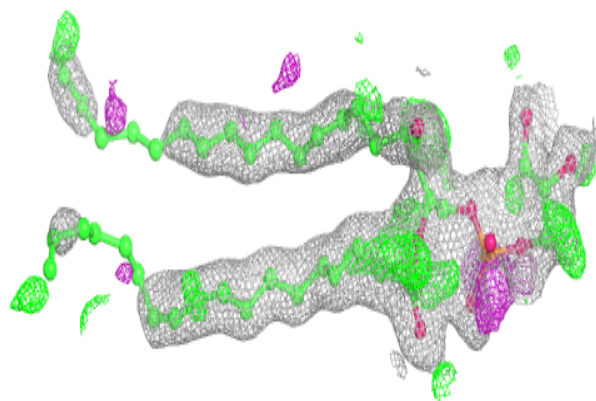
**Electron density around TGL L 101:**

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

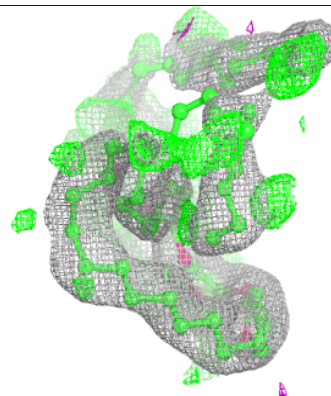
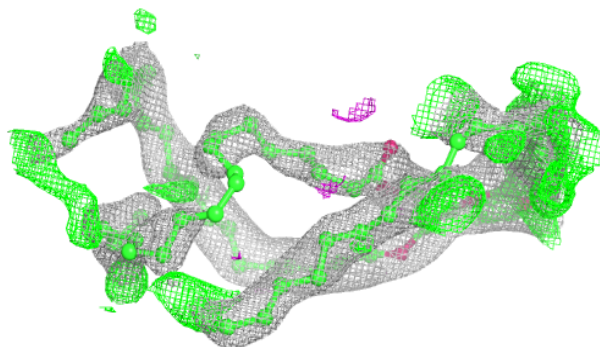
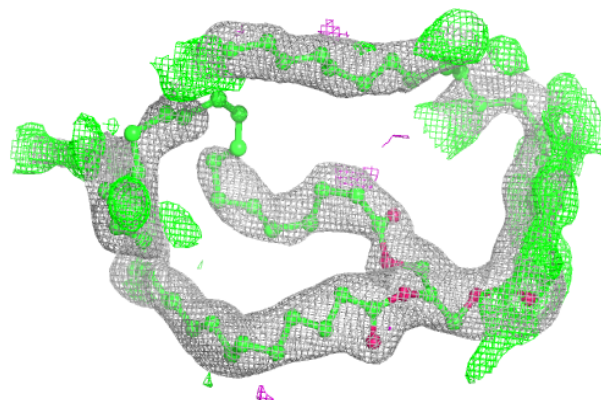


**Electron density around PGV A 608:**

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

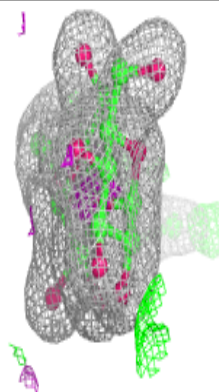
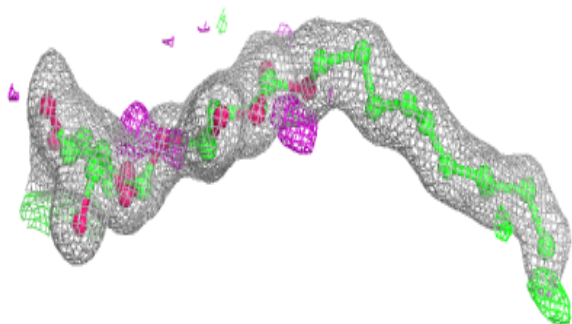
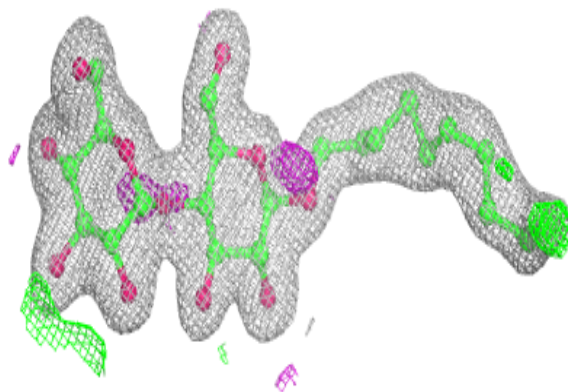
**Electron density around TGL B 301:**

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

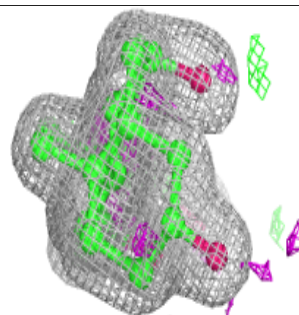
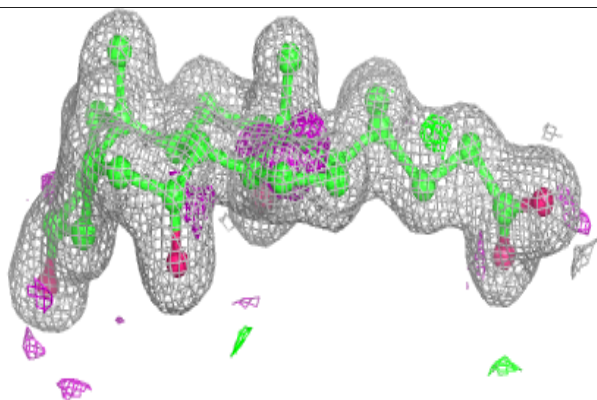
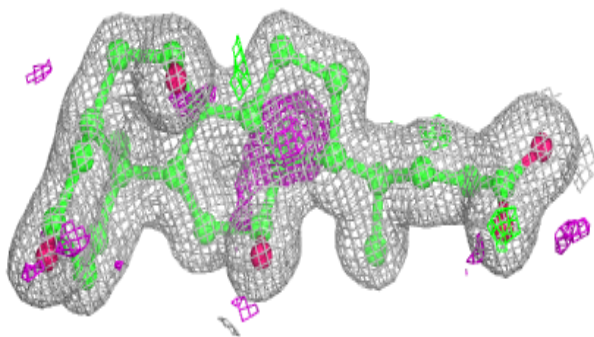


**Electron density around DMU M 101:**

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

**Electron density around CHD C 305:**

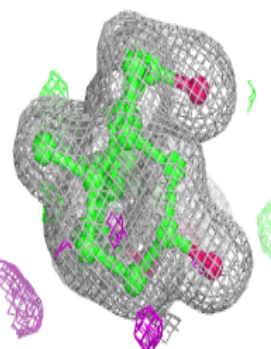
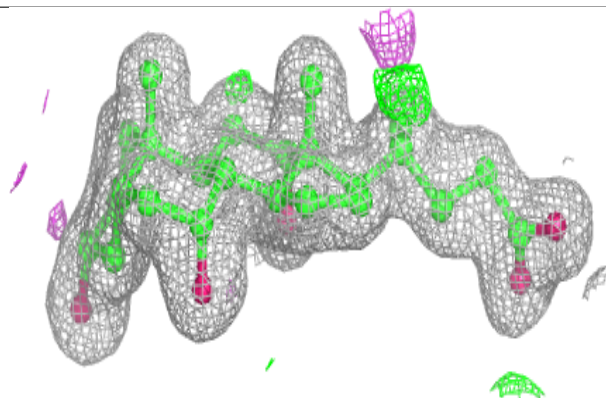
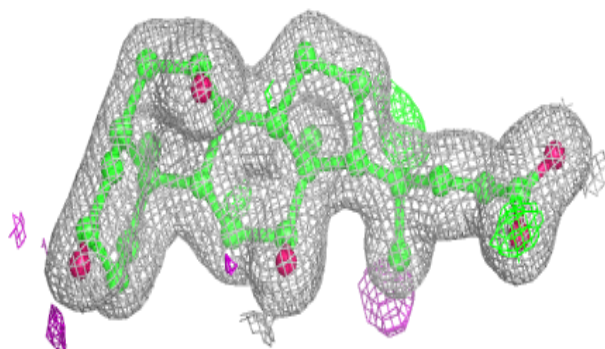
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



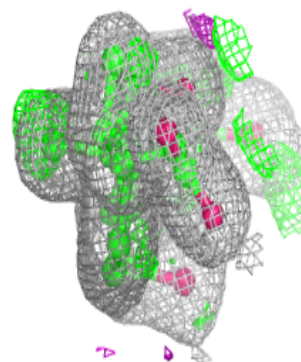
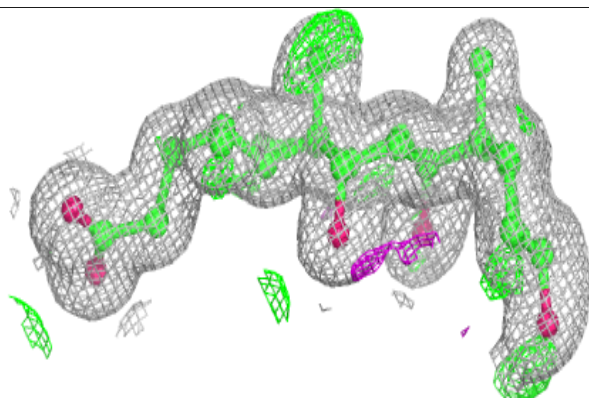
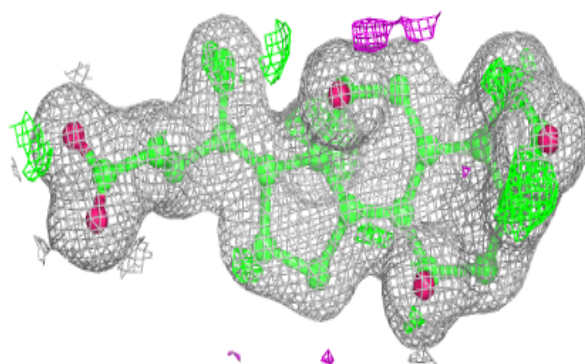


**Electron density around CHD P 307:**

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

**Electron density around CHD O 302:**

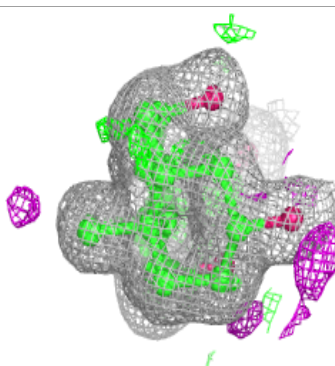
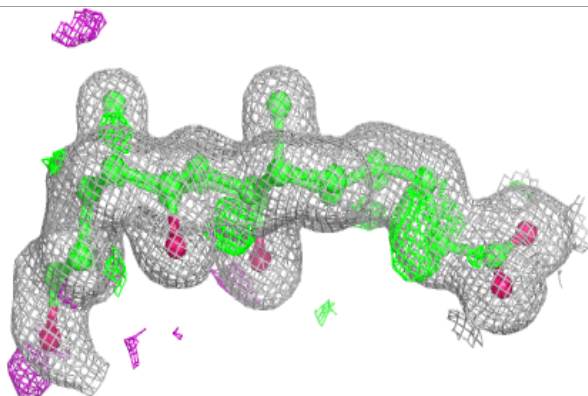
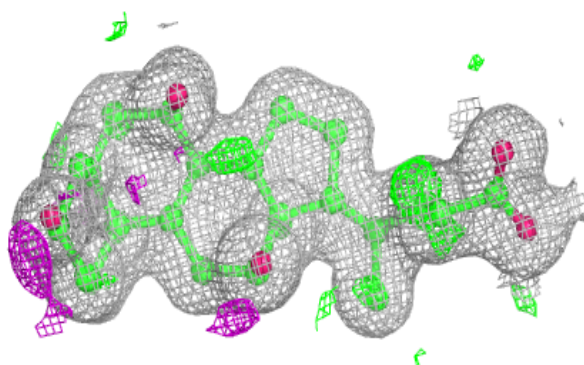
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



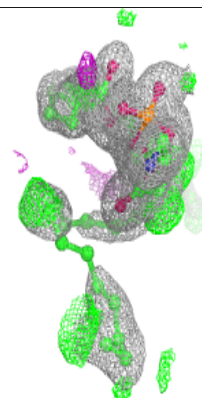
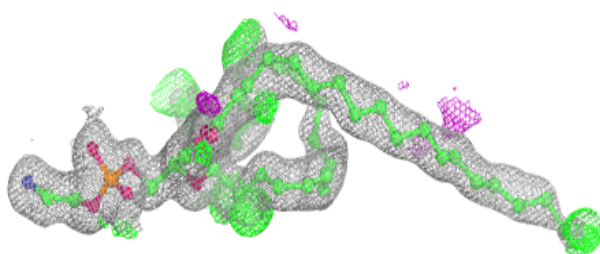
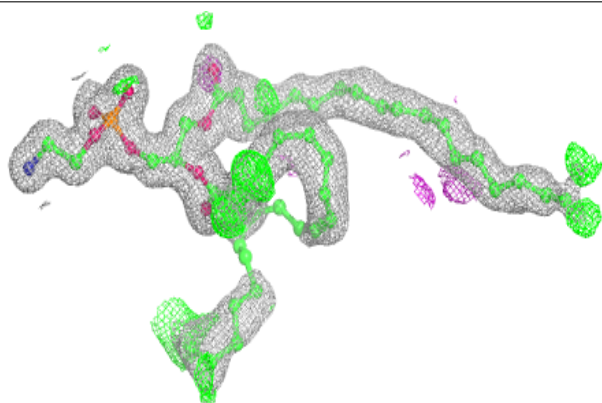


**Electron density around CHD B 303:**

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

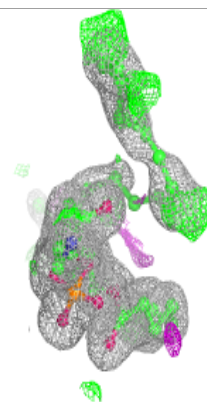
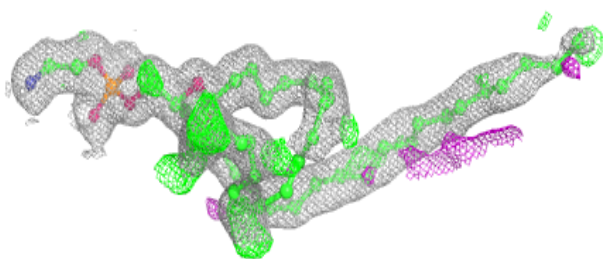
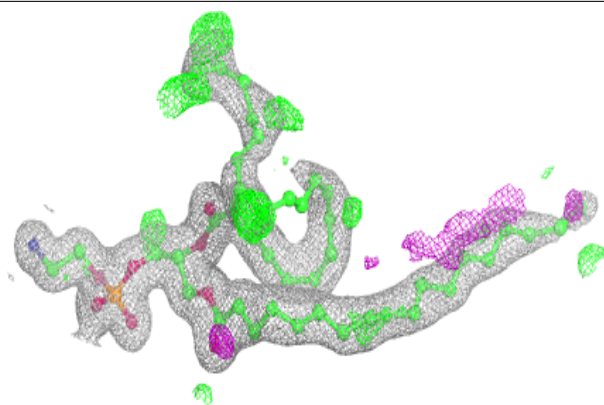
**Electron density around PEK T 101:**

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

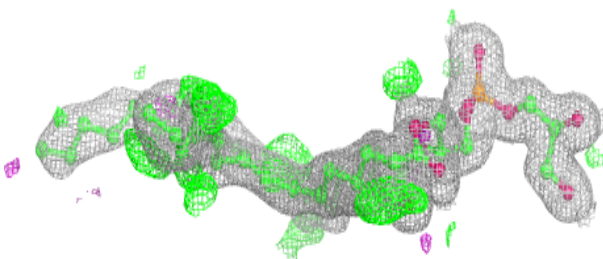
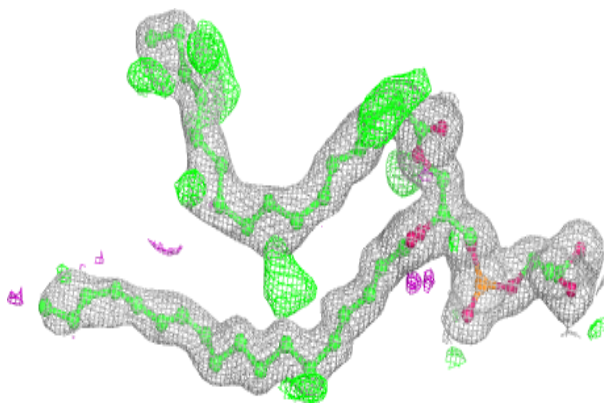


**Electron density around PEK G 101:**

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

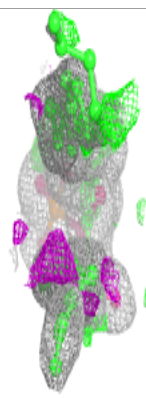
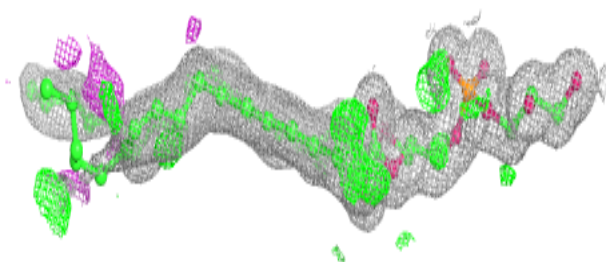
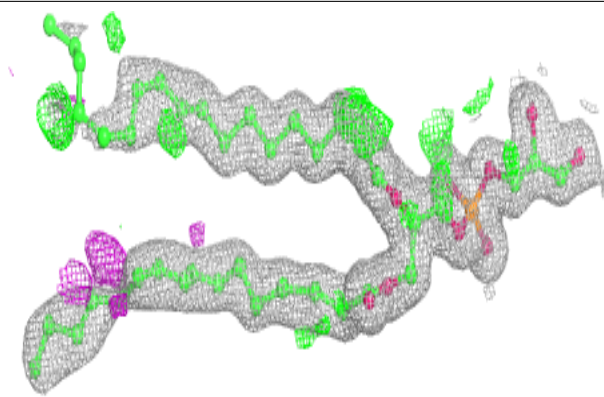
**Electron density around PGV N 607:**

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

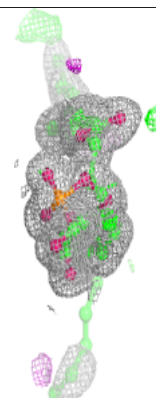
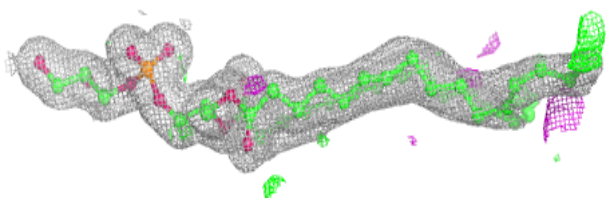
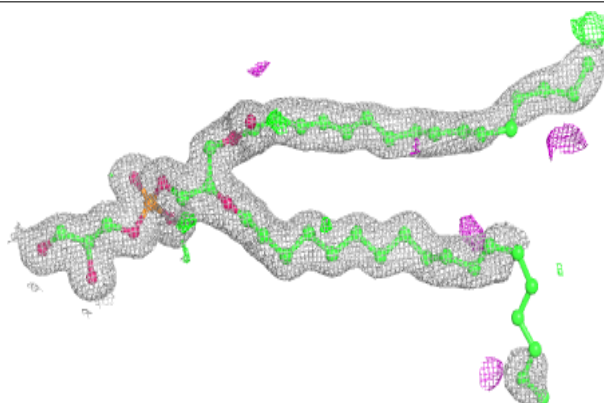


**Electron density around PGV C 302:**

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

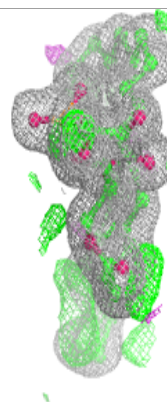
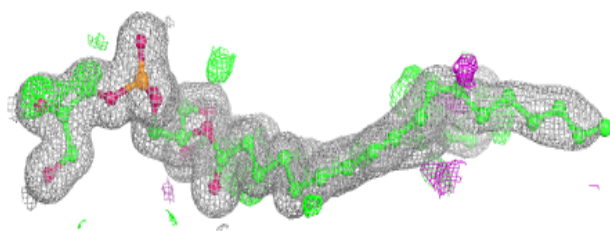
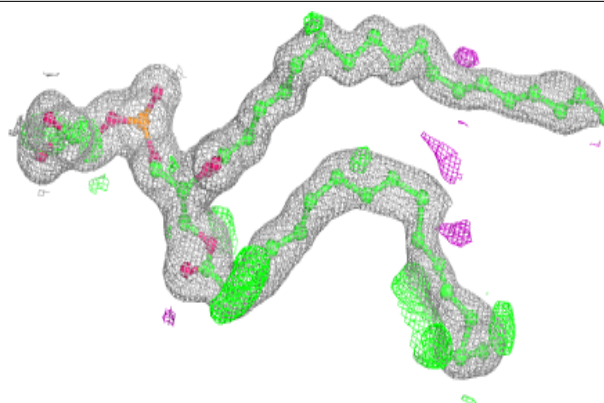
**Electron density around PGV P 303:**

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

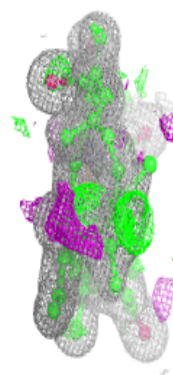
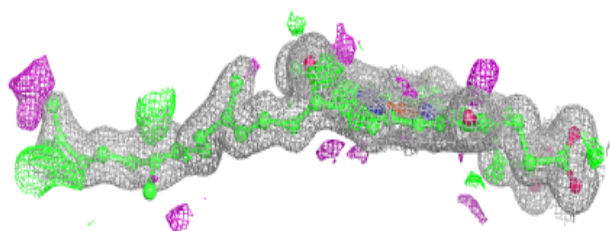
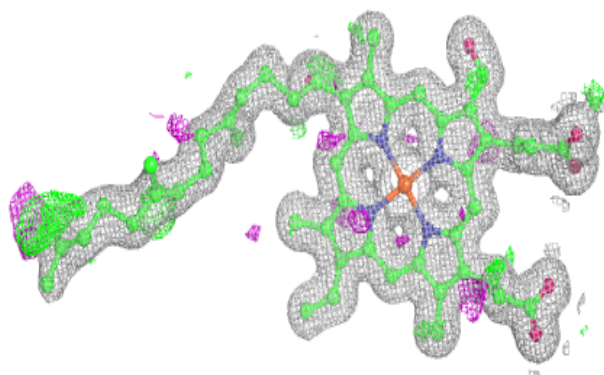


**Electron density around PGV A 607:**

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

**Electron density around HEA A 601:**

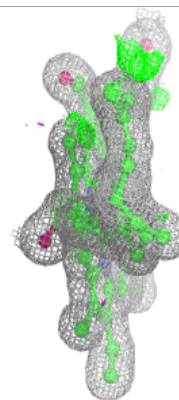
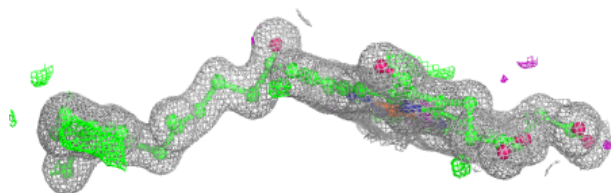
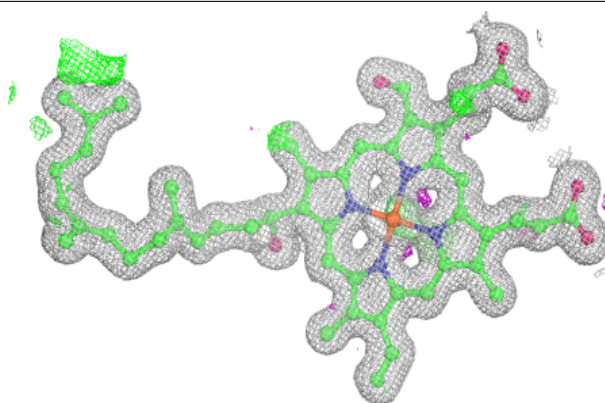
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



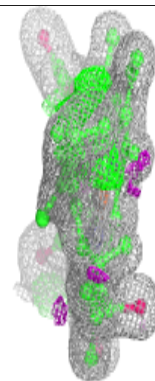
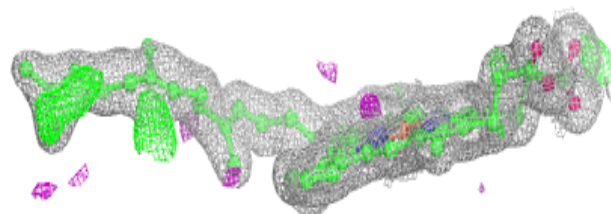
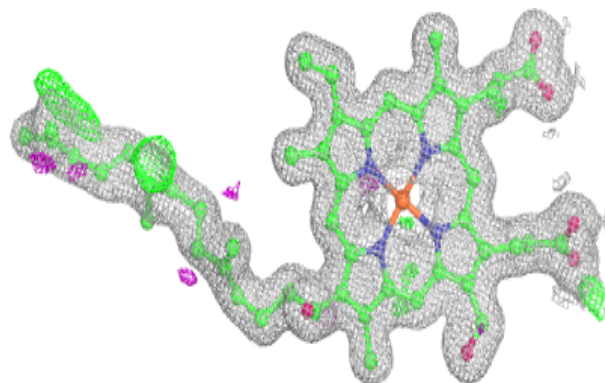


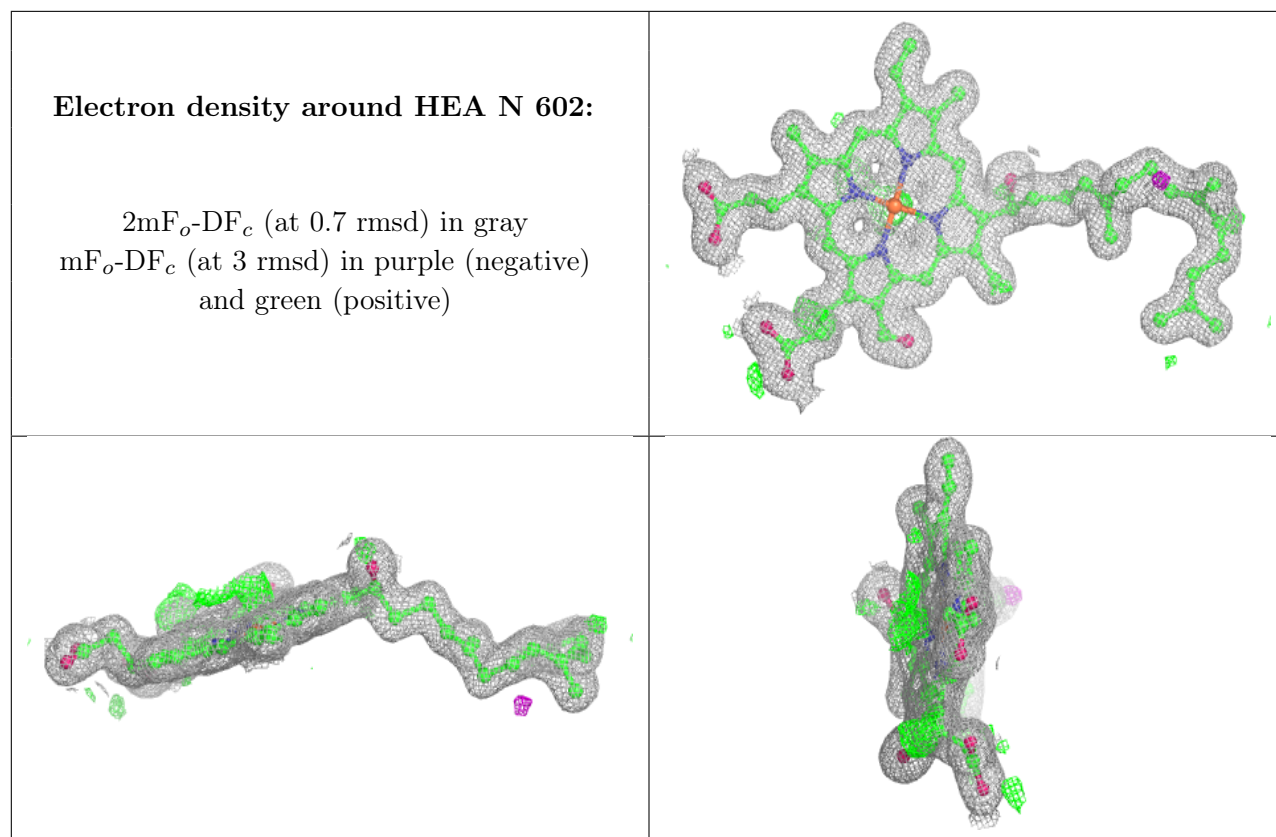
**Electron density around HEA A 602:**

$2mF_o-DF_c$  (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)





## 6.5 Other polymers [i](#)

There are no such residues in this entry.