



wwPDB X-ray Structure Validation Summary Report ⓘ

May 14, 2020 – 11:04 pm BST

PDB ID : 5XDX
Title : Bovine heart cytochrome c oxidase in the reduced state with pH 7.3 at 1.99 angstrom resolution
Authors : Luo, F.J.; Shimada, A.; Hagimoto, N.; Shimada, S.; Shinzawa-Itoh, K.; Yamashita, E.; Yoshikawa, S.; Tsukihara, T.
Deposited on : 2017-03-30
Resolution : 1.99 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) 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 : 2.11
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) : 2.11

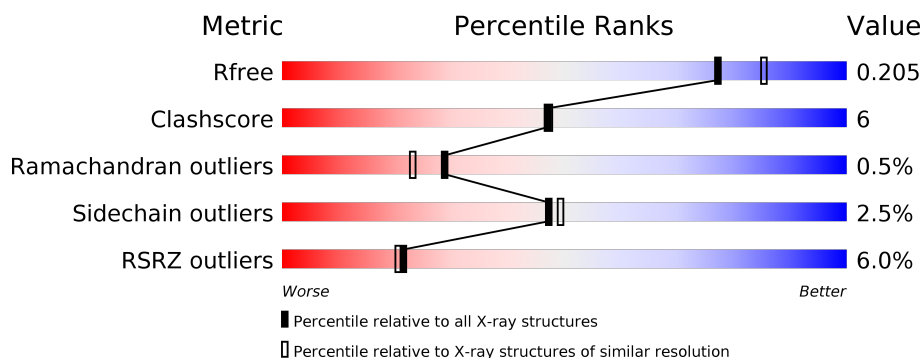
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.99 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	514	<div> <div style="width: 100%;"></div> <div> % 88% 12% • </div> </div>
1	N	514	<div> <div style="width: 100%;"></div> <div> % 87% 12% • </div> </div>
2	B	227	<div> <div style="width: 100%;"></div> <div> 4% 85% 15% • </div> </div>
2	O	227	<div> <div style="width: 100%;"></div> <div> 7% 82% 17% • </div> </div>
3	C	260	<div> <div style="width: 100%;"></div> <div> 3% 85% 13% • </div> </div>
3	P	260	<div> <div style="width: 100%;"></div> <div> 2% 90% 10% • </div> </div>

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Mol	Chain	Length	Quality of chain
4	D	147	
4	Q	147	
5	E	109	
5	R	109	
6	F	94	
6	S	94	
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
14	HEA	A	601	X	-	-	-
14	HEA	A	602	X	-	-	-
14	HEA	N	601	X	-	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
14	HEA	N	602	X	-	-	-
19	PGV	N	606	-	-	-	X
20	EDO	N	613	-	-	X	-
25	DMU	C	313	-	-	-	X
25	DMU	P	309	-	-	-	X
9	SAC	V	1	-	-	-	X

2 Entry composition

There are 28 unique types of molecules in this entry. The entry contains 33247 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	20	0
			4128	2774	626	685	43			
1	N	514	Total	C	N	O	S	0	21	0
			4132	2773	627	687	45			

- 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	7	0
			1854	1209	281	343	21			
2	O	227	Total	C	N	O	S	0	6	0
			1849	1205	281	343	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	13	0
			2173	1461	338	358	16			
3	P	259	Total	C	N	O	S	0	12	0
			2168	1457	337	357	17			

- 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	0	0
			1195	777	196	218	4			
4	Q	144	Total	C	N	O	S	0	0	0
			1195	777	196	218	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	1	0
			857	548	144	162	3			
5	R	105	Total	C	N	O	S	0	1	0
			857	548	144	162	3			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	94	Total	C	N	O	S	0	1	0
			721	448	127	140	6			
6	S	94	Total	C	N	O	S	0	1	0
			721	448	127	140	6			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
7	G	84	Total 675	C 431	N 129	O 113	P 1	S 1	0	0	0
7	T	84	Total 675	C 431	N 129	O 113	P 1	S 1	0	0	0

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

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

- Molecule 9 is a protein called Cytochrome c oxidase polypeptide VIc.

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 VIIa-heart.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	J	58	Total	C	N	O	S	0	2	0
			471	305	78	84	4			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	W	58	Total	C	N	O	S	0	2	0
			471	305	78	84	4			

- Molecule 11 is a protein called Cytochrome c oxidase subunit VIIb.

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	0	0
			384	250	65	67	2			

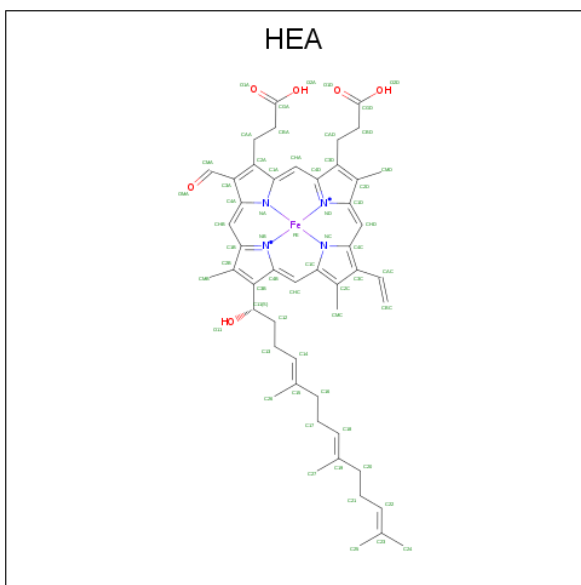
- Molecule 12 is a protein called Cytochrome c oxidase subunit VIIc.

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
			385	258	64	60	3			

- Molecule 13 is a protein called Cytochrome c oxidase subunit VIII-heart.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
13	M	43	Total	C	N	O	0	1	0
			340	228	53	59			
13	Z	43	Total	C	N	O	0	1	0
			340	228	53	59			

- Molecule 14 is HEME-A (three-letter code: HEA) (formula: C₄₉H₅₆FeN₄O₆).



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		

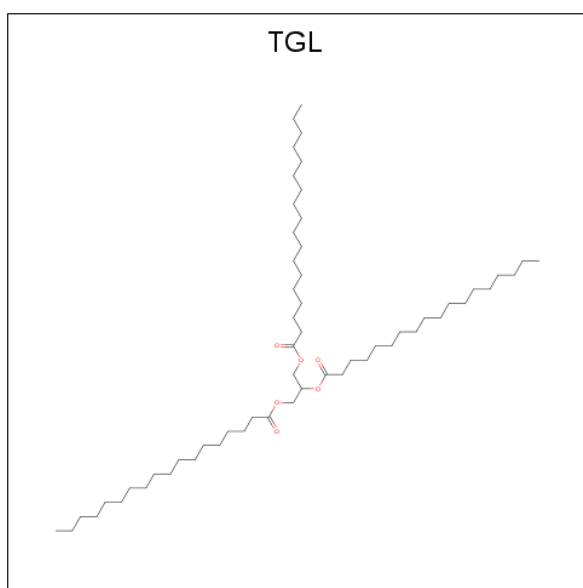
- Molecule 16 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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).

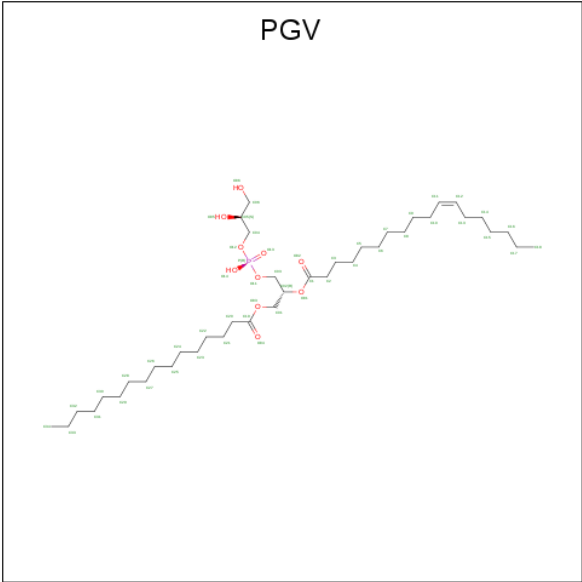
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
17	P	1	Total Na 1 1	0	0
17	A	1	Total Na 1 1	0	0
17	C	1	Total Na 1 1	0	0
17	N	1	Total Na 1 1	0	0

- Molecule 18 is TRISTEAROYLGLYCEROL (three-letter code: TGL) (formula: C₅₇H₁₁₀O₆).



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

- 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: C₄₀H₇₇O₁₀P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
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	M	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	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		

- Molecule 20 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



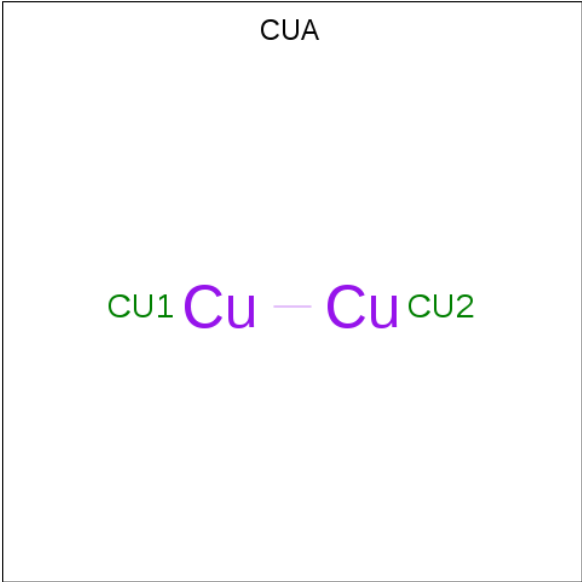
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
20	A	1	Total	C	O	0	0
			4	2	2		
20	A	1	Total	C	O	0	0
			4	2	2		
20	A	1	Total	C	O	0	0
			4	2	2		
20	A	1	Total	C	O	0	0
			4	2	2		
20	A	1	Total	C	O	0	0
			4	2	2		
20	A	1	Total	C	O	0	0
			4	2	2		
20	A	1	Total	C	O	0	0
			4	2	2		
20	B	1	Total	C	O	0	0
			4	2	2		
20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

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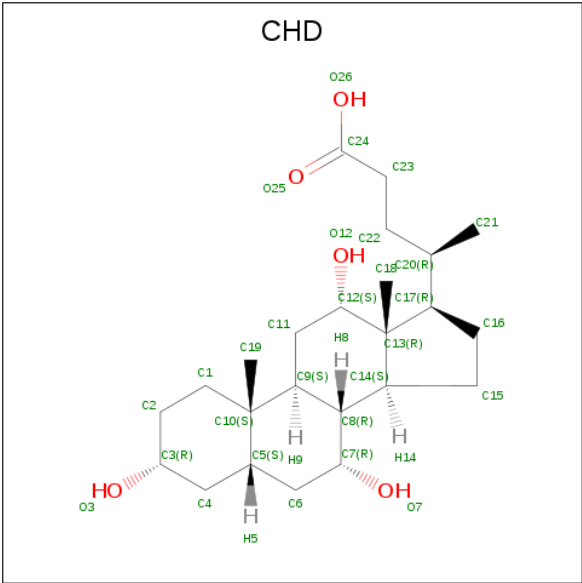
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
20	D	1	Total 4	C 2	O 2	0	0
20	E	1	Total 4	C 2	O 2	0	0
20	F	1	Total 4	C 2	O 2	0	0
20	F	1	Total 4	C 2	O 2	0	0
20	F	1	Total 4	C 2	O 2	0	0
20	N	1	Total 4	C 2	O 2	0	0
20	N	1	Total 4	C 2	O 2	0	0
20	N	1	Total 4	C 2	O 2	0	0
20	N	1	Total 4	C 2	O 2	0	0
20	N	1	Total 4	C 2	O 2	0	0
20	N	1	Total 4	C 2	O 2	0	0
20	N	1	Total 4	C 2	O 2	0	0
20	N	1	Total 4	C 2	O 2	0	0
20	N	1	Total 4	C 2	O 2	0	0
20	N	1	Total 4	C 2	O 2	0	0
20	N	1	Total 4	C 2	O 2	0	0
20	P	1	Total 4	C 2	O 2	0	0
20	R	1	Total 4	C 2	O 2	0	0
20	S	1	Total 4	C 2	O 2	0	0
20	S	1	Total 4	C 2	O 2	0	0

- Molecule 21 is DINUCLEAR COPPER ION (three-letter code: CUA) (formula: Cu₂).



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: C₂₄H₄₀O₅).



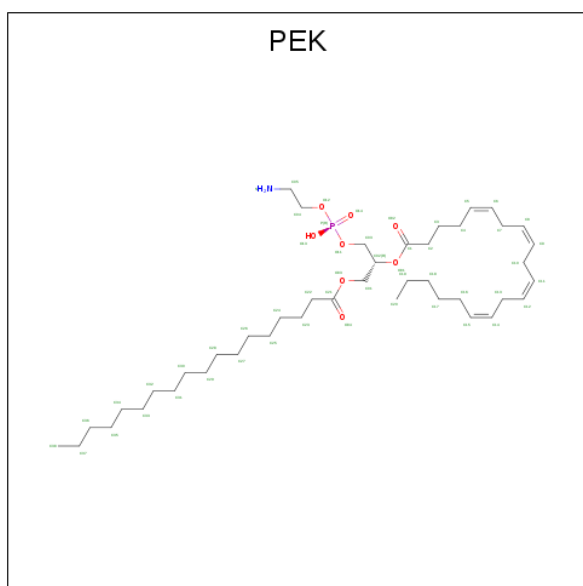
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
22	B	1	Total	C	O	0	0
			29	24	5		
22	C	1	Total	C	O	0	0
			29	24	5		

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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 (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: C₄₃H₇₈NO₈P).



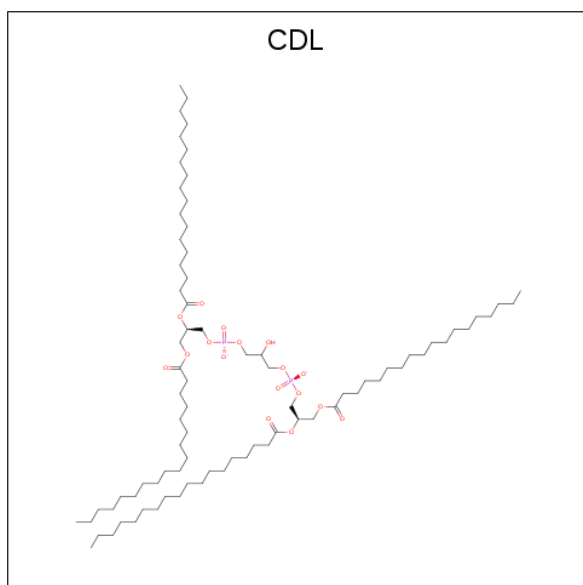
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
23	C	1	Total	C	N	O	P	0	0
			53	43	1	8	1		
23	C	1	Total	C	N	O	P	0	0
			53	43	1	8	1		
23	G	1	Total	C	N	O	P	0	0
			53	43	1	8	1		
23	T	1	Total	C	N	O	P	0	0
			53	43	1	8	1		

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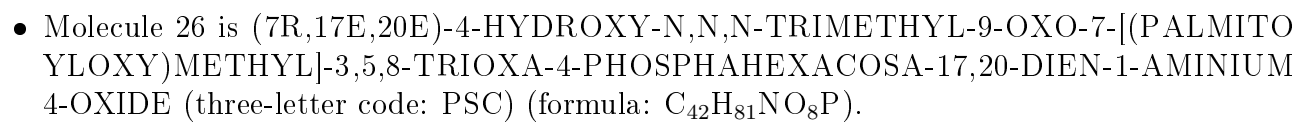
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
23	T	1	Total	C	N	O	P	0	0
			53	43	1	8	1		
23	T	1	Total	C	N	O	P	0	0
			53	43	1	8	1		

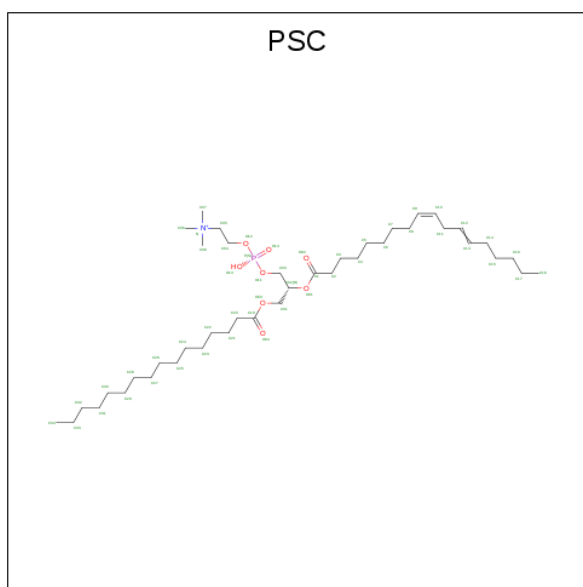
- Molecule 24 is CARDIOLIPIN (three-letter code: CDL) (formula: $C_{81}H_{156}O_{17}P_2$).



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

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





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

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

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

- Molecule 28 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
28	A	245	Total	O	0	0
			245	245		
28	B	177	Total	O	0	1
			178	178		
28	C	103	Total	O	0	0
			103	103		
28	D	130	Total	O	0	0
			130	130		
28	E	92	Total	O	0	0
			92	92		

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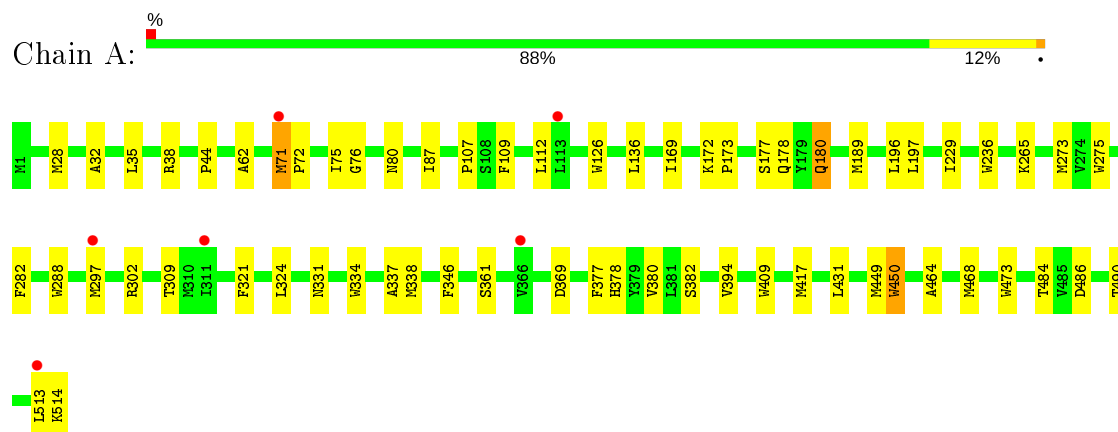
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
28	F	92	Total 92	O 92	0	0
28	G	58	Total 58	O 58	0	0
28	H	56	Total 56	O 56	0	0
28	I	48	Total 48	O 48	0	0
28	J	28	Total 28	O 28	0	0
28	K	42	Total 42	O 42	0	0
28	L	26	Total 26	O 26	0	0
28	M	27	Total 27	O 27	0	0
28	N	214	Total 214	O 214	0	0
28	O	128	Total 128	O 128	0	0
28	P	99	Total 99	O 99	0	0
28	Q	58	Total 58	O 58	0	0
28	R	62	Total 62	O 62	0	0
28	S	64	Total 64	O 64	0	0
28	T	51	Total 51	O 51	0	0
28	U	45	Total 45	O 45	0	0
28	V	31	Total 31	O 31	0	0
28	W	19	Total 19	O 19	0	0
28	X	11	Total 11	O 11	0	0
28	Y	21	Total 21	O 21	0	0
28	Z	14	Total 14	O 14	0	0

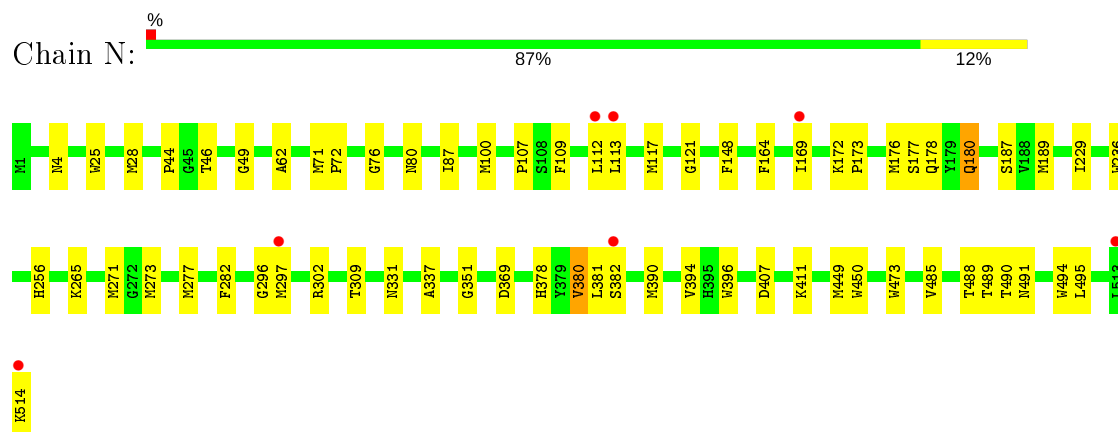
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

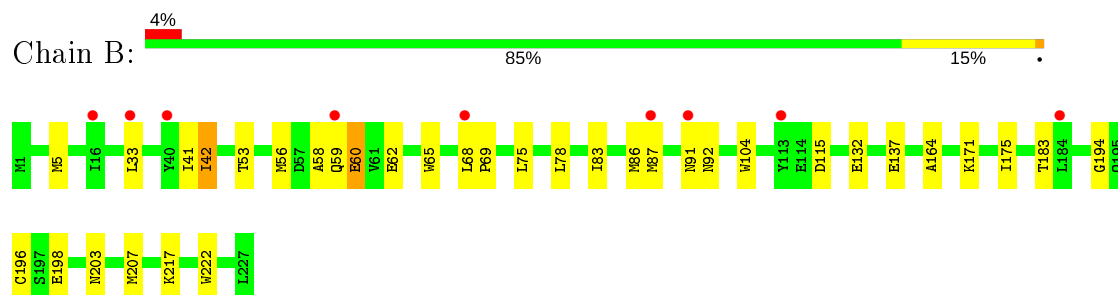
• Molecule 1: Cytochrome c oxidase subunit 1



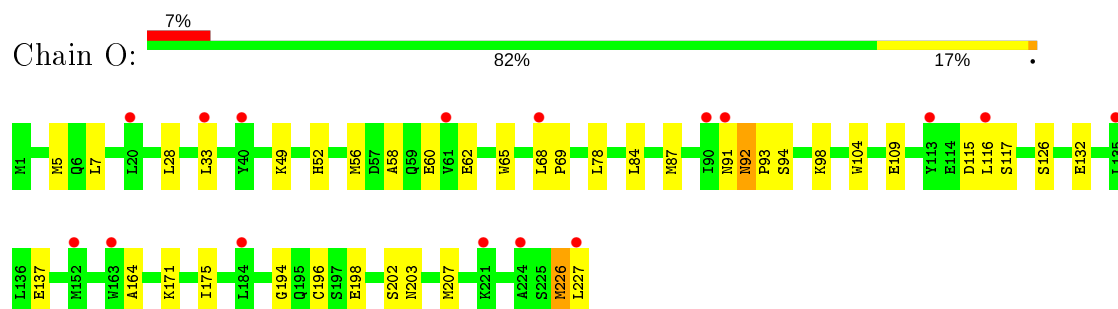
• Molecule 1: Cytochrome c oxidase subunit 1



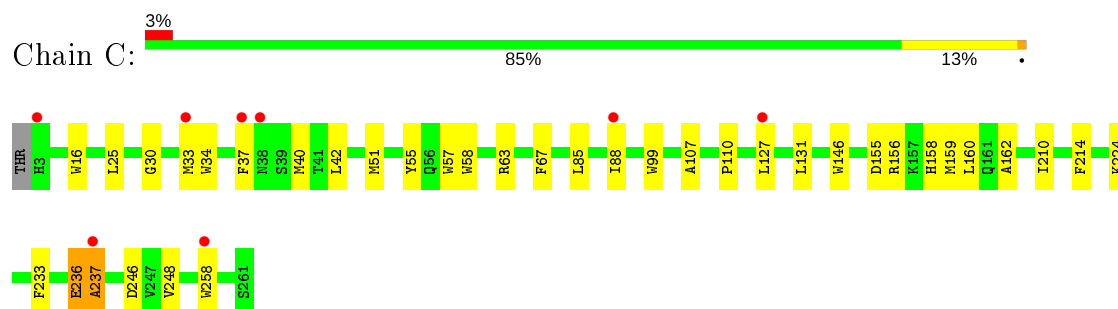
• Molecule 2: Cytochrome c oxidase subunit 2



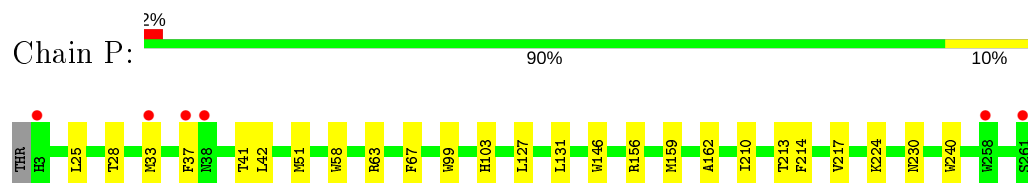
- Molecule 2: Cytochrome c oxidase subunit 2



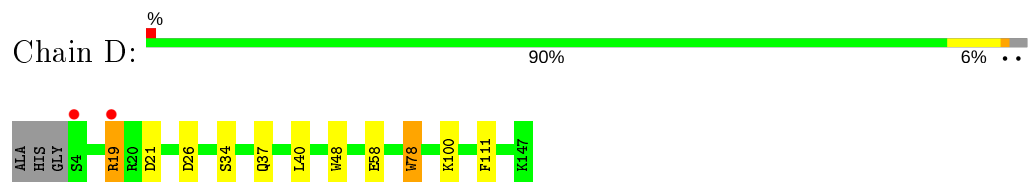
- Molecule 3: Cytochrome c oxidase subunit 3



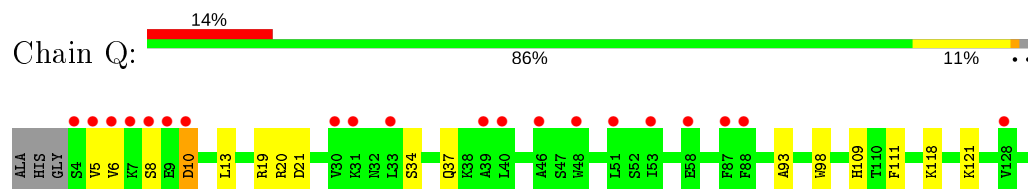
- Molecule 3: Cytochrome c oxidase subunit 3



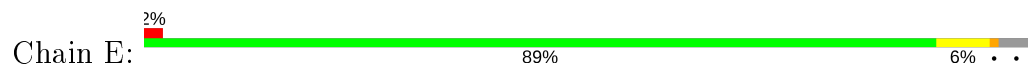
- Molecule 4: Cytochrome c oxidase subunit 4 isoform 1, mitochondrial

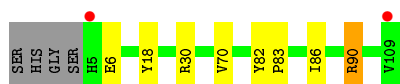


- Molecule 4: Cytochrome c oxidase subunit 4 isoform 1, mitochondrial

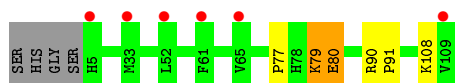
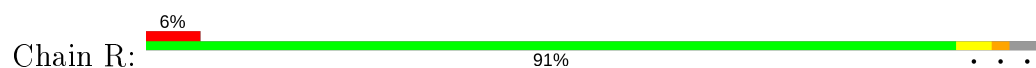


- Molecule 5: Cytochrome c oxidase subunit 5A, mitochondrial

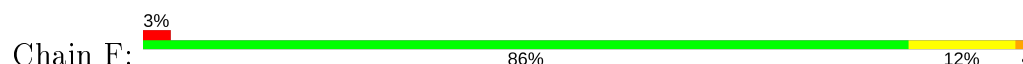




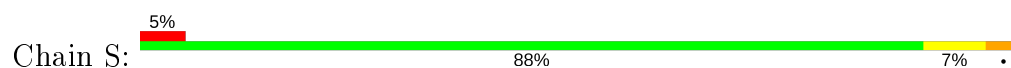
- Molecule 5: Cytochrome c oxidase subunit 5A, mitochondrial



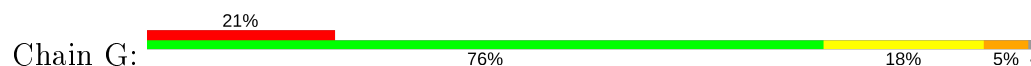
- Molecule 6: Cytochrome c oxidase subunit 5B, mitochondrial



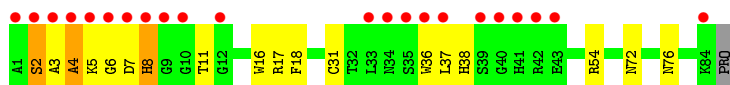
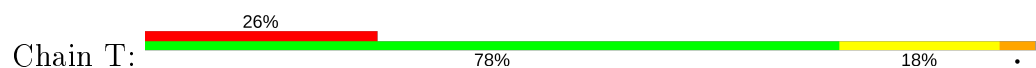
- Molecule 6: Cytochrome c oxidase subunit 5B, mitochondrial



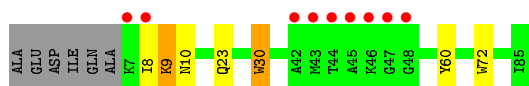
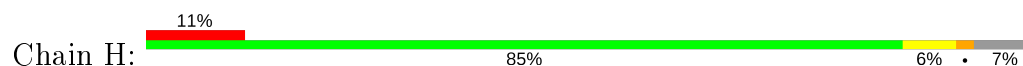
- Molecule 7: Cytochrome c oxidase subunit 6A2, mitochondrial



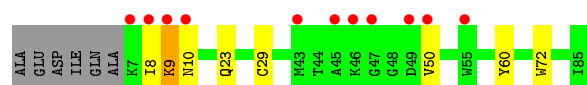
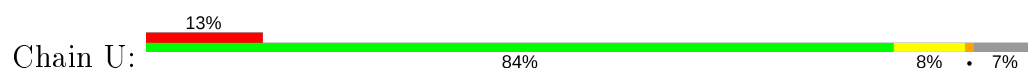
- Molecule 7: Cytochrome c oxidase subunit 6A2, mitochondrial



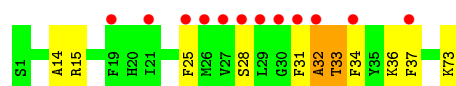
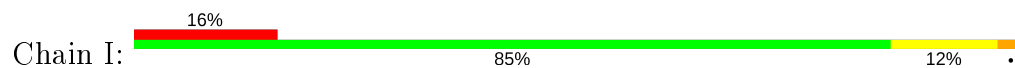
- Molecule 8: Cytochrome c oxidase subunit 6B1



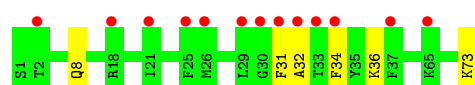
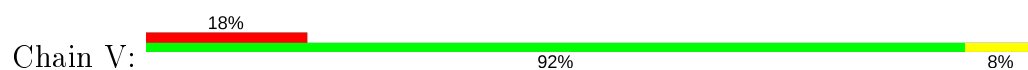
- Molecule 8: Cytochrome c oxidase subunit 6B1



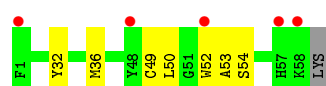
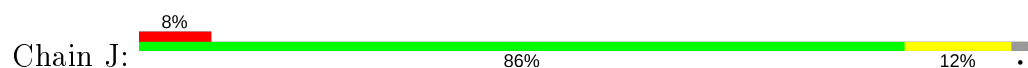
- Molecule 9: Cytochrome c oxidase polypeptide VIc



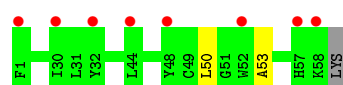
- Molecule 9: Cytochrome c oxidase polypeptide VIc



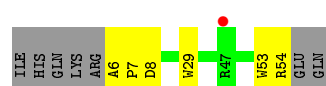
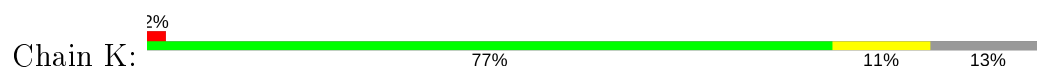
- Molecule 10: Cytochrome c oxidase subunit VIIa-heart



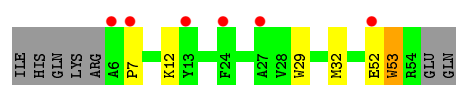
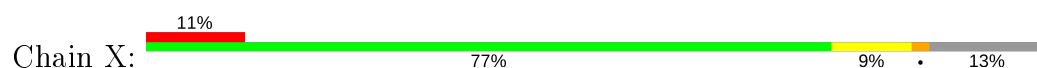
- Molecule 10: Cytochrome c oxidase subunit VIIa-heart



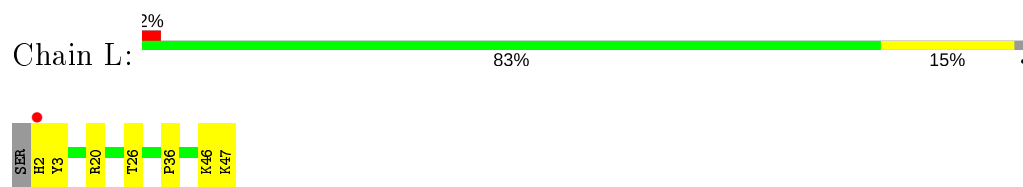
- Molecule 11: Cytochrome c oxidase subunit VIIb



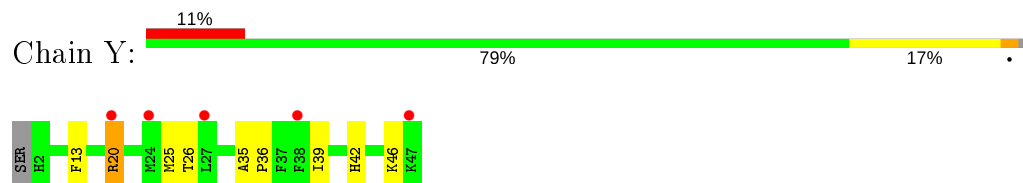
- Molecule 11: Cytochrome c oxidase subunit VIIb



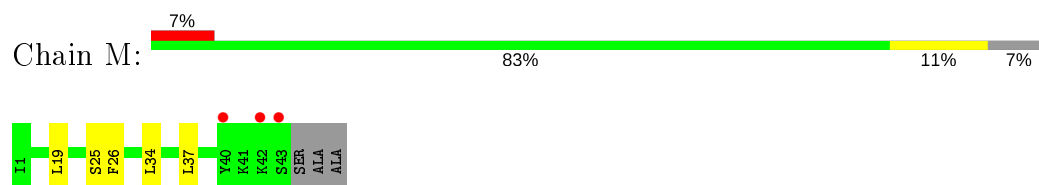
- Molecule 12: Cytochrome c oxidase subunit VIIc



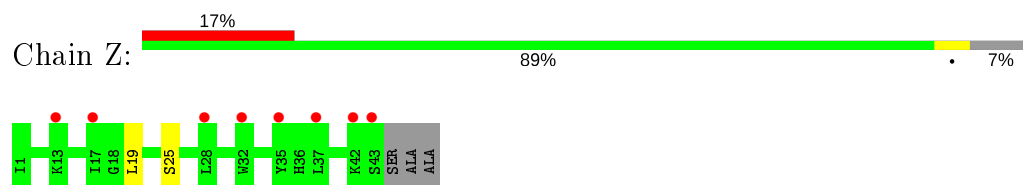
- Molecule 12: Cytochrome c oxidase subunit VIIc



- Molecule 13: Cytochrome c oxidase subunit VIII-heart



- Molecule 13: Cytochrome c oxidase subunit VIII-heart



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	183.04Å 205.69Å 177.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 1.99 108.39 – 1.99	Depositor EDS
% Data completeness (in resolution range)	99.7 (40.00-1.99) 99.7 (108.39-1.99)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.30 (at 1.98Å)	Xtriage
Refinement program	REFMAC 5.8.0048	Depositor
R, R_{free}	0.182 , 0.203 0.183 , 0.205	Depositor DCC
R_{free} test set	23139 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	36.0	Xtriage
Anisotropy	0.540	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 62.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.003 for l,-k,h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	33247	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.13% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.66	6/4315 (0.1%)	0.61	2/5889 (0.0%)
1	N	0.63	4/4322 (0.1%)	0.59	0/5896
2	B	0.61	3/1911 (0.2%)	0.66	0/2602
2	O	0.53	2/1903 (0.1%)	0.61	0/2592
3	C	0.71	5/2299 (0.2%)	0.58	1/3140 (0.0%)
3	P	0.69	3/2291 (0.1%)	0.53	0/3128
4	D	0.63	2/1229 (0.2%)	0.54	0/1658
4	Q	0.58	2/1229 (0.2%)	0.53	0/1658
5	E	0.52	0/879	0.55	0/1192
5	R	0.49	0/879	0.54	0/1192
6	F	0.54	0/740	0.57	0/1003
6	S	0.54	0/740	0.55	0/1003
7	G	0.68	1/690 (0.1%)	0.60	0/937
7	T	0.69	2/690 (0.3%)	0.61	0/937
8	H	0.63	2/682 (0.3%)	0.55	0/921
8	U	0.62	1/682 (0.1%)	0.57	0/921
9	I	0.45	0/605	0.56	0/802
9	V	0.38	0/605	0.50	0/802
10	J	0.52	0/488	0.52	0/658
10	W	0.52	0/488	0.50	0/658
11	K	0.70	2/398 (0.5%)	0.53	0/546
11	X	0.67	2/398 (0.5%)	0.50	0/546
12	L	0.60	0/393	0.55	0/526
12	Y	0.54	0/401	0.49	0/536
13	M	0.58	0/353	0.53	0/481
13	Z	0.54	0/353	0.50	0/481
All	All	0.62	37/29963 (0.1%)	0.57	3/40705 (0.0%)

The worst 5 of 37 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	198	GLU	C-O	6.34	1.35	1.23
1	A	126	TRP	CD2-CE2	5.79	1.48	1.41
2	B	65	TRP	CD2-CE2	5.54	1.48	1.41
8	H	72	TRP	CD2-CE2	5.51	1.48	1.41
11	K	29	TRP	CD2-CE2	5.46	1.47	1.41

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	71[A]	MET	CG-SD-CE	-5.71	91.06	100.20
1	A	71[B]	MET	CG-SD-CE	-5.71	91.06	100.20
3	C	236	GLU	C-N-CA	-5.35	108.33	121.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4128	0	4171	59	0
1	N	4132	0	4169	75	0
2	B	1854	0	1888	23	0
2	O	1849	0	1879	21	0
3	C	2173	0	2130	34	0
3	P	2168	0	2122	23	0
4	D	1195	0	1183	14	0
4	Q	1195	0	1183	11	0
5	E	857	0	854	4	0
5	R	857	0	854	5	0
6	F	721	0	706	13	0
6	S	721	0	706	11	0
7	G	675	0	644	16	0
7	T	675	0	643	15	0
8	H	662	0	623	6	0
8	U	662	0	623	2	0
9	I	601	0	613	9	0
9	V	601	0	613	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	J	471	0	474	9	0
10	W	471	0	474	3	0
11	K	384	0	366	2	0
11	X	384	0	366	5	0
12	L	380	0	380	9	0
12	Y	385	0	389	10	0
13	M	340	0	363	5	0
13	Z	340	0	363	2	0
14	A	120	0	108	6	0
14	N	120	0	108	7	0
15	A	1	0	0	0	0
15	N	1	0	0	0	0
16	A	1	0	0	0	0
16	N	1	0	0	0	0
17	A	1	0	0	0	0
17	C	1	0	0	0	0
17	N	1	0	0	0	0
17	P	1	0	0	0	0
18	A	63	0	110	3	0
18	D	63	0	110	4	0
18	L	63	0	110	4	0
18	O	63	0	110	2	0
18	Q	63	0	110	0	0
18	Y	63	0	110	5	0
19	A	51	0	76	2	0
19	C	102	0	152	4	0
19	M	51	0	76	2	0
19	N	102	0	152	4	0
19	P	102	0	152	4	0
20	A	36	0	54	5	0
20	B	4	0	6	0	0
20	C	12	0	18	0	0
20	D	8	0	12	2	0
20	E	4	0	6	1	0
20	F	12	0	18	0	0
20	N	40	0	60	9	0
20	P	4	0	6	0	0
20	R	4	0	6	0	0
20	S	8	0	12	1	0
21	B	2	0	0	0	0
21	O	2	0	0	0	0
22	B	29	0	39	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
22	C	58	0	78	3	0
22	J	29	0	38	0	0
22	O	29	0	39	1	0
22	P	58	0	78	3	0
22	W	29	0	39	1	0
23	C	106	0	154	5	0
23	G	53	0	77	0	0
23	T	159	0	231	7	0
24	C	100	0	156	10	0
24	G	100	0	156	5	0
24	P	100	0	156	8	0
24	T	100	0	156	13	0
25	C	66	0	84	12	0
25	M	33	0	42	0	0
25	P	66	0	84	9	0
25	Z	33	0	42	0	0
26	E	52	0	80	10	0
26	O	52	0	80	6	0
27	F	1	0	0	0	0
27	S	1	0	0	0	0
28	A	245	0	0	12	0
28	B	178	0	0	5	0
28	C	103	0	0	2	0
28	D	130	0	0	2	0
28	E	92	0	0	0	0
28	F	92	0	0	1	0
28	G	58	0	0	0	0
28	H	56	0	0	2	0
28	I	48	0	0	0	0
28	J	28	0	0	1	0
28	K	42	0	0	1	0
28	L	26	0	0	0	0
28	M	27	0	0	0	0
28	N	214	0	0	6	0
28	O	128	0	0	1	0
28	P	99	0	0	1	0
28	Q	58	0	0	1	0
28	R	62	0	0	0	0
28	S	64	0	0	0	0
28	T	51	0	0	0	0
28	U	45	0	0	0	0
28	V	31	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
28	W	19	0	0	0	0
28	X	11	0	0	3	0
28	Y	21	0	0	0	0
28	Z	14	0	0	0	0
All	All	33247	0	32270	397	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

The worst 5 of 397 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)
3:C:236:GLU:O	3:C:237:ALA:CB	1.79	1.26
1:N:113[B]:LEU:HG	1:N:117[B]:MET:CE	1.69	1.22
20:A:615:EDO:H22	28:A:920:HOH:O	1.42	1.19
3:C:33[B]:MET:SD	25:C:313:DMU:H11	1.83	1.18
1:A:28[B]:MET:HA	1:A:28[B]:MET:CE	1.83	1.09

There are no symmetry-related clashes.

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	532/514 (104%)	520 (98%)	12 (2%)	0	100	100
1	N	533/514 (104%)	520 (98%)	13 (2%)	0	100	100
2	B	232/227 (102%)	227 (98%)	4 (2%)	1 (0%)	34	30
2	O	231/227 (102%)	225 (97%)	5 (2%)	1 (0%)	34	30
3	C	270/260 (104%)	263 (97%)	6 (2%)	1 (0%)	34	30
3	P	269/260 (104%)	262 (97%)	7 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	D	142/147 (97%)	139 (98%)	3 (2%)	0	100	100
4	Q	142/147 (97%)	136 (96%)	5 (4%)	1 (1%)	22	16
5	E	104/109 (95%)	103 (99%)	0	1 (1%)	15	9
5	R	104/109 (95%)	104 (100%)	0	0	100	100
6	F	93/94 (99%)	91 (98%)	2 (2%)	0	100	100
6	S	93/94 (99%)	91 (98%)	2 (2%)	0	100	100
7	G	81/85 (95%)	71 (88%)	6 (7%)	4 (5%)	2	0
7	T	81/85 (95%)	69 (85%)	7 (9%)	5 (6%)	1	0
8	H	77/85 (91%)	74 (96%)	2 (3%)	1 (1%)	12	6
8	U	77/85 (91%)	72 (94%)	5 (6%)	0	100	100
9	I	71/73 (97%)	68 (96%)	2 (3%)	1 (1%)	11	5
9	V	71/73 (97%)	70 (99%)	1 (1%)	0	100	100
10	J	58/59 (98%)	57 (98%)	1 (2%)	0	100	100
10	W	58/59 (98%)	57 (98%)	1 (2%)	0	100	100
11	K	47/56 (84%)	46 (98%)	1 (2%)	0	100	100
11	X	47/56 (84%)	46 (98%)	1 (2%)	0	100	100
12	L	44/47 (94%)	42 (96%)	2 (4%)	0	100	100
12	Y	45/47 (96%)	43 (96%)	1 (2%)	1 (2%)	6	2
13	M	42/46 (91%)	41 (98%)	1 (2%)	0	100	100
13	Z	42/46 (91%)	41 (98%)	1 (2%)	0	100	100
All	All	3586/3604 (100%)	3478 (97%)	91 (2%)	17 (0%)	29	23

5 of 17 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	237	ALA
7	G	4	ALA
7	G	8	HIS
4	Q	8	SER
7	T	4	ALA

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar

resolution.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	446/426 (105%)	439 (98%)	7 (2%)	62	67
1	N	447/426 (105%)	441 (99%)	6 (1%)	69	74
2	B	217/210 (103%)	208 (96%)	9 (4%)	30	28
2	O	216/210 (103%)	207 (96%)	9 (4%)	30	27
3	C	237/225 (105%)	234 (99%)	3 (1%)	69	74
3	P	236/225 (105%)	233 (99%)	3 (1%)	69	74
4	D	128/129 (99%)	127 (99%)	1 (1%)	81	86
4	Q	128/129 (99%)	125 (98%)	3 (2%)	50	53
5	E	93/95 (98%)	91 (98%)	2 (2%)	52	55
5	R	93/95 (98%)	91 (98%)	2 (2%)	52	55
6	F	79/78 (101%)	77 (98%)	2 (2%)	47	49
6	S	79/78 (101%)	74 (94%)	5 (6%)	18	13
7	G	67/68 (98%)	60 (90%)	7 (10%)	7	4
7	T	67/68 (98%)	63 (94%)	4 (6%)	19	14
8	H	71/75 (95%)	69 (97%)	2 (3%)	43	44
8	U	71/75 (95%)	66 (93%)	5 (7%)	15	10
9	I	57/57 (100%)	54 (95%)	3 (5%)	22	18
9	V	57/57 (100%)	56 (98%)	1 (2%)	59	63
10	J	51/50 (102%)	51 (100%)	0	100	100
10	W	51/50 (102%)	50 (98%)	1 (2%)	55	58
11	K	39/46 (85%)	38 (97%)	1 (3%)	46	48
11	X	39/46 (85%)	39 (100%)	0	100	100
12	L	39/40 (98%)	39 (100%)	0	100	100
12	Y	40/40 (100%)	39 (98%)	1 (2%)	47	49
13	M	38/38 (100%)	38 (100%)	0	100	100
13	Z	38/38 (100%)	38 (100%)	0	100	100
All	All	3124/3074 (102%)	3047 (98%)	77 (2%)	47	49

5 of 77 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
9	I	25	PHE
2	O	33	LEU
8	U	29	CYS
9	I	33	THR
1	N	369	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 37 such sidechains are listed below:

Mol	Chain	Res	Type
7	G	76	ASN
2	O	181	GLN
11	X	35	GLN
1	N	180	GLN
2	O	10	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

8 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
9	SAC	V	1	9	7,8,9	2.09	2 (28%)	8,9,11	1.21	0
7	TPO	G	11	7	8,10,11	1.39	1 (12%)	10,14,16	0.83	0
2	FME	O	1	2	8,9,10	0.59	0	7,9,11	1.12	0
7	TPO	T	11	7	8,10,11	1.40	1 (12%)	10,14,16	0.87	0
9	SAC	I	1	9	7,8,9	2.00	2 (28%)	8,9,11	1.11	0
1	FME	A	1	1	8,9,10	0.57	0	7,9,11	1.71	2 (28%)
1	FME	N	1	1	8,9,10	0.51	0	7,9,11	1.33	2 (28%)
2	FME	B	1	2	8,9,10	0.76	0	7,9,11	1.28	1 (14%)

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
9	SAC	V	1	9	-	2/7/8/10	-
7	TPO	G	11	7	-	4/9/11/13	-
2	FME	O	1	2	-	0/7/9/11	-
7	TPO	T	11	7	-	5/9/11/13	-
9	SAC	I	1	9	-	2/7/8/10	-
1	FME	A	1	1	-	4/7/9/11	-
1	FME	N	1	1	-	2/7/9/11	-
2	FME	B	1	2	-	0/7/9/11	-

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	I	1	SAC	OAC-C1A	4.62	1.33	1.23
9	V	1	SAC	OAC-C1A	4.59	1.33	1.23
9	V	1	SAC	CA-N	2.93	1.50	1.46
7	G	11	TPO	P-O1P	2.80	1.59	1.50
7	T	11	TPO	P-O1P	2.76	1.59	1.50

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1	FME	C-CA-N	3.18	115.47	109.73
1	N	1	FME	C-CA-N	2.22	113.74	109.73
1	N	1	FME	O-C-CA	-2.08	119.32	124.78
2	B	1	FME	O-C-CA	-2.07	119.36	124.78
1	A	1	FME	O-C-CA	-2.03	119.45	124.78

There are no chirality outliers.

5 of 19 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	G	11	TPO	N-CA-CB-CG2
7	G	11	TPO	N-CA-CB-OG1
7	G	11	TPO	C-CA-CB-CG2
7	T	11	TPO	N-CA-CB-CG2
7	T	11	TPO	N-CA-CB-OG1

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	T	11	TPO	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 89 ligands modelled in this entry, 10 are monoatomic - leaving 79 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	$\# Z > 2$
20	EDO	A	615	-	3,3,3	0.42	0	2,2,2	0.43	0
18	TGL	Y	101	-	62,62,62	1.25	6 (9%)	65,65,65	1.02	3 (4%)
26	PSC	E	201	-	51,51,51	1.11	3 (5%)	57,59,59	1.01	2 (3%)
23	PEK	T	103	-	52,52,52	0.97	2 (3%)	55,57,57	0.87	3 (5%)
18	TGL	O	302	-	62,62,62	1.25	6 (9%)	65,65,65	1.10	3 (4%)
20	EDO	A	609	-	3,3,3	0.55	0	2,2,2	0.51	0
20	EDO	S	102	-	3,3,3	0.46	0	2,2,2	0.39	0
20	EDO	A	608	-	3,3,3	0.42	0	2,2,2	0.40	0
22	CHD	J	101	-	29,32,32	0.48	0	48,51,51	1.62	9 (18%)
24	CDL	C	307	-	99,99,99	1.33	12 (12%)	105,111,111	1.10	4 (3%)
19	PGV	A	607	-	50,50,50	0.88	2 (4%)	53,56,56	0.91	3 (5%)
21	CUA	O	303	2	0,1,1	0.00	-	-	-	-
22	CHD	C	308	-	29,32,32	0.52	0	48,51,51	1.62	11 (22%)
19	PGV	C	305	-	50,50,50	0.85	2 (4%)	53,56,56	0.81	1 (1%)
20	EDO	N	616	-	3,3,3	0.46	0	2,2,2	0.32	0
20	EDO	D	202	-	3,3,3	0.50	0	2,2,2	0.25	0
20	EDO	E	202	-	3,3,3	0.43	0	2,2,2	0.33	0
23	PEK	T	101	-	52,52,52	0.95	2 (3%)	55,57,57	0.93	2 (3%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
23	PEK	C	303	-	52,52,52	0.83	2 (3%)	55,57,57	0.86	1 (1%)
24	CDL	P	305	-	99,99,99	1.32	12 (12%)	105,111,111	1.10	7 (6%)
20	EDO	P	307	-	3,3,3	0.46	0	2,2,2	0.33	0
20	EDO	C	310	-	3,3,3	0.48	0	2,2,2	0.35	0
23	PEK	C	304	-	52,52,52	0.96	2 (3%)	55,57,57	0.94	2 (3%)
25	DMU	P	309	-	34,34,34	0.59	1 (2%)	45,45,45	0.83	1 (2%)
14	HEA	N	601	1	44,67,67	1.00	2 (4%)	37,103,103	1.68	12 (32%)
20	EDO	F	104	-	3,3,3	0.48	0	2,2,2	0.48	0
20	EDO	N	617	-	3,3,3	0.41	0	2,2,2	0.42	0
14	HEA	N	602	1	44,67,67	0.92	1 (2%)	37,103,103	1.53	7 (18%)
19	PGV	N	607	-	50,50,50	0.89	2 (4%)	53,56,56	0.86	2 (3%)
25	DMU	C	312	-	34,34,34	0.57	1 (2%)	45,45,45	1.25	5 (11%)
18	TGL	L	101	-	62,62,62	1.26	6 (9%)	65,65,65	1.06	4 (6%)
25	DMU	C	313	-	34,34,34	0.60	1 (2%)	45,45,45	1.03	3 (6%)
19	PGV	P	303	-	50,50,50	0.90	2 (4%)	53,56,56	0.83	2 (3%)
20	EDO	B	303	-	3,3,3	0.41	0	2,2,2	0.39	0
21	CUA	B	301	2	0,1,1	0.00	-	-	-	-
20	EDO	A	614	-	3,3,3	0.36	0	2,2,2	0.40	0
25	DMU	M	102	-	34,34,34	0.47	0	45,45,45	0.70	0
20	EDO	F	102	-	3,3,3	0.41	0	2,2,2	0.43	0
20	EDO	A	610	-	3,3,3	0.47	0	2,2,2	0.50	0
20	EDO	N	610	-	3,3,3	0.48	0	2,2,2	0.37	0
20	EDO	N	615	-	3,3,3	0.39	0	2,2,2	0.41	0
23	PEK	T	102	-	52,52,52	0.89	3 (5%)	55,57,57	0.86	2 (3%)
24	CDL	G	101	-	99,99,99	1.32	12 (12%)	105,111,111	1.07	5 (4%)
19	PGV	N	606	-	50,50,50	0.97	2 (4%)	53,56,56	0.97	2 (3%)
22	CHD	C	301	-	29,32,32	0.59	0	48,51,51	0.98	2 (4%)
20	EDO	R	201	-	3,3,3	0.52	0	2,2,2	0.41	0
25	DMU	P	308	-	34,34,34	0.59	1 (2%)	45,45,45	1.01	2 (4%)
18	TGL	D	201	-	62,62,62	1.26	6 (9%)	65,65,65	0.96	3 (4%)
20	EDO	A	611	-	3,3,3	0.43	0	2,2,2	0.29	0
20	EDO	A	612	-	3,3,3	0.52	0	2,2,2	0.41	0
20	EDO	C	309	-	3,3,3	0.46	0	2,2,2	0.32	0
20	EDO	F	103	-	3,3,3	0.61	0	2,2,2	0.42	0
20	EDO	A	613	-	3,3,3	0.51	0	2,2,2	0.46	0
24	CDL	T	104	-	99,99,99	1.32	12 (12%)	105,111,111	1.10	5 (4%)
22	CHD	O	301	-	29,32,32	0.58	0	48,51,51	1.17	4 (8%)
20	EDO	N	609	-	3,3,3	0.40	0	2,2,2	0.43	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
22	CHD	B	302	-	29,32,32	0.62	0	48,51,51	1.09	2 (4%)
14	HEA	A	601	1	44,67,67	1.00	2 (4%)	37,103,103	1.89	14 (37%)
20	EDO	C	311	-	3,3,3	0.58	0	2,2,2	0.23	0
22	CHD	P	301	-	29,32,32	0.55	0	48,51,51	0.99	3 (6%)
20	EDO	D	203	-	3,3,3	0.50	0	2,2,2	0.39	0
19	PGV	P	304	-	50,50,50	0.97	2 (4%)	53,56,56	0.93	2 (3%)
25	DMU	Z	101	-	34,34,34	0.46	0	45,45,45	0.71	2 (4%)
20	EDO	N	613	-	3,3,3	0.35	0	2,2,2	0.61	0
22	CHD	P	306	-	29,32,32	0.49	0	48,51,51	1.72	13 (27%)
20	EDO	N	608	-	3,3,3	0.51	0	2,2,2	0.25	0
19	PGV	M	101	-	50,50,50	0.98	2 (4%)	53,56,56	1.00	2 (3%)
19	PGV	C	306	-	50,50,50	0.96	2 (4%)	53,56,56	0.96	2 (3%)
26	PSC	O	304	-	51,51,51	1.11	3 (5%)	57,59,59	1.01	2 (3%)
23	PEK	G	102	-	52,52,52	0.94	2 (3%)	55,57,57	0.95	2 (3%)
20	EDO	N	612	-	3,3,3	0.58	0	2,2,2	0.18	0
18	TGL	A	606	-	62,62,62	1.25	6 (9%)	65,65,65	1.41	8 (12%)
14	HEA	A	602	1	44,67,67	0.91	1 (2%)	37,103,103	1.31	8 (21%)
20	EDO	A	616	-	3,3,3	0.43	0	2,2,2	0.18	0
20	EDO	S	103	-	3,3,3	0.48	0	2,2,2	0.13	0
20	EDO	N	611	-	3,3,3	0.42	0	2,2,2	0.25	0
22	CHD	W	101	-	29,32,32	0.47	0	48,51,51	1.53	9 (18%)
18	TGL	Q	201	-	62,62,62	1.25	6 (9%)	65,65,65	1.11	6 (9%)
20	EDO	N	614	-	3,3,3	0.48	0	2,2,2	0.12	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
20	EDO	A	615	-	-	0/1/1/1	-
18	TGL	Y	101	-	-	36/65/65/65	-
26	PSC	E	201	-	-	32/55/55/55	-
23	PEK	T	103	-	-	29/56/56/56	-
18	TGL	O	302	-	-	34/65/65/65	-
20	EDO	A	609	-	-	0/1/1/1	-
20	EDO	S	102	-	-	0/1/1/1	-
20	EDO	A	608	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
22	CHD	J	101	-	-	3/7/74/74	0/4/4/4
24	CDL	C	307	-	-	59/110/110/110	-
19	PGV	A	607	-	-	11/55/55/55	-
22	CHD	C	308	-	-	7/7/74/74	0/4/4/4
19	PGV	C	305	-	-	16/55/55/55	-
20	EDO	N	616	-	-	0/1/1/1	-
20	EDO	D	202	-	-	1/1/1/1	-
20	EDO	E	202	-	-	0/1/1/1	-
23	PEK	T	101	-	-	34/56/56/56	-
23	PEK	C	303	-	-	17/56/56/56	-
24	CDL	P	305	-	-	64/110/110/110	-
20	EDO	P	307	-	-	0/1/1/1	-
20	EDO	C	310	-	-	0/1/1/1	-
23	PEK	C	304	-	-	31/56/56/56	-
25	DMU	P	309	-	-	7/19/59/59	0/2/2/2
14	HEA	N	601	1	3/3/7/16	0/24/76/76	-
20	EDO	F	104	-	-	1/1/1/1	-
20	EDO	N	617	-	-	0/1/1/1	-
14	HEA	N	602	1	2/2/7/16	3/24/76/76	-
19	PGV	N	607	-	-	11/55/55/55	-
25	DMU	C	312	-	-	6/19/59/59	0/2/2/2
18	TGL	L	101	-	-	36/65/65/65	-
25	DMU	C	313	-	-	8/19/59/59	0/2/2/2
19	PGV	P	303	-	-	16/55/55/55	-
20	EDO	B	303	-	-	1/1/1/1	-
20	EDO	A	614	-	-	1/1/1/1	-
25	DMU	M	102	-	-	4/19/59/59	0/2/2/2
20	EDO	F	102	-	-	0/1/1/1	-
20	EDO	A	610	-	-	0/1/1/1	-
20	EDO	N	610	-	-	0/1/1/1	-
20	EDO	N	615	-	-	1/1/1/1	-
23	PEK	T	102	-	-	20/56/56/56	-
24	CDL	G	101	-	-	63/110/110/110	-
19	PGV	N	606	-	-	31/55/55/55	-
22	CHD	C	301	-	-	0/7/74/74	0/4/4/4
20	EDO	R	201	-	-	0/1/1/1	-
25	DMU	P	308	-	-	9/19/59/59	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
18	TGL	D	201	-	-	39/65/65/65	-
20	EDO	A	611	-	-	0/1/1/1	-
20	EDO	A	612	-	-	0/1/1/1	-
20	EDO	C	309	-	-	0/1/1/1	-
20	EDO	F	103	-	-	0/1/1/1	-
20	EDO	A	613	-	-	0/1/1/1	-
24	CDL	T	104	-	-	61/110/110/110	-
22	CHD	O	301	-	-	0/7/74/74	0/4/4/4
20	EDO	N	609	-	-	1/1/1/1	-
22	CHD	B	302	-	-	0/7/74/74	0/4/4/4
14	HEA	A	601	1	3/3/7/16	1/24/76/76	-
20	EDO	N	612	-	-	0/1/1/1	-
20	EDO	C	311	-	-	0/1/1/1	-
22	CHD	P	301	-	-	0/7/74/74	0/4/4/4
20	EDO	D	203	-	-	0/1/1/1	-
19	PGV	P	304	-	-	32/55/55/55	-
25	DMU	Z	101	-	-	7/19/59/59	0/2/2/2
20	EDO	N	613	-	-	1/1/1/1	-
22	CHD	P	306	-	-	6/7/74/74	0/4/4/4
20	EDO	N	608	-	-	1/1/1/1	-
19	PGV	M	101	-	-	31/55/55/55	-
19	PGV	C	306	-	-	25/55/55/55	-
26	PSC	O	304	-	-	34/55/55/55	-
23	PEK	G	102	-	-	34/56/56/56	-
14	HEA	A	602	1	2/2/7/16	3/24/76/76	-
18	TGL	A	606	-	-	37/65/65/65	-
20	EDO	A	616	-	-	1/1/1/1	-
20	EDO	S	103	-	-	0/1/1/1	-
20	EDO	N	611	-	-	0/1/1/1	-
22	CHD	W	101	-	-	4/7/74/74	0/4/4/4
18	TGL	Q	201	-	-	35/65/65/65	-
20	EDO	N	614	-	-	0/1/1/1	-

The worst 5 of 129 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
18	D	201	TGL	OG1-CA1	4.73	1.47	1.33
18	A	606	TGL	OG2-CB1	4.66	1.47	1.34
24	C	307	CDL	OA8-CA7	4.63	1.46	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
24	C	307	CDL	OB8-CB7	4.60	1.46	1.33
18	L	101	TGL	OG2-CB1	4.60	1.47	1.34

The worst 5 of 187 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
22	J	101	CHD	C17-C13-C14	-4.98	95.08	100.09
18	A	606	TGL	OG2-CB1-CB2	4.60	121.41	111.50
24	T	104	CDL	OB6-CB5-C51	4.42	121.02	111.50
24	C	307	CDL	OA6-CA5-C11	4.26	120.69	111.50
24	G	101	CDL	OB6-CB5-C51	4.24	120.64	111.50

5 of 10 chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
14	A	601	HEA	ND
14	A	601	HEA	NA
14	A	601	HEA	NB
14	N	601	HEA	ND
14	N	601	HEA	NA

5 of 945 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
18	O	302	TGL	OB1-CB1-OG2-CG2
23	T	101	PEK	O12-C04-C05-N
23	T	101	PEK	C2-C1-O01-C02
23	T	101	PEK	C7-C8-C9-C10
23	T	101	PEK	C12-C13-C14-C15

There are no ring outliers.

49 monomers are involved in 153 short contacts:

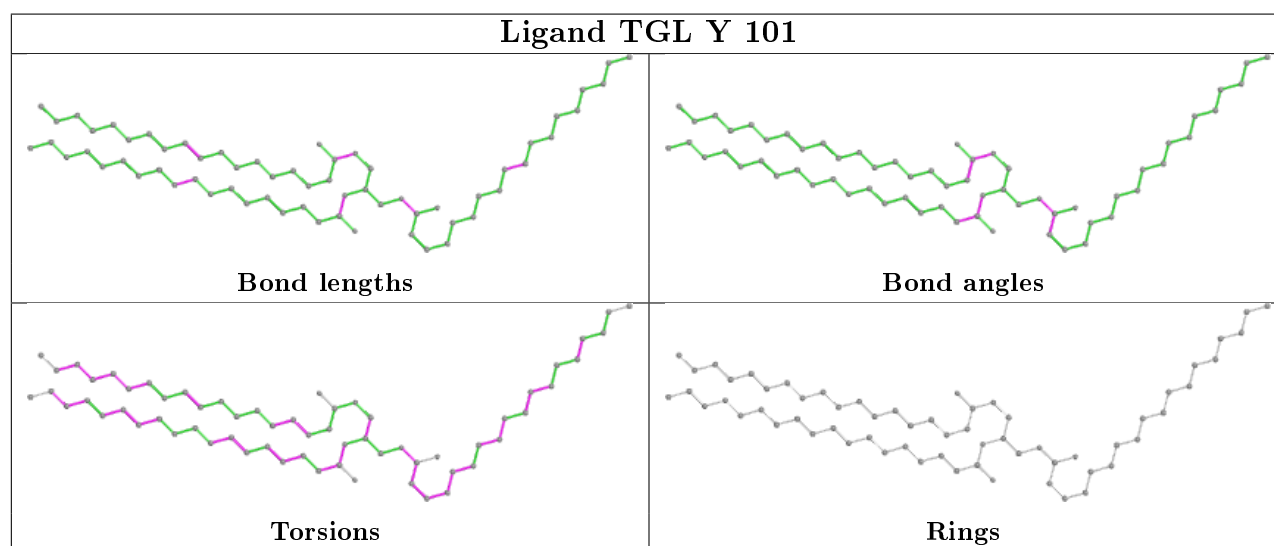
Mol	Chain	Res	Type	Clashes	Symm-Clashes
20	A	615	EDO	1	0
18	Y	101	TGL	5	0
26	E	201	PSC	10	0
23	T	103	PEK	2	0
18	O	302	TGL	2	0
20	A	608	EDO	1	0
24	C	307	CDL	10	0

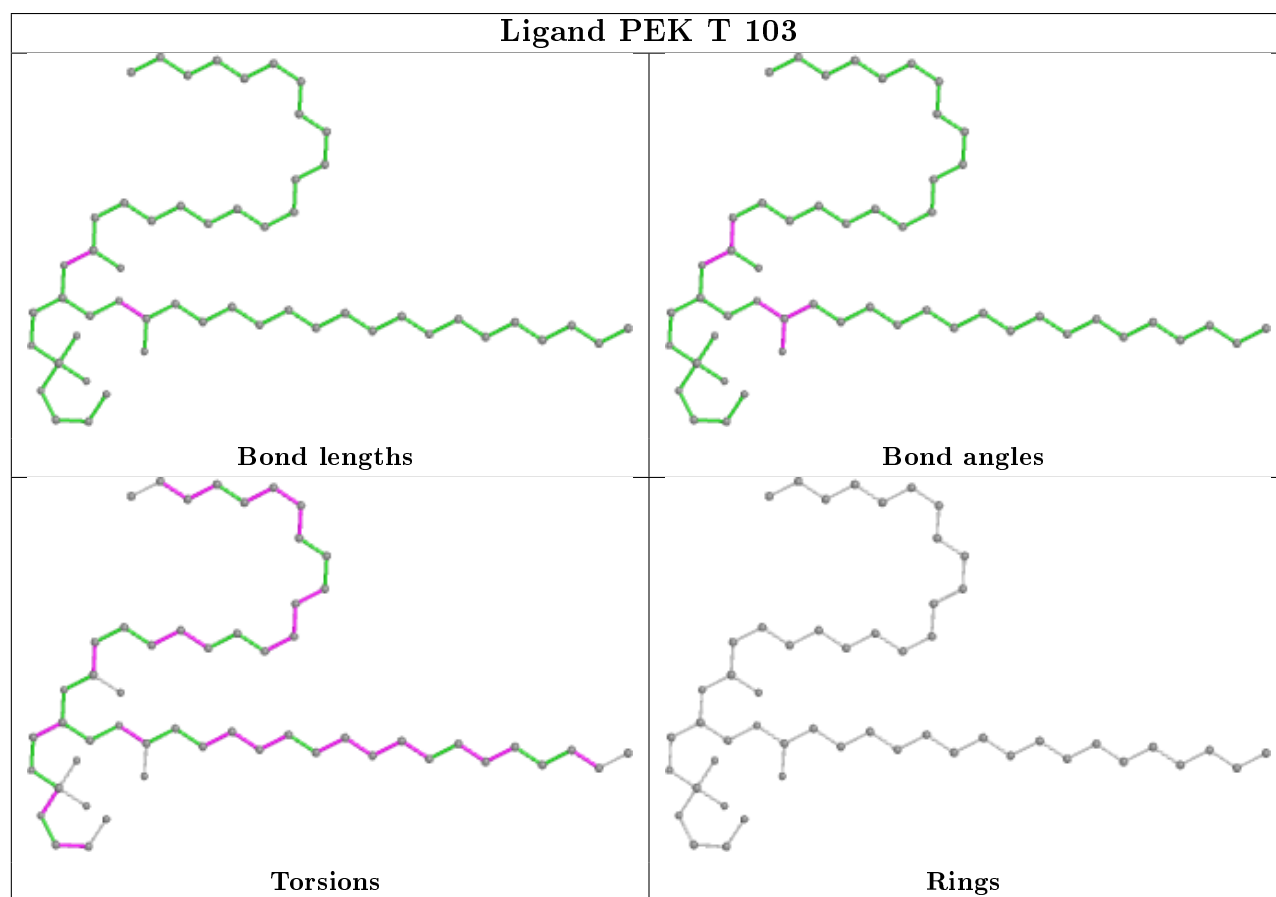
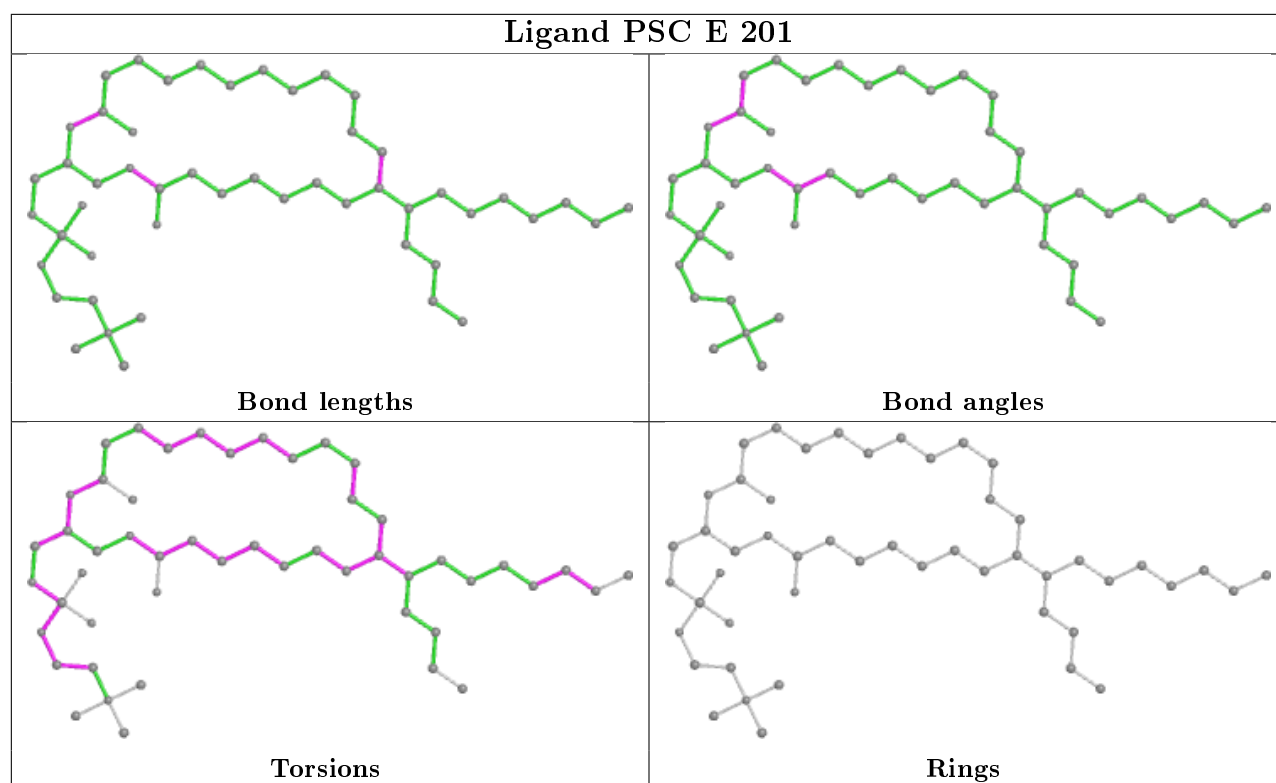
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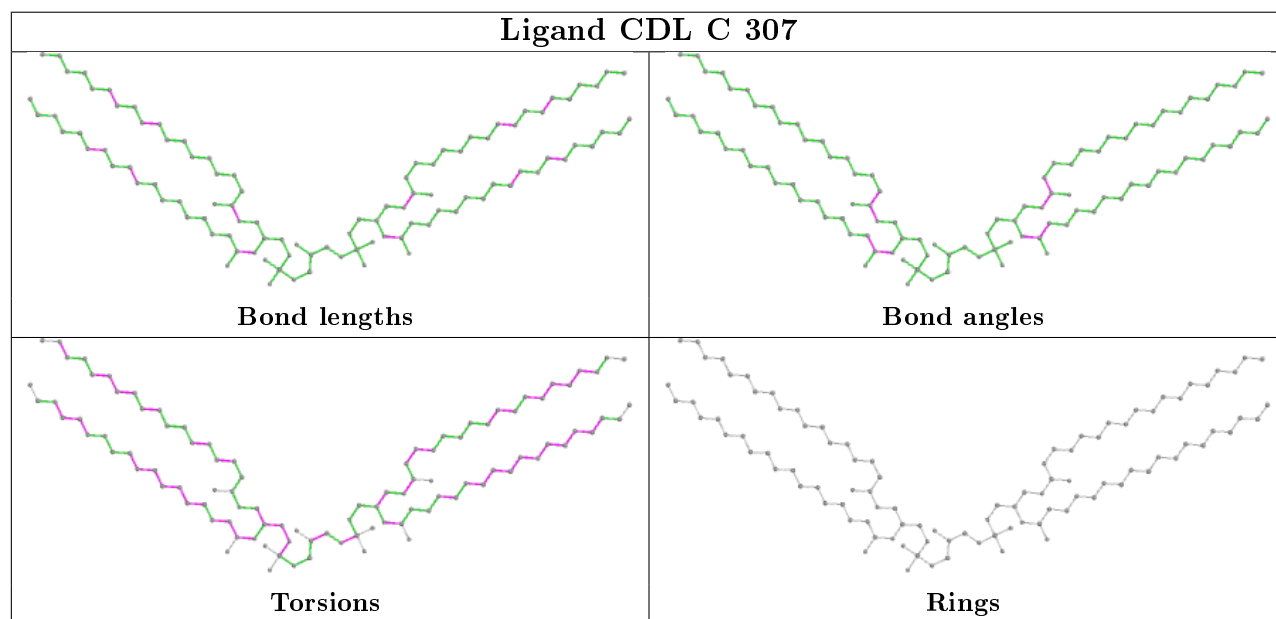
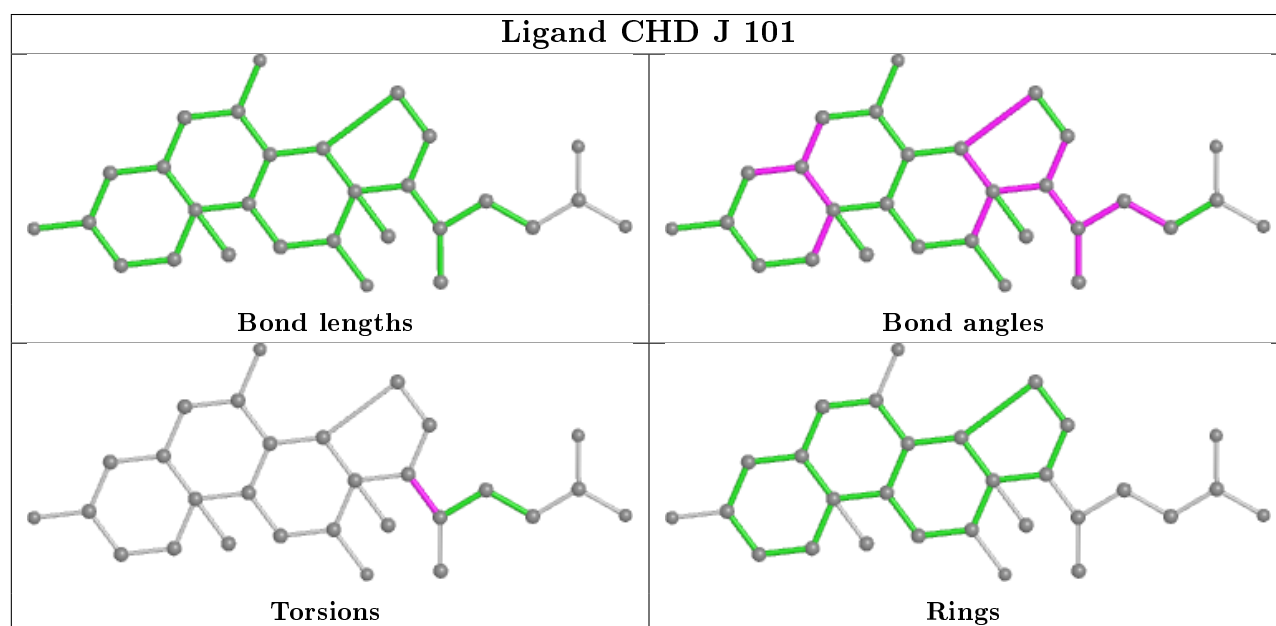
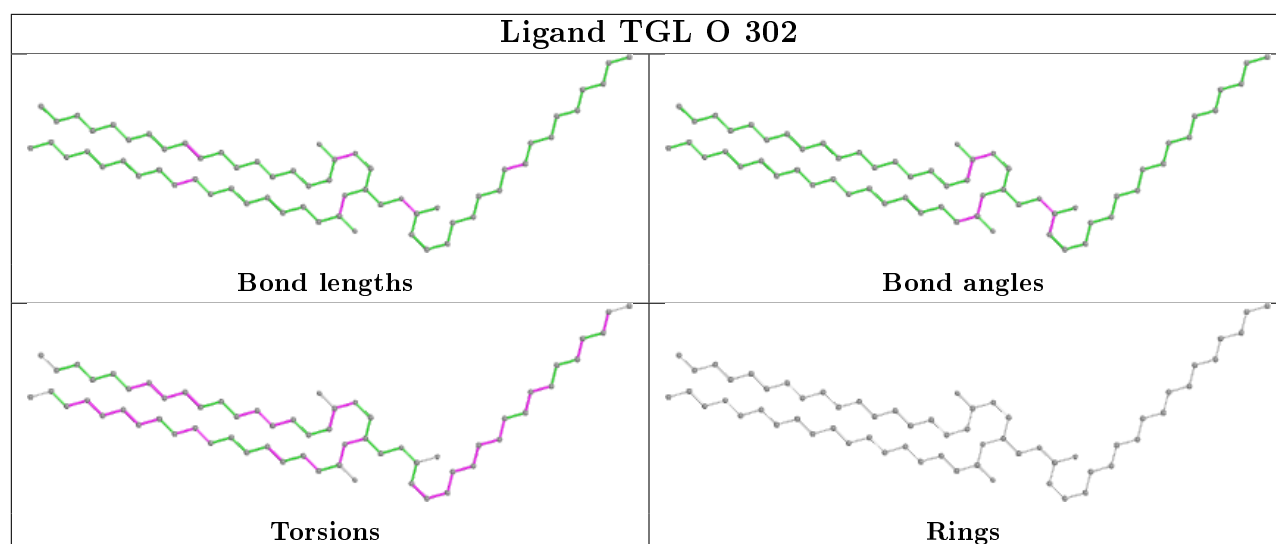
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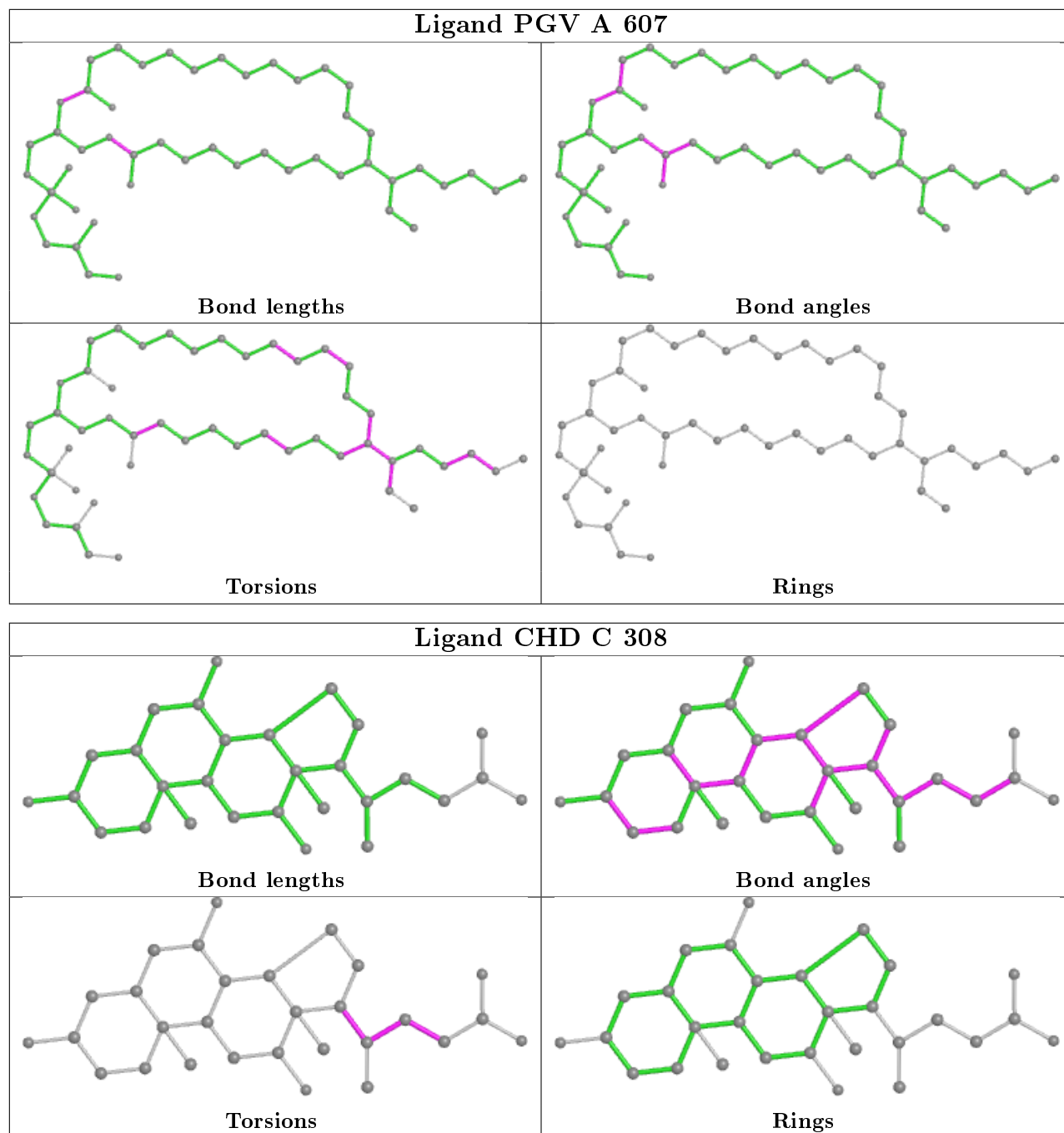
Mol	Chain	Res	Type	Clashes	Symm-Clashes
19	A	607	PGV	2	0
22	C	308	CHD	3	0
19	C	305	PGV	3	0
20	D	202	EDO	2	0
20	E	202	EDO	1	0
23	T	101	PEK	2	0
23	C	303	PEK	3	0
24	P	305	CDL	8	0
23	C	304	PEK	2	0
25	P	309	DMU	8	0
14	N	601	HEA	2	0
20	N	617	EDO	1	0
14	N	602	HEA	5	0
19	N	607	PGV	2	0
18	L	101	TGL	4	0
25	C	313	DMU	12	0
19	P	303	PGV	3	0
20	N	615	EDO	2	0
23	T	102	PEK	3	0
24	G	101	CDL	5	0
19	N	606	PGV	2	0
25	P	308	DMU	1	0
18	D	201	TGL	4	0
20	A	611	EDO	2	0
24	T	104	CDL	13	0
22	O	301	CHD	1	0
14	A	601	HEA	2	0
22	P	301	CHD	2	0
19	P	304	PGV	1	0
20	N	613	EDO	4	0
22	P	306	CHD	1	0
20	N	608	EDO	1	0
19	M	101	PGV	2	0
19	C	306	PGV	1	0
26	O	304	PSC	6	0
20	N	612	EDO	1	0
18	A	606	TGL	3	0
14	A	602	HEA	4	0
20	A	616	EDO	1	0
20	S	103	EDO	1	0
22	W	101	CHD	1	0
20	N	614	EDO	1	0

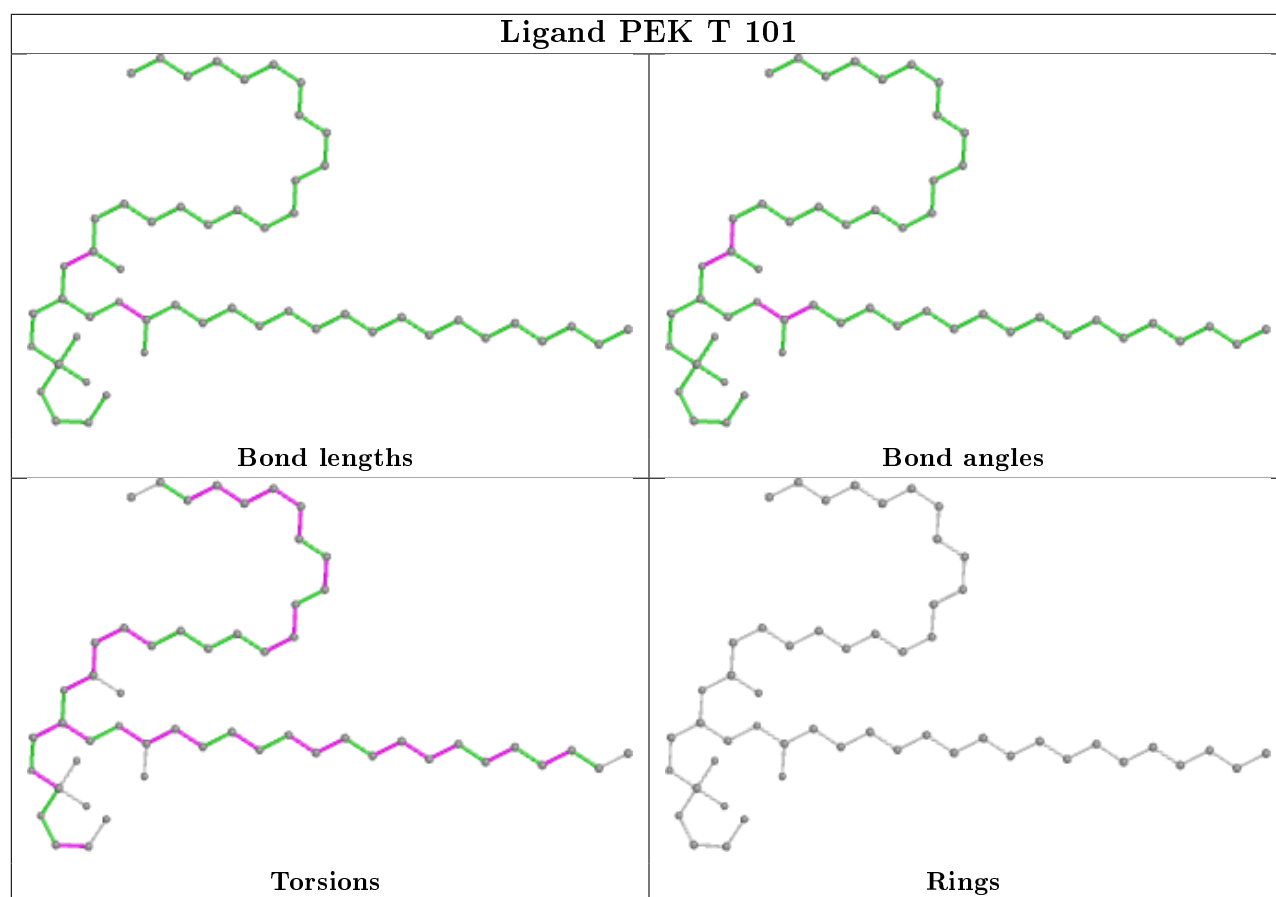
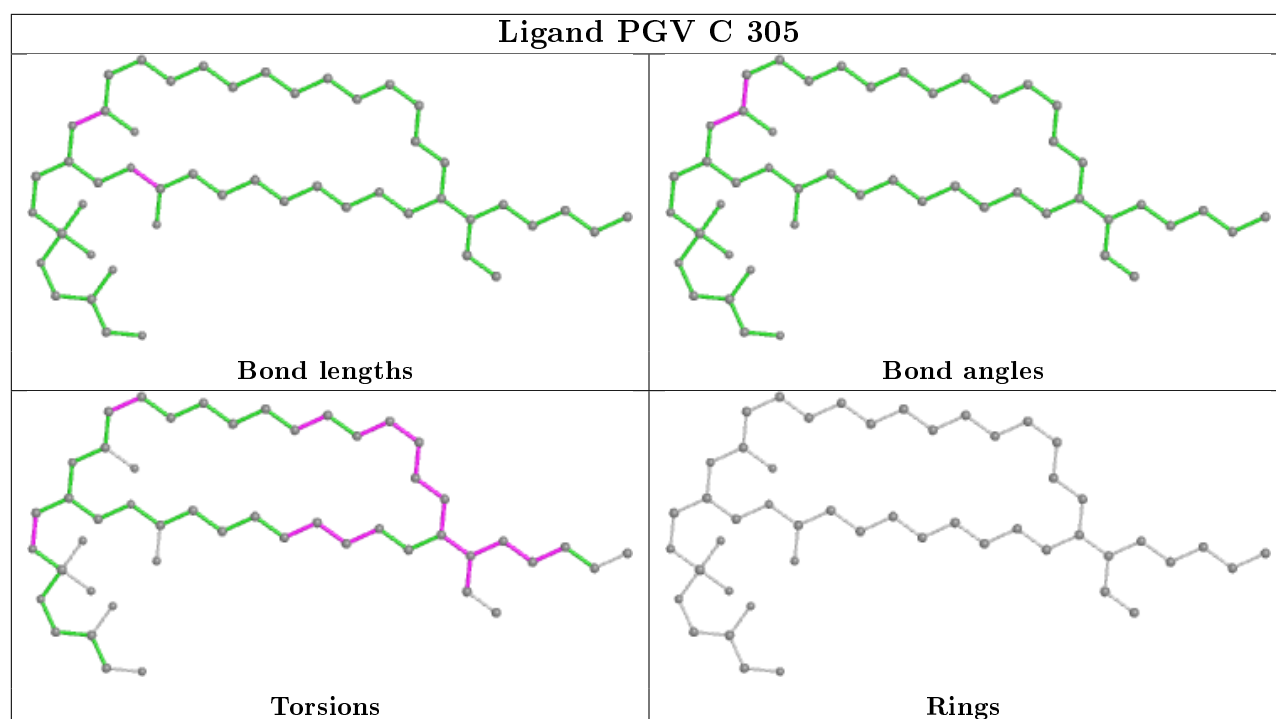
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

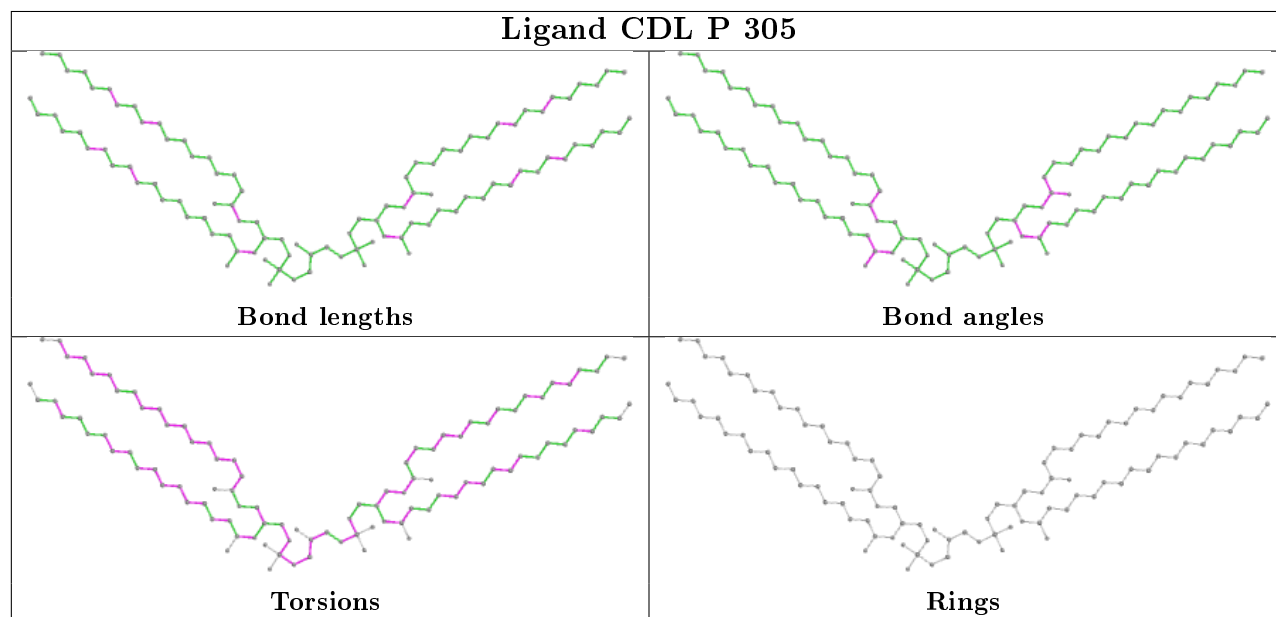
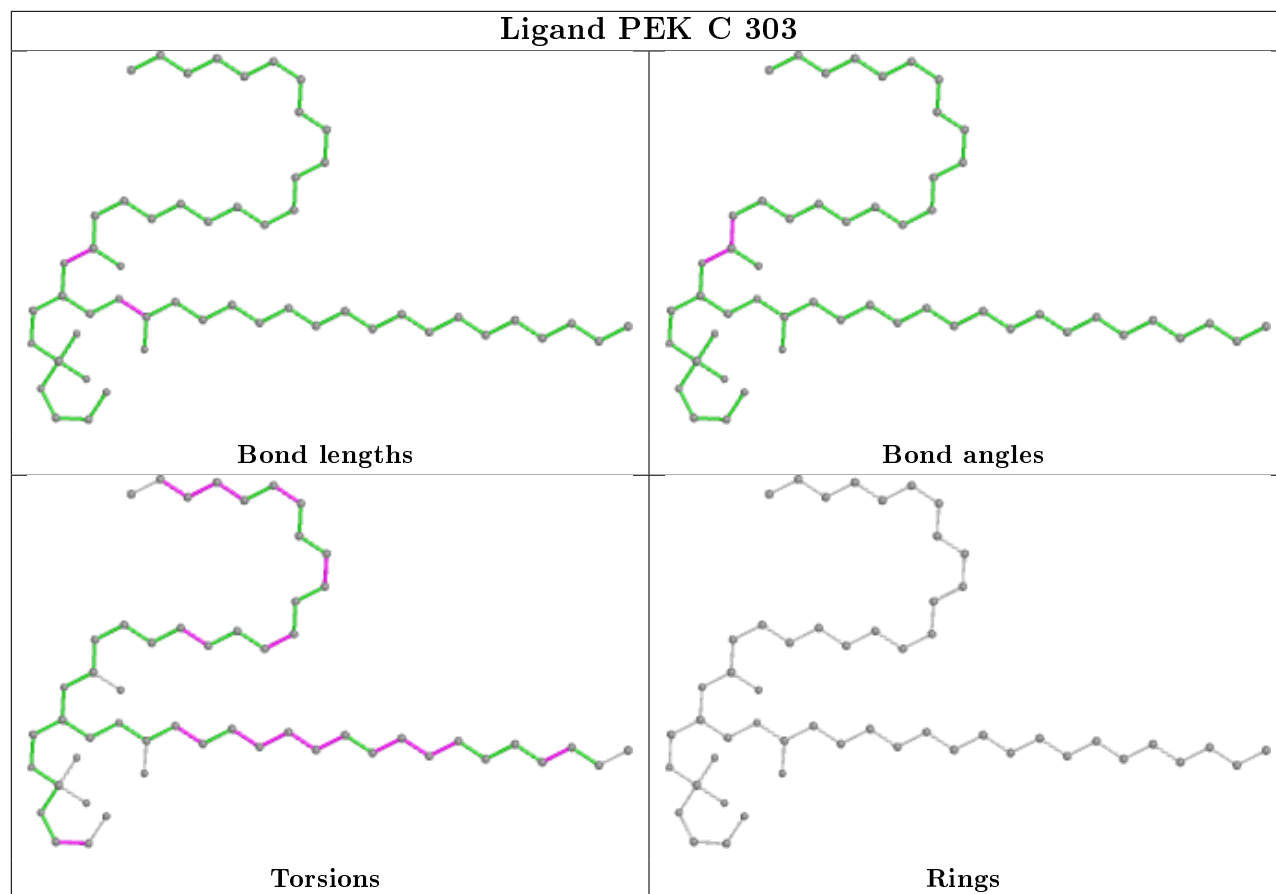


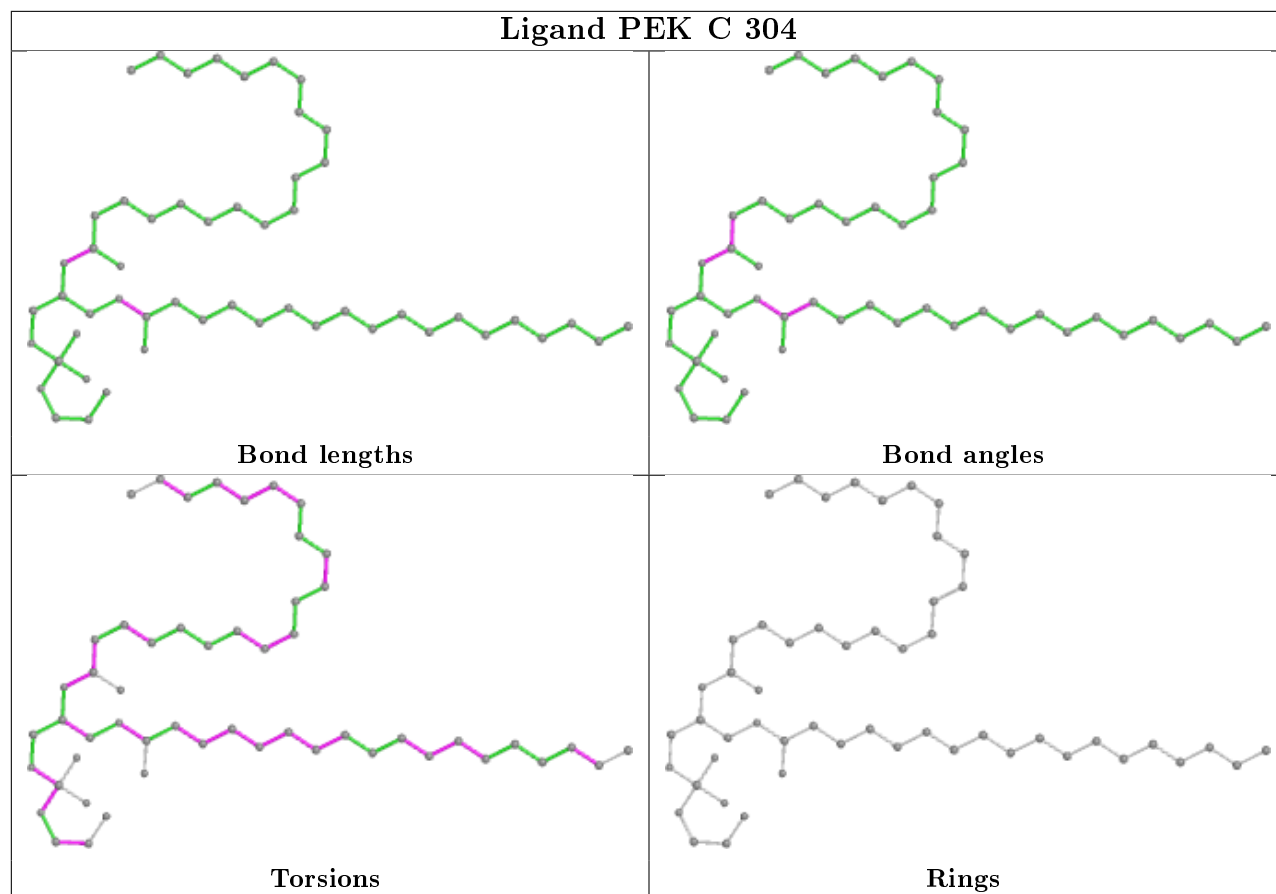


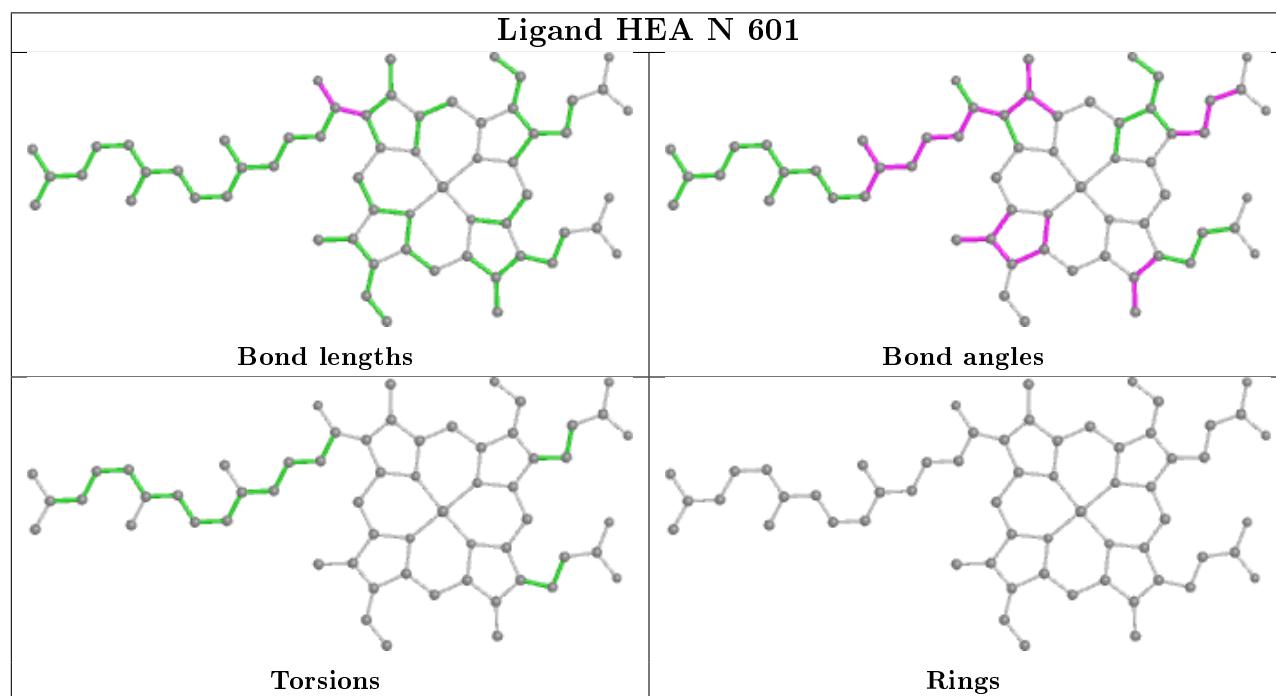
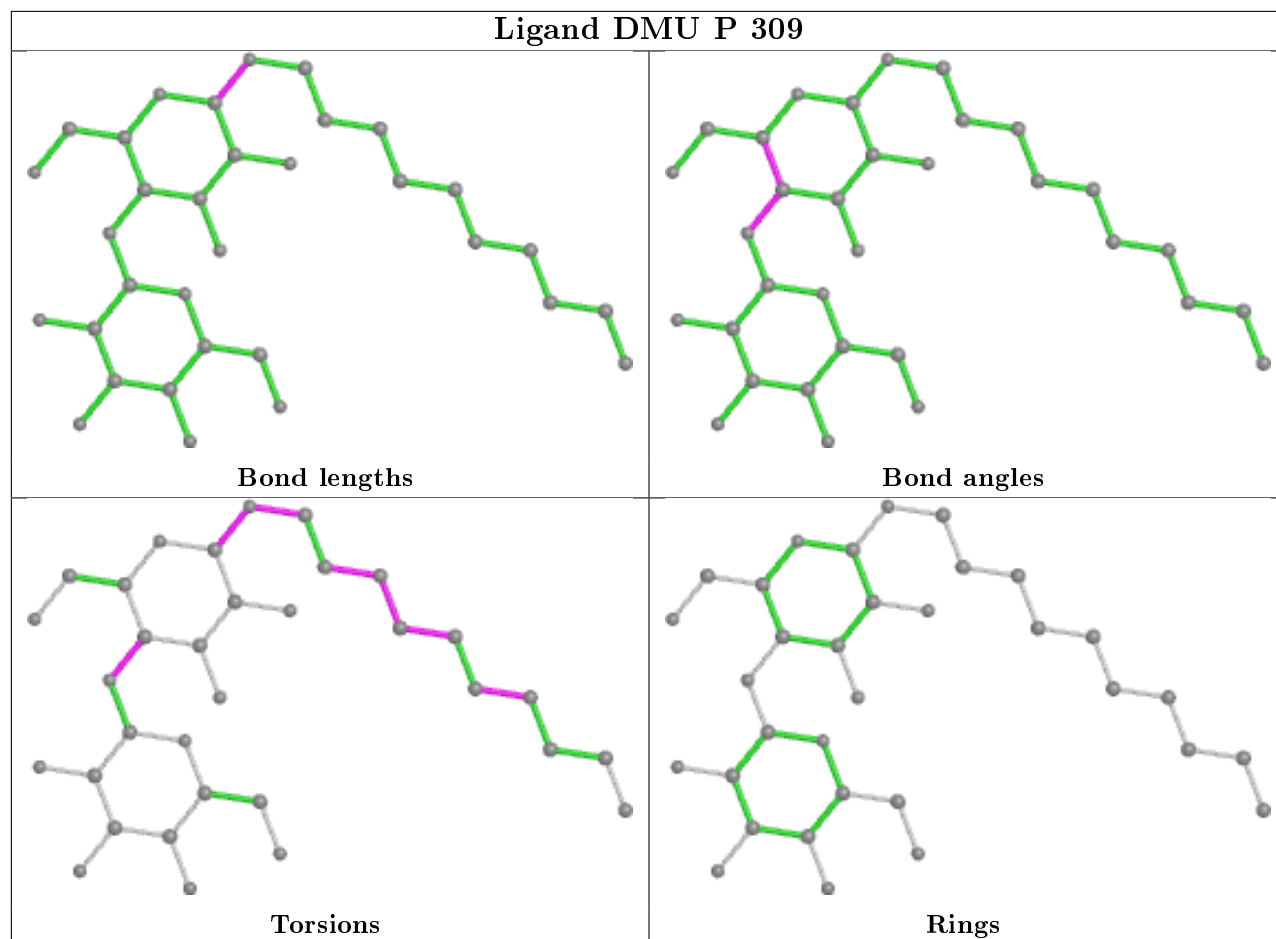


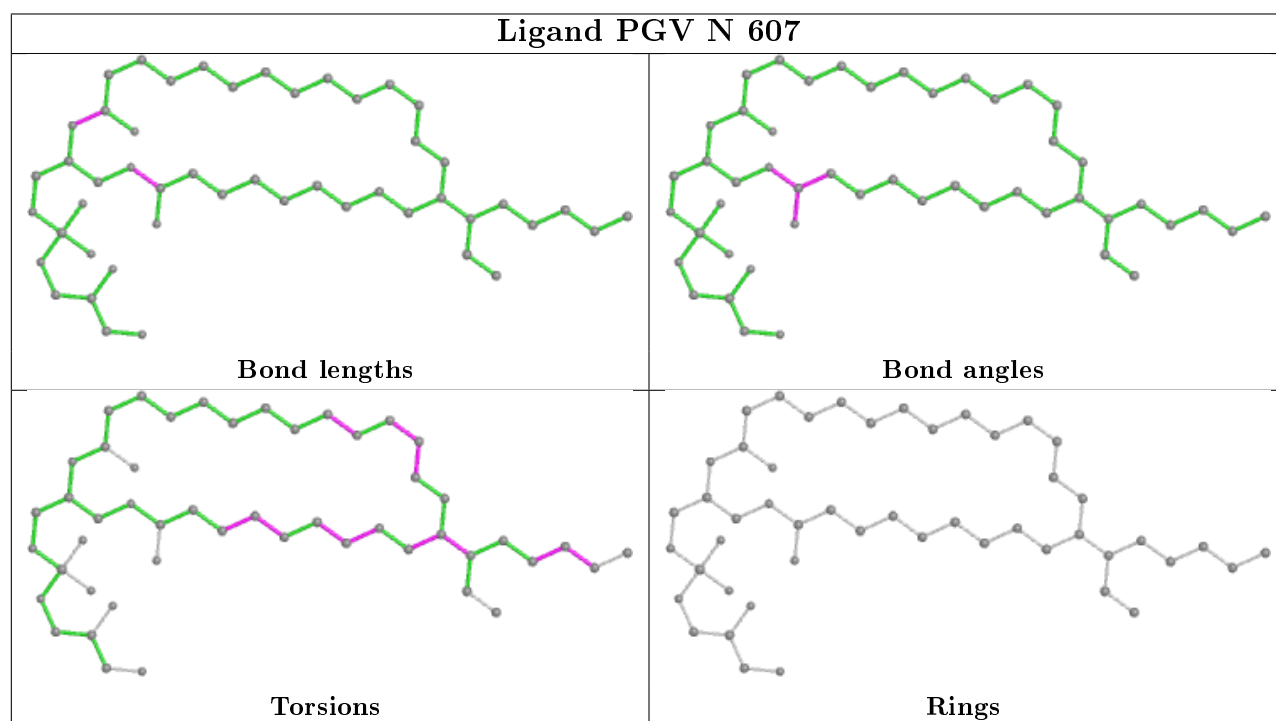
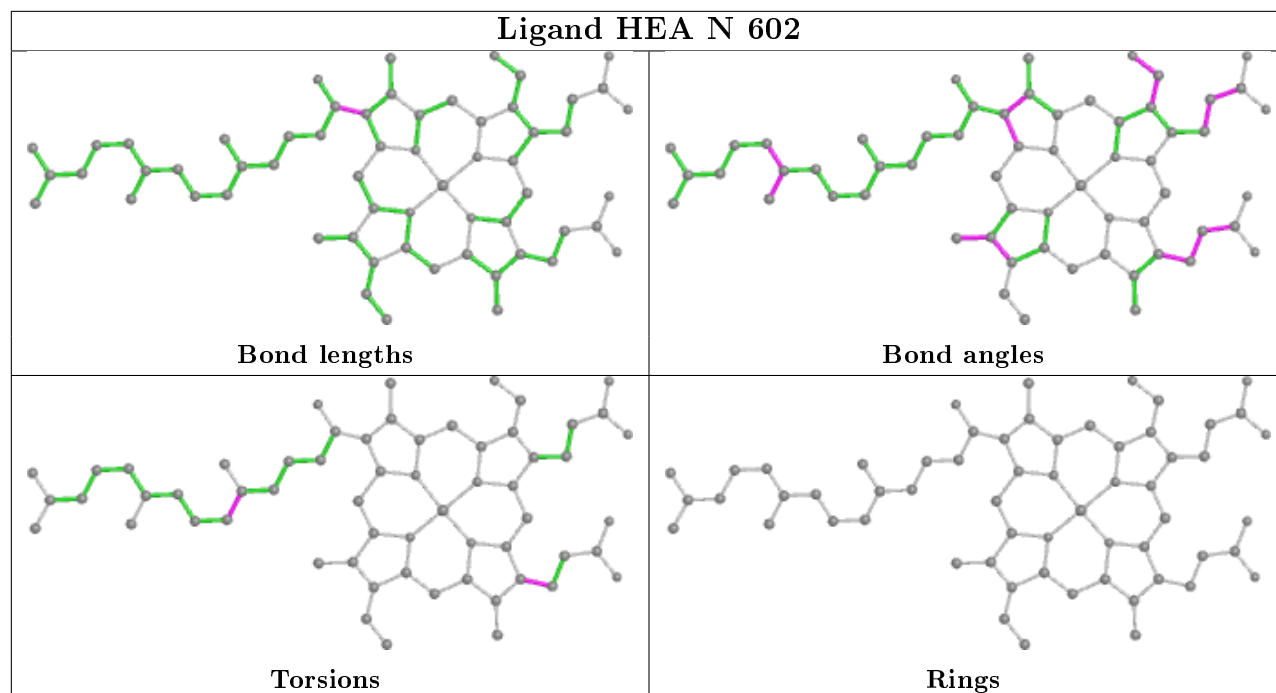


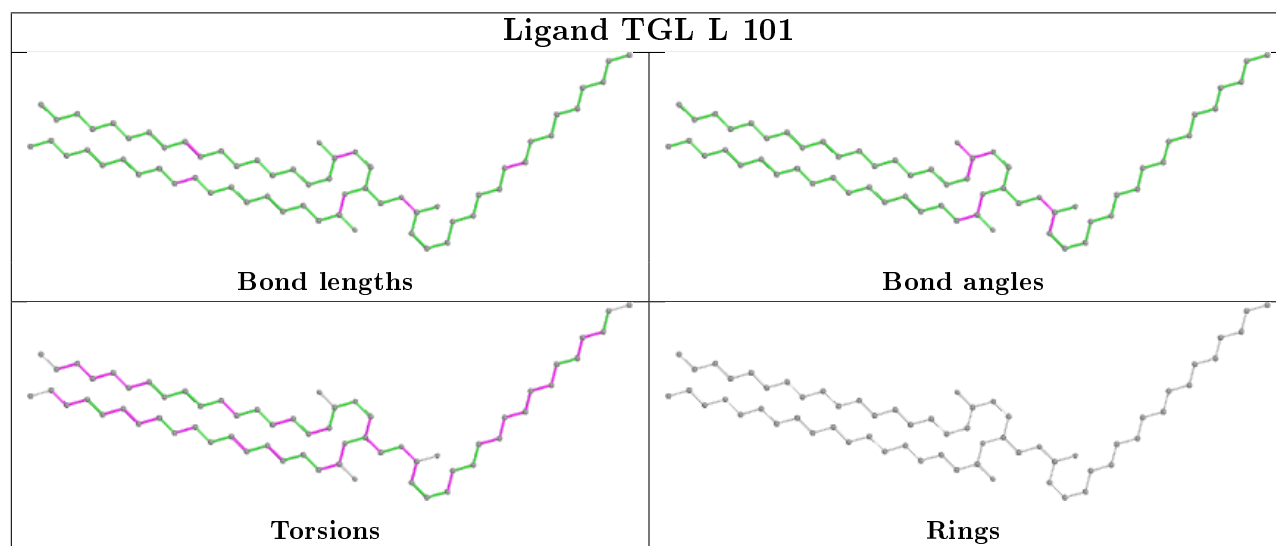
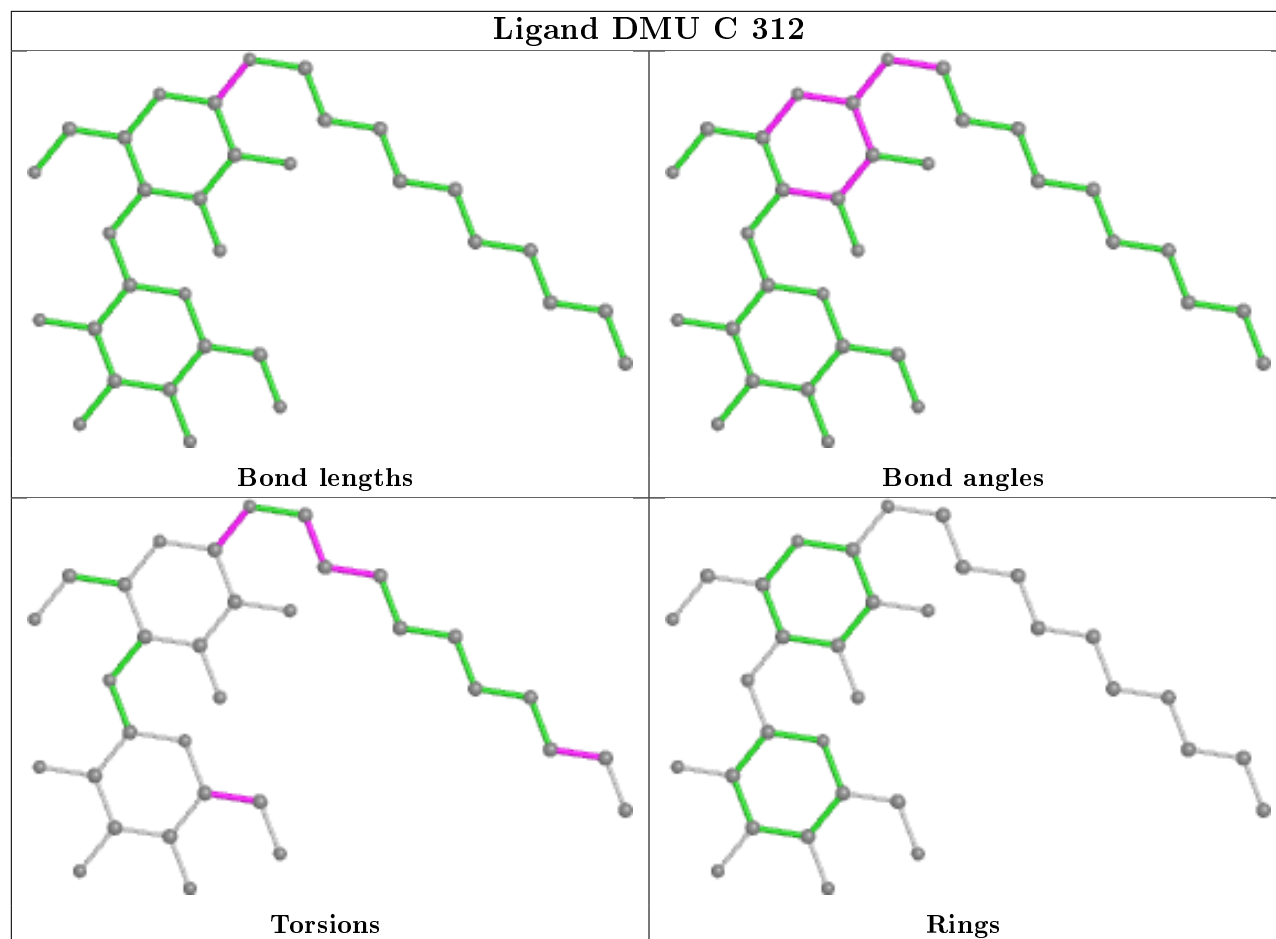


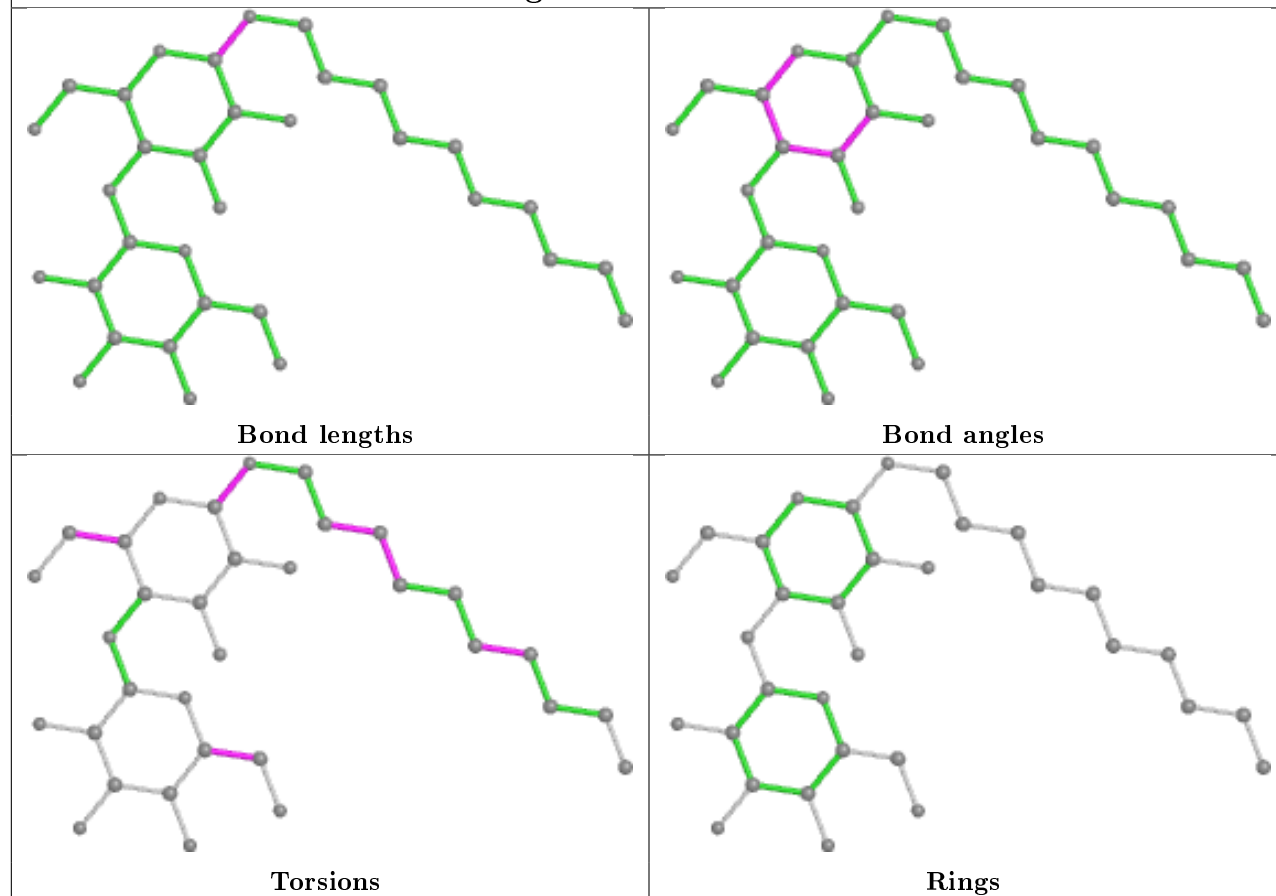
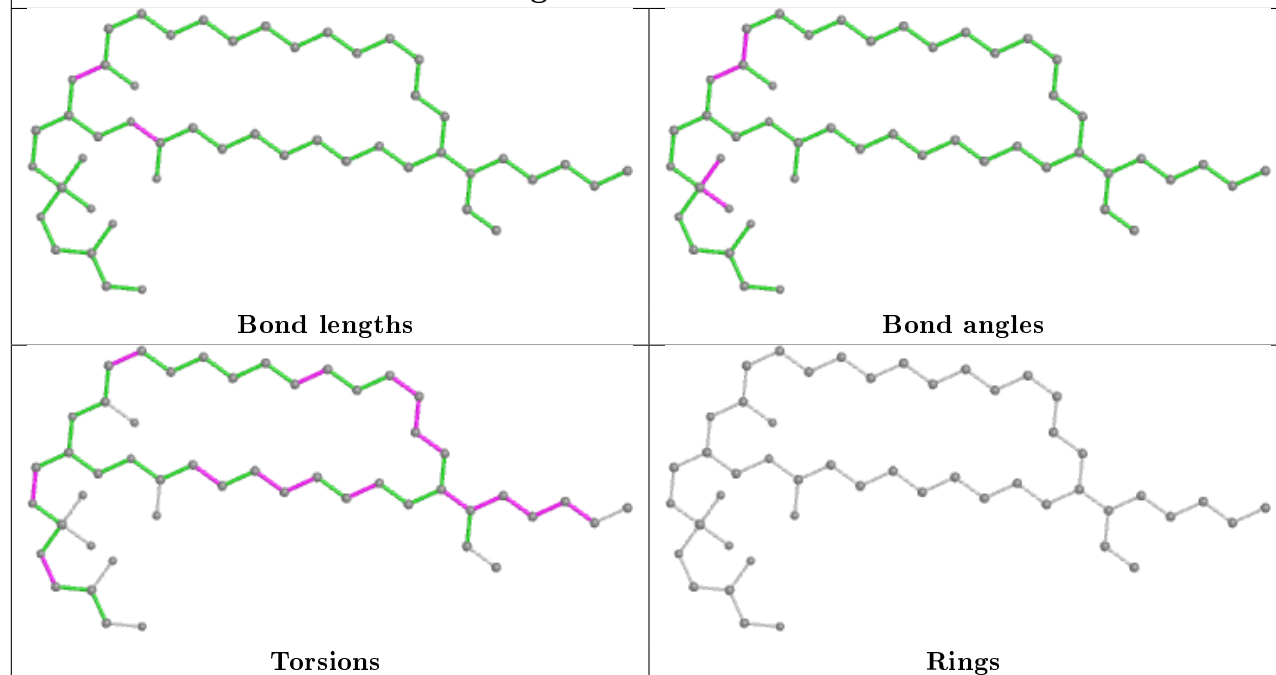


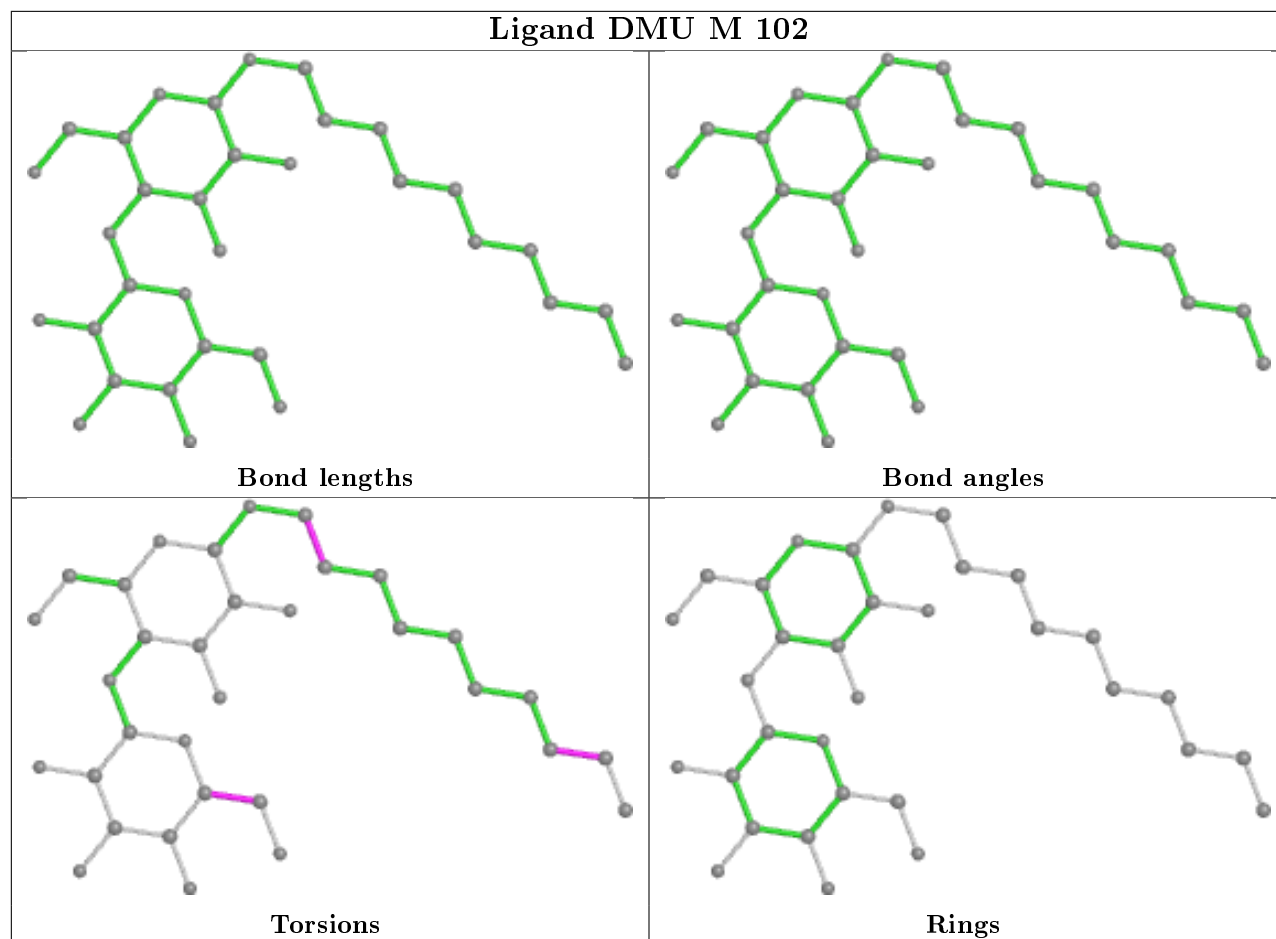


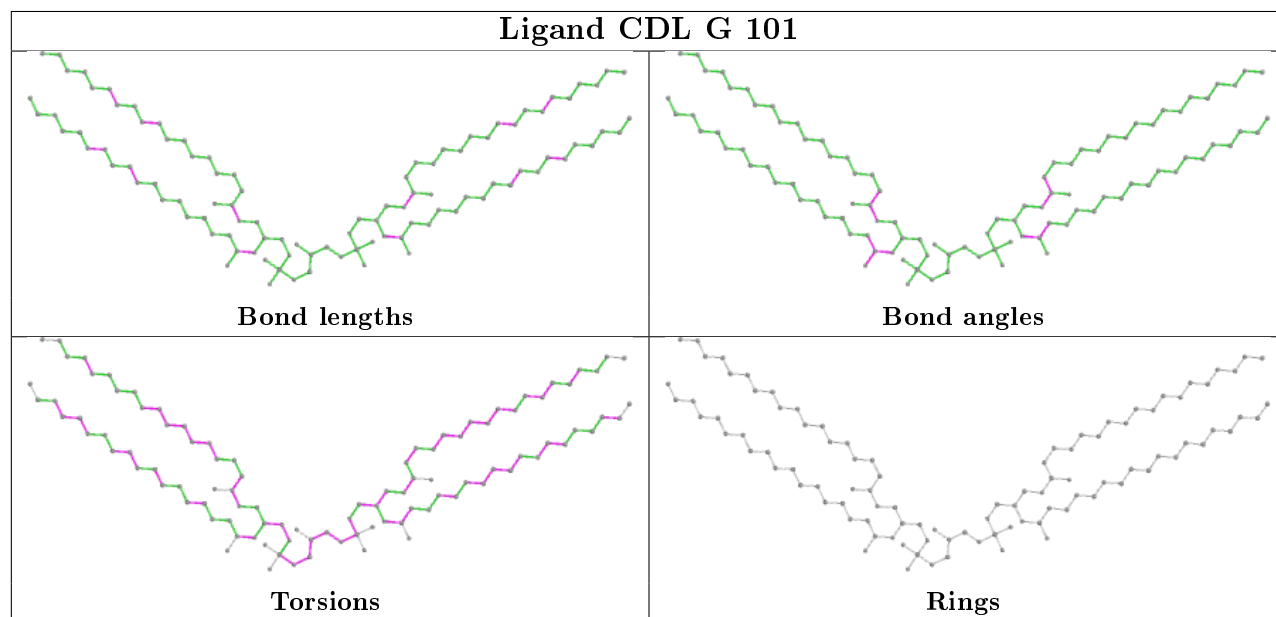
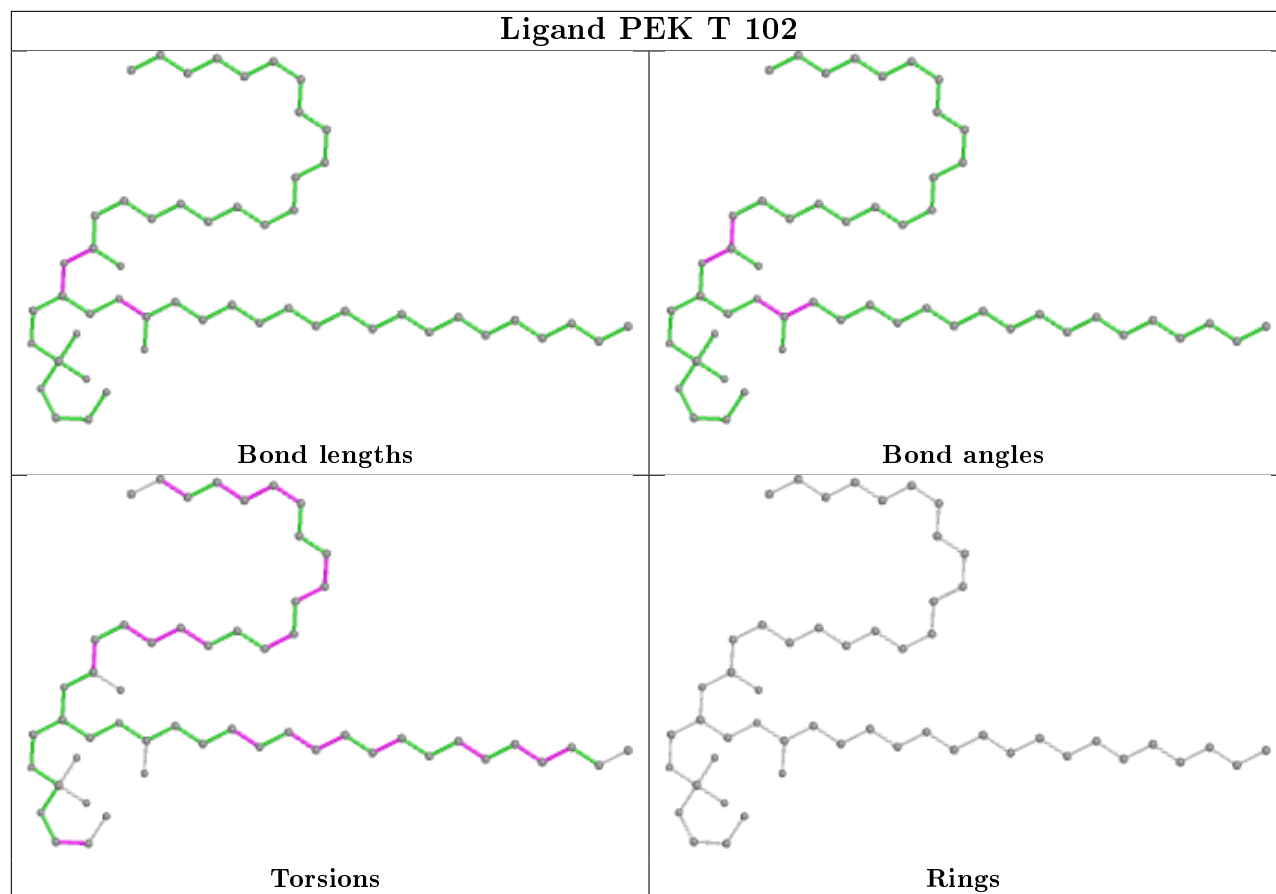


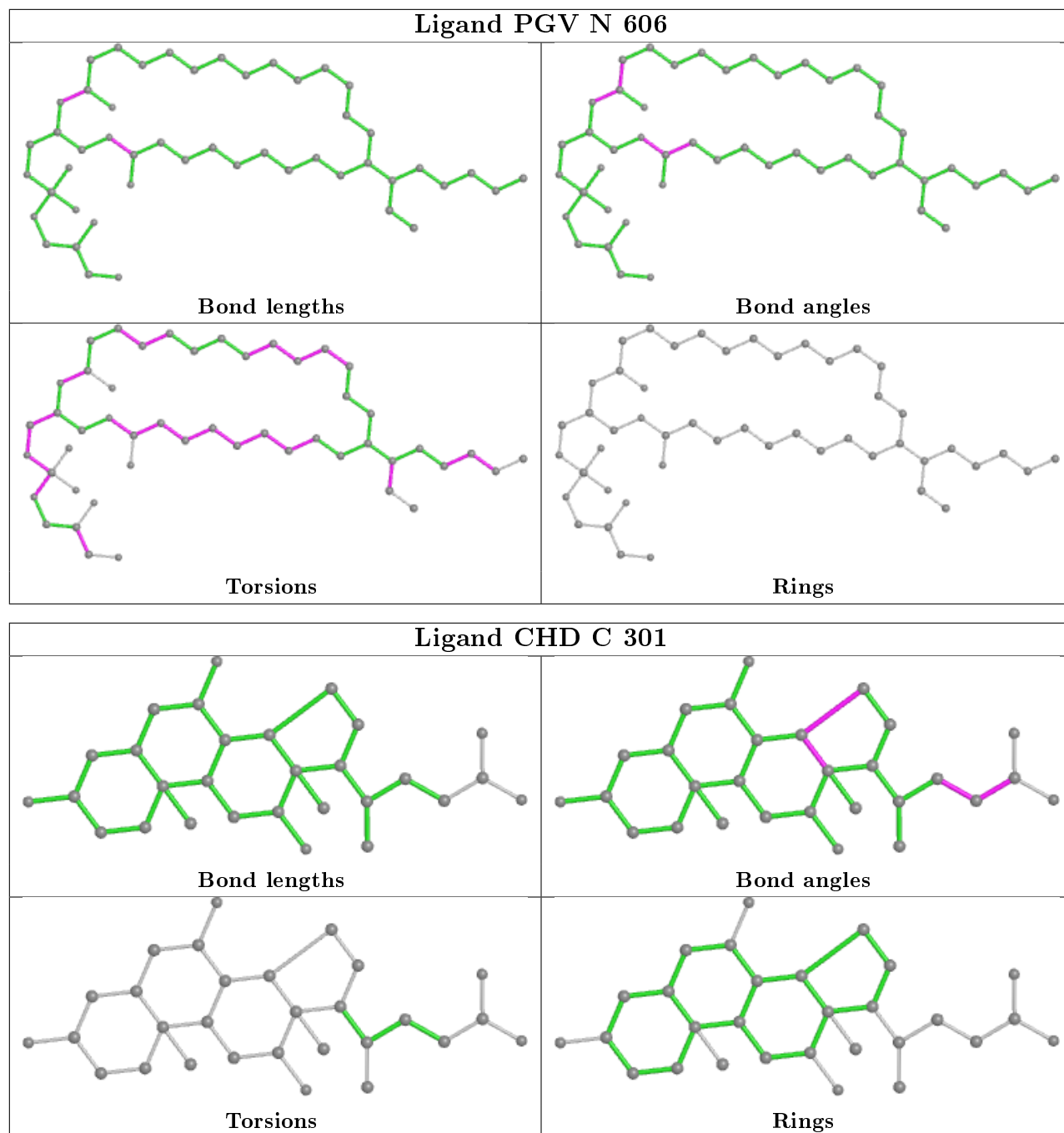


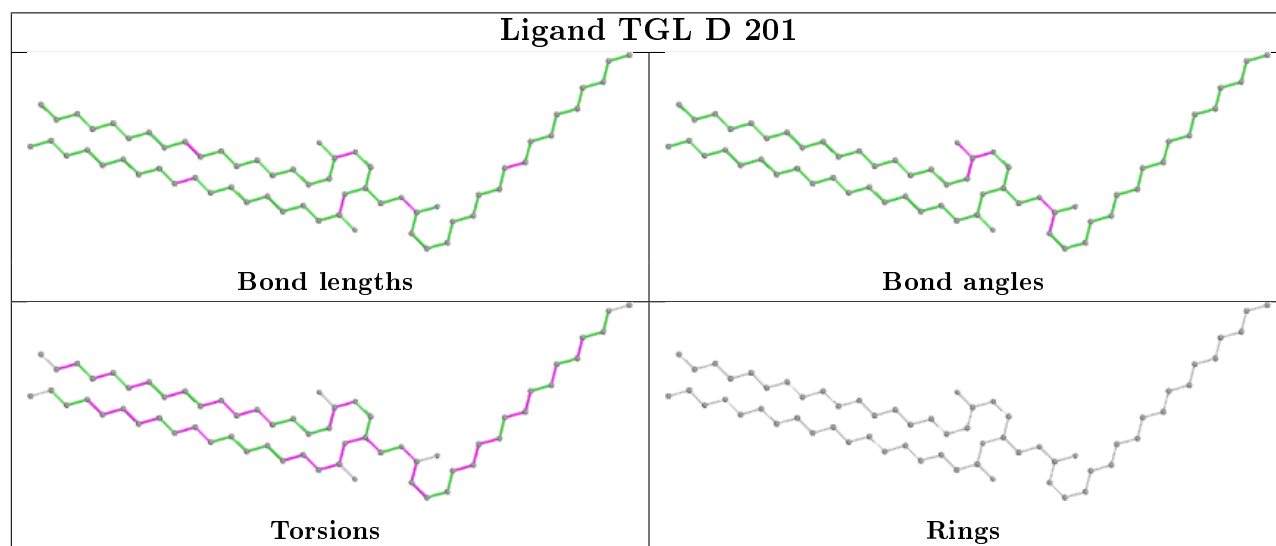
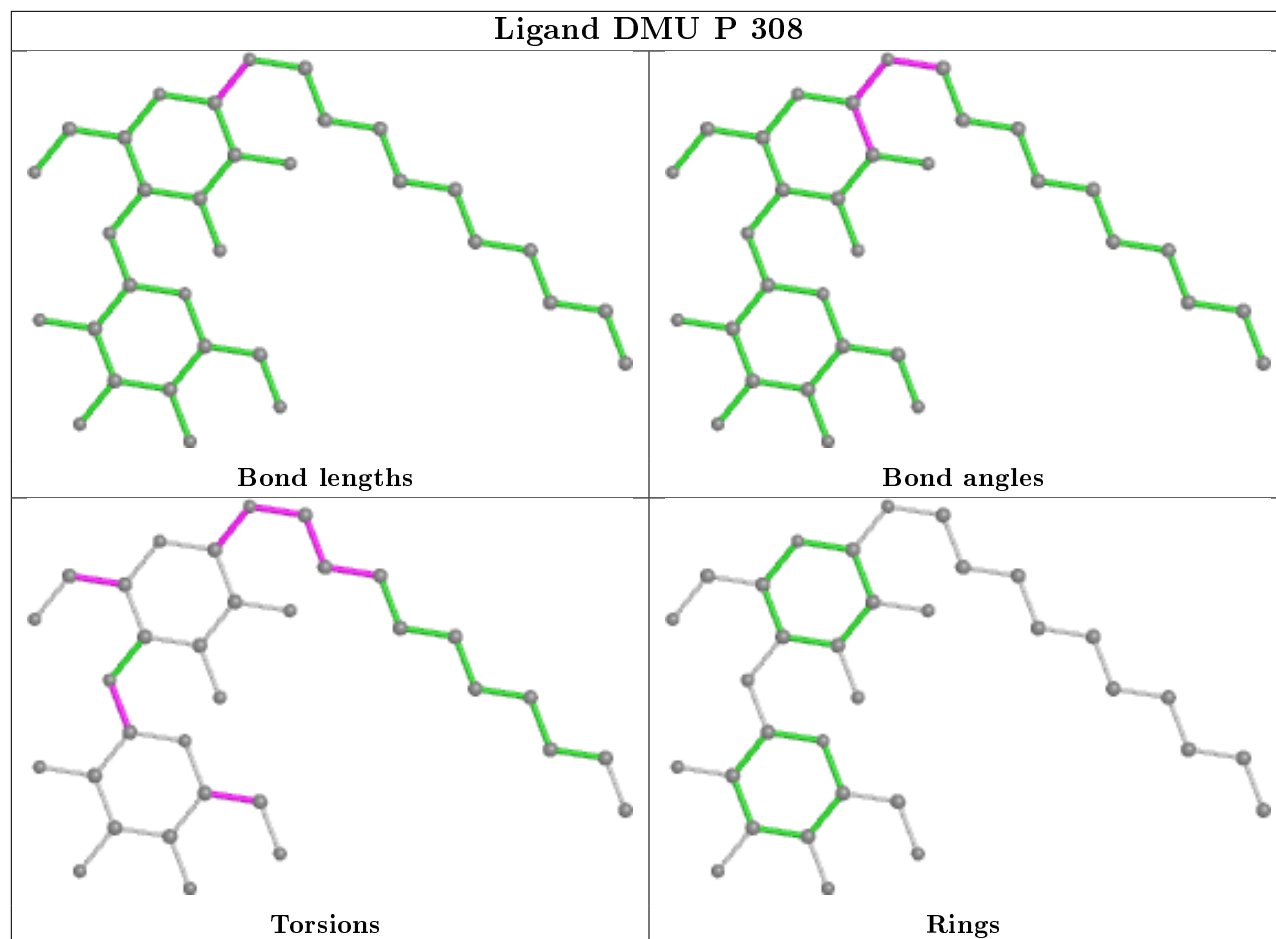


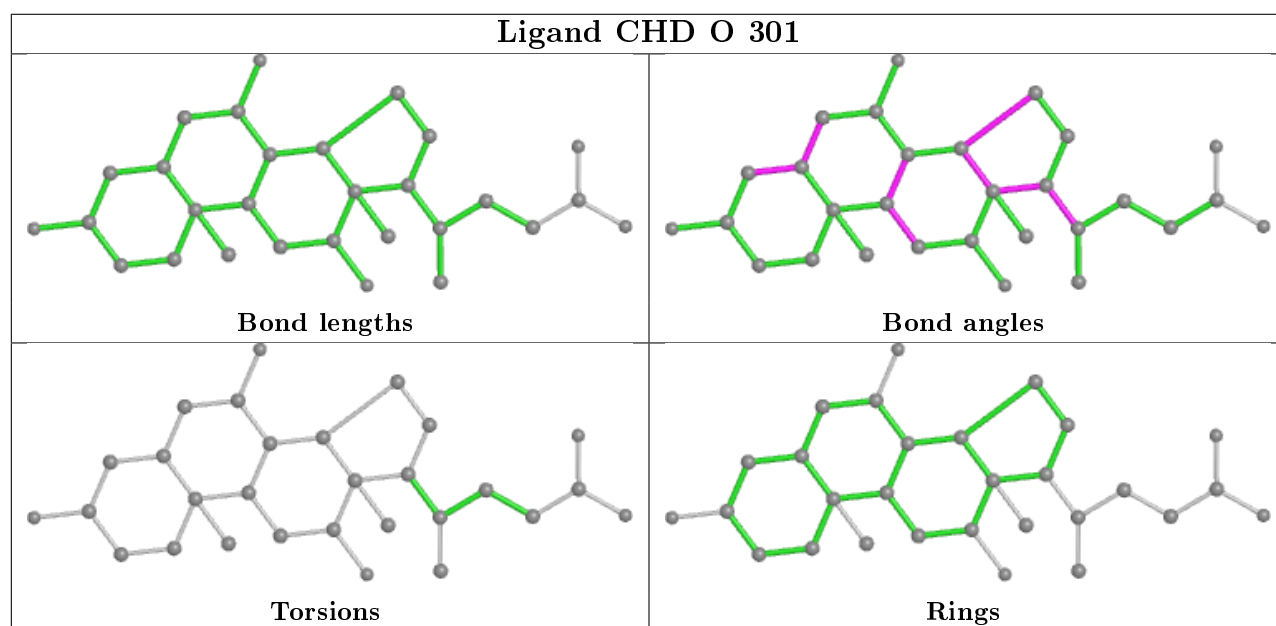
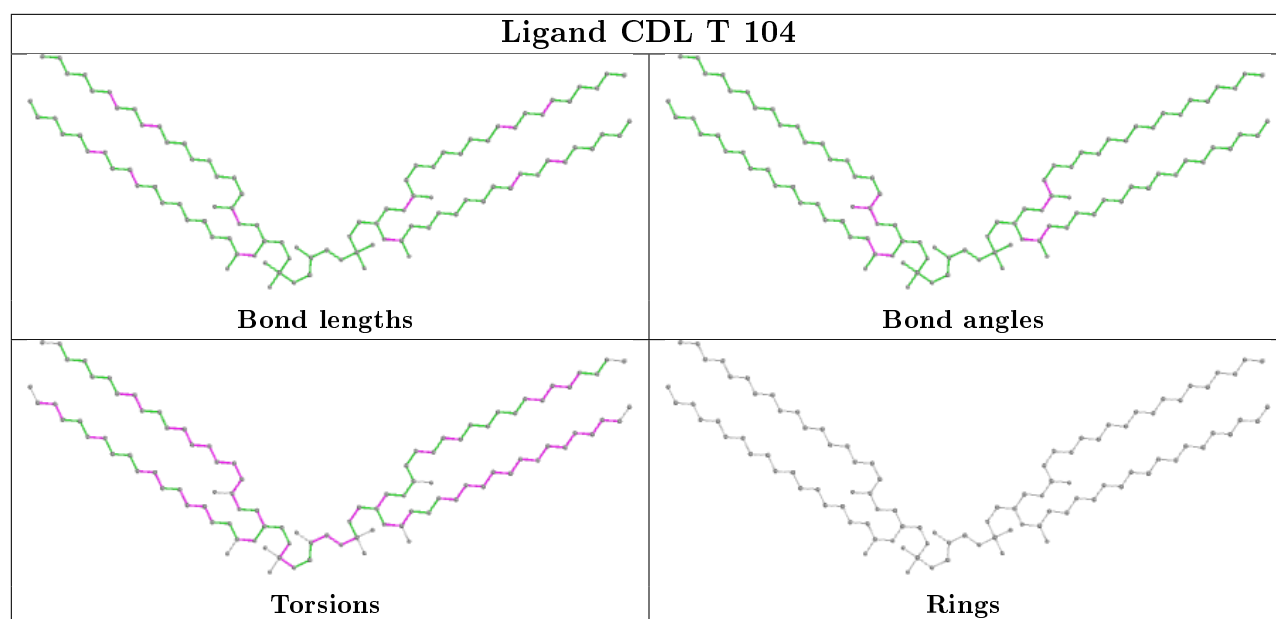
Ligand DMU C 313**Ligand PGV P 303**

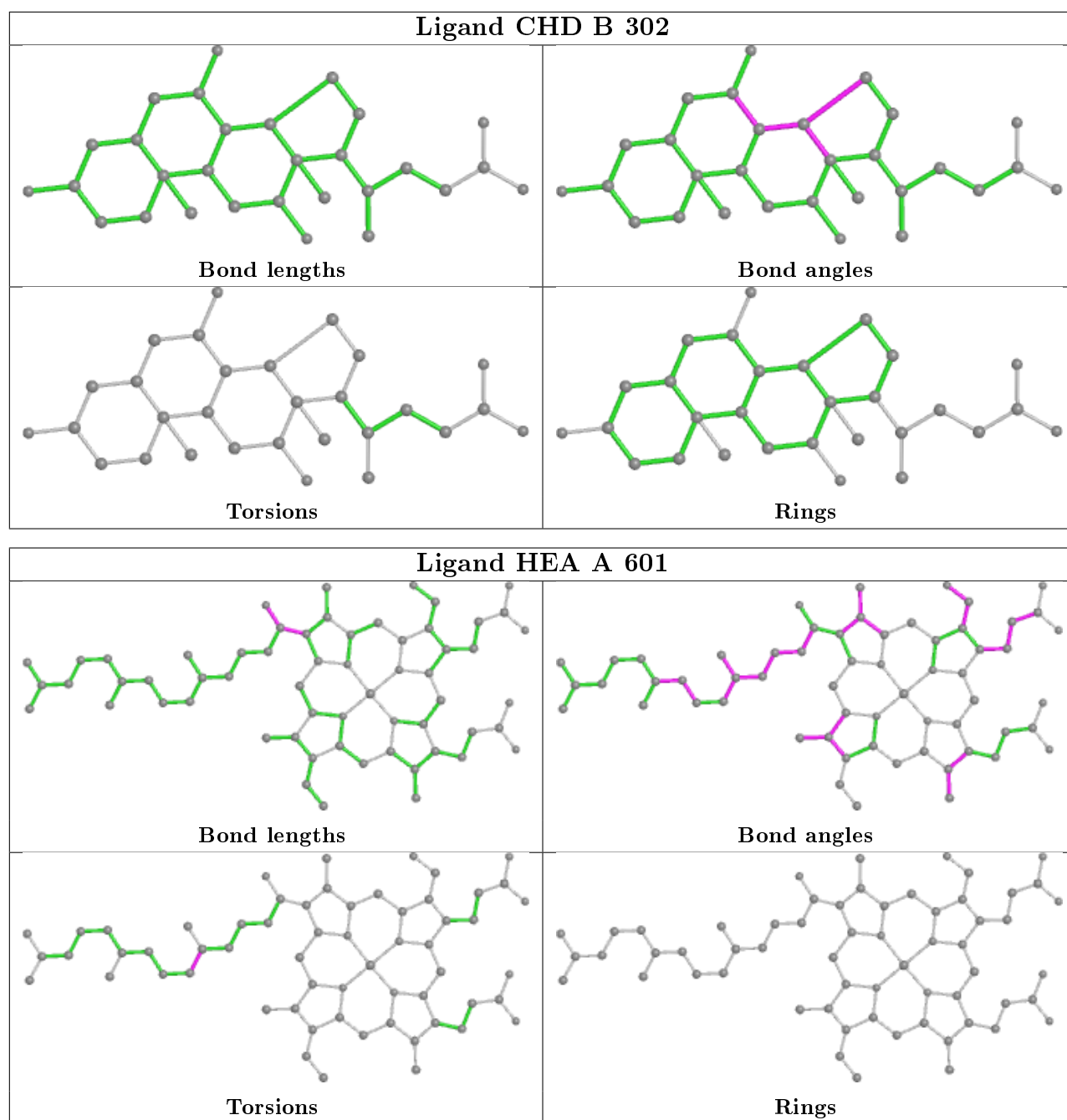


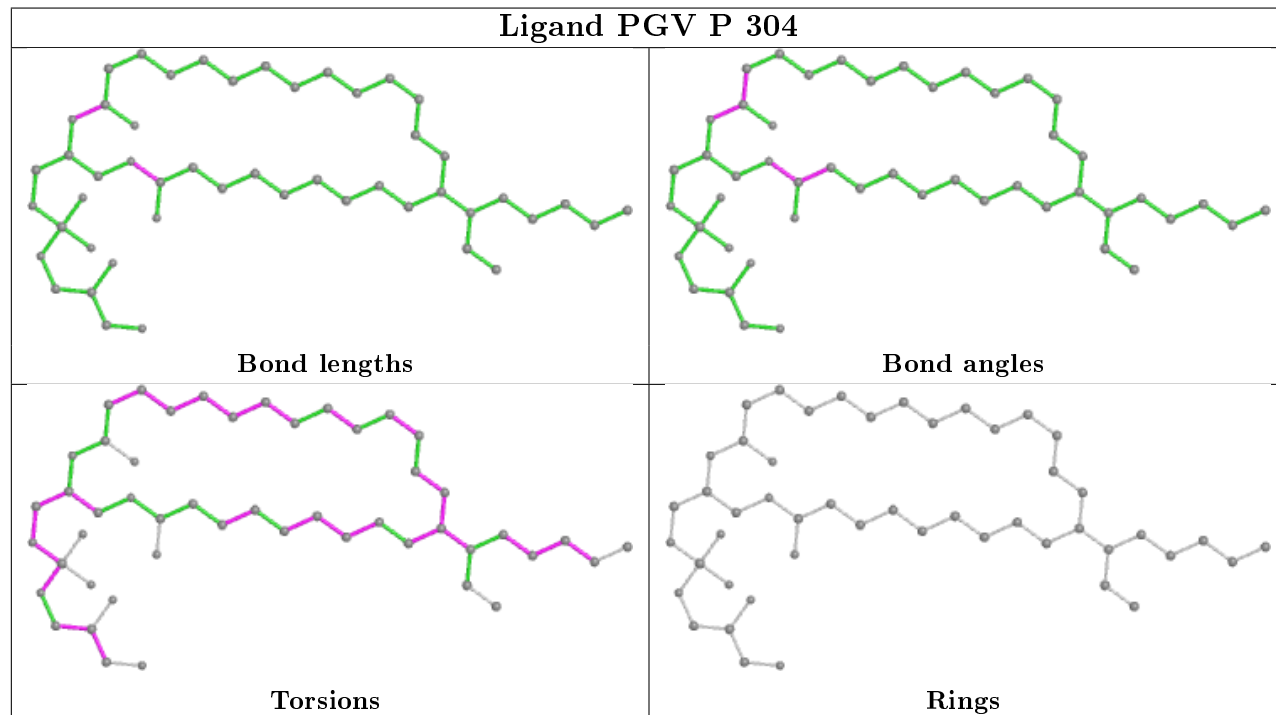
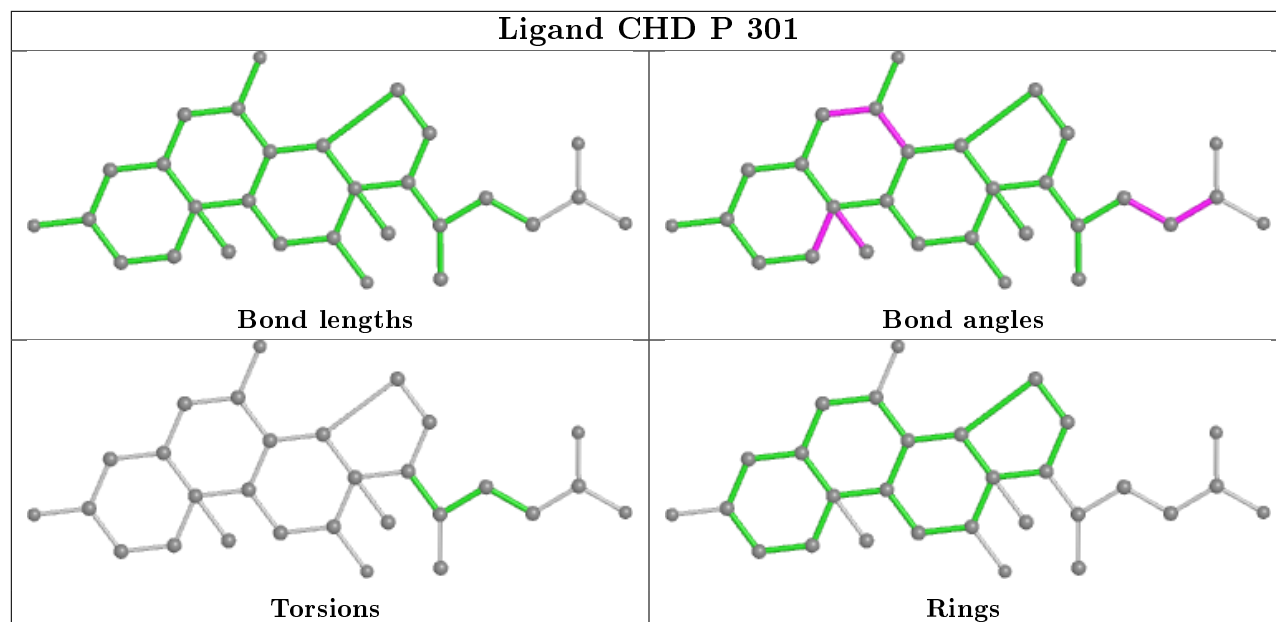


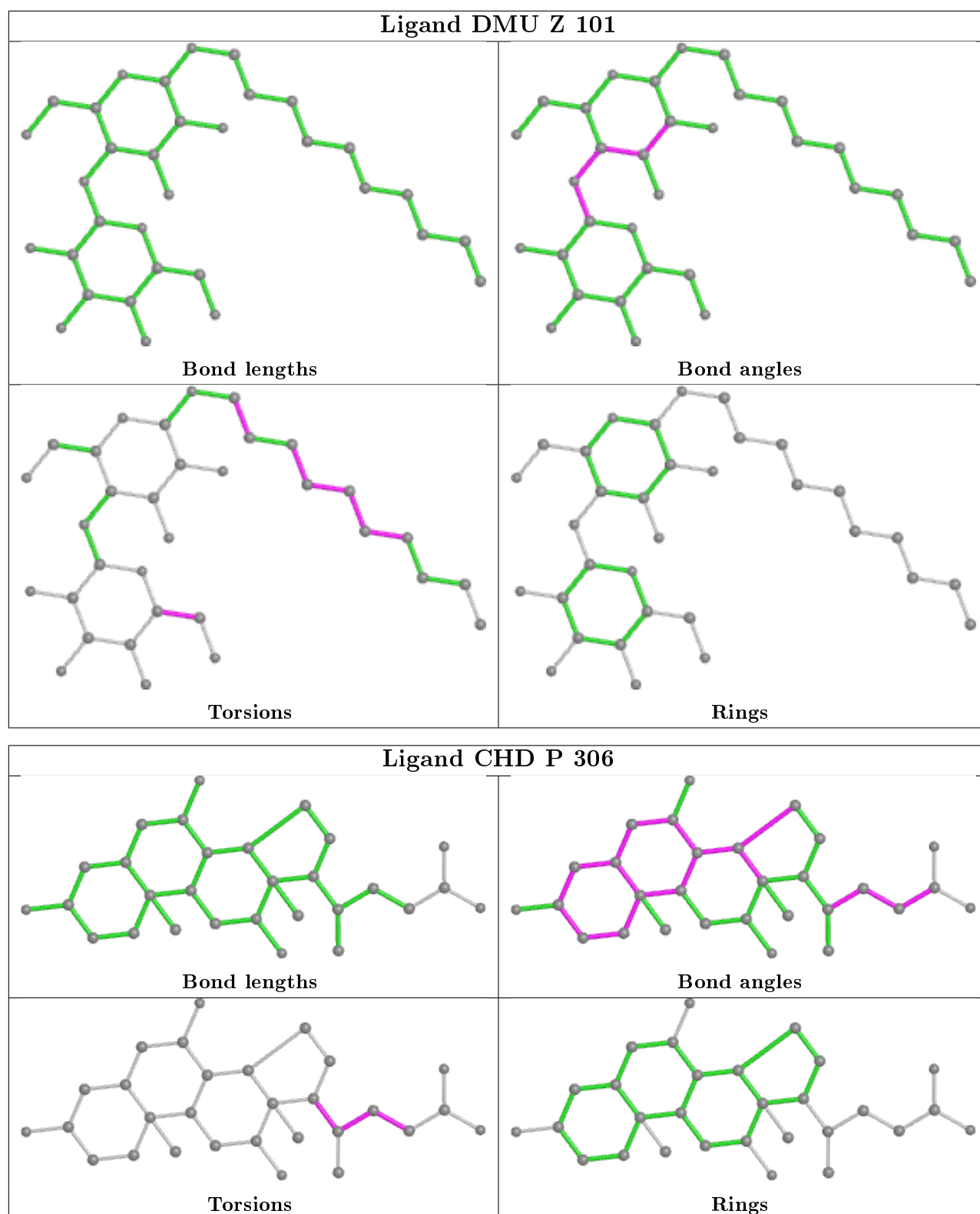




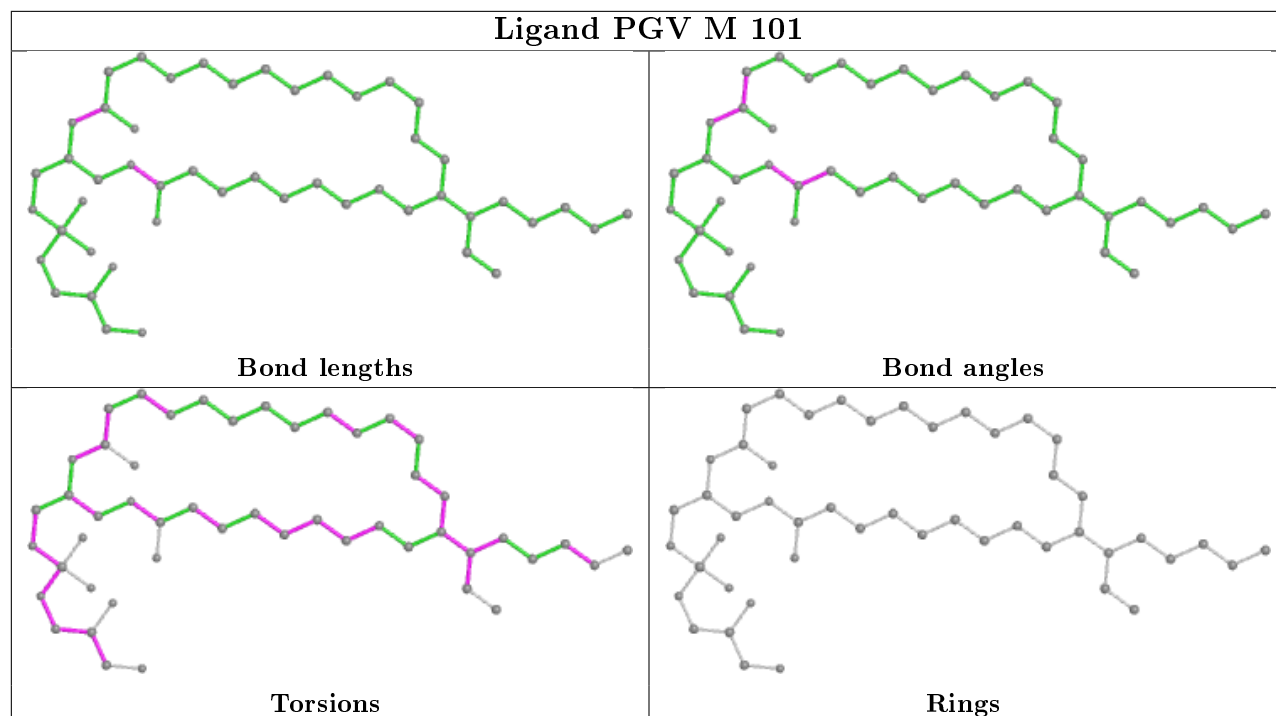




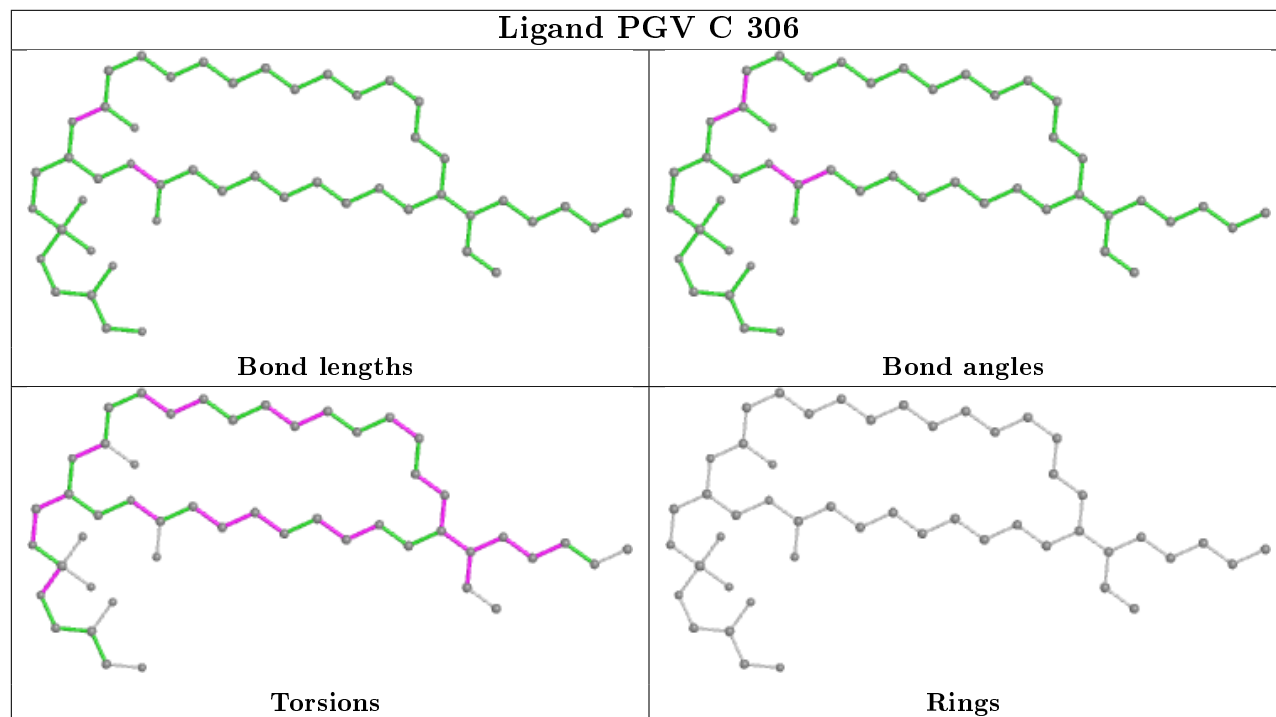


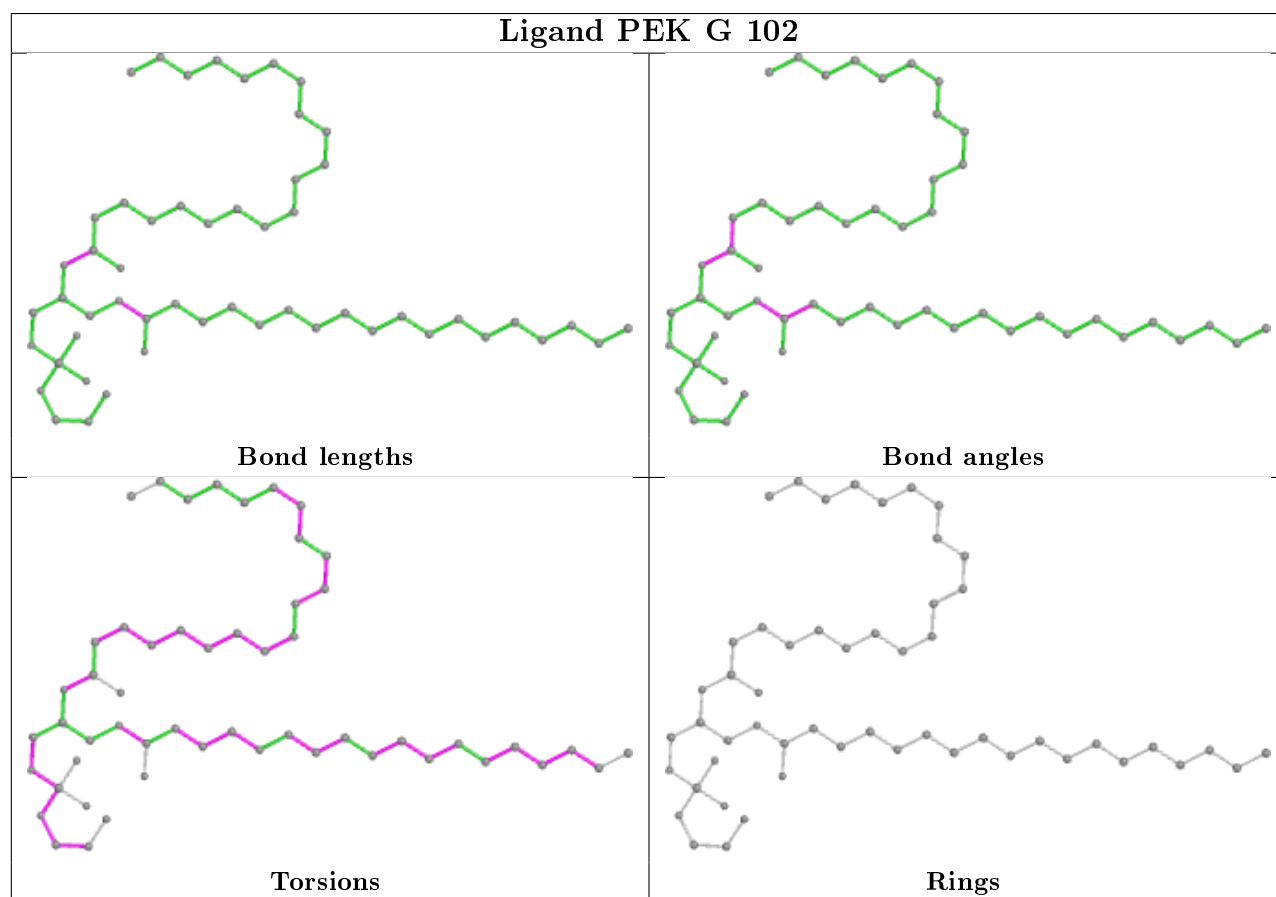
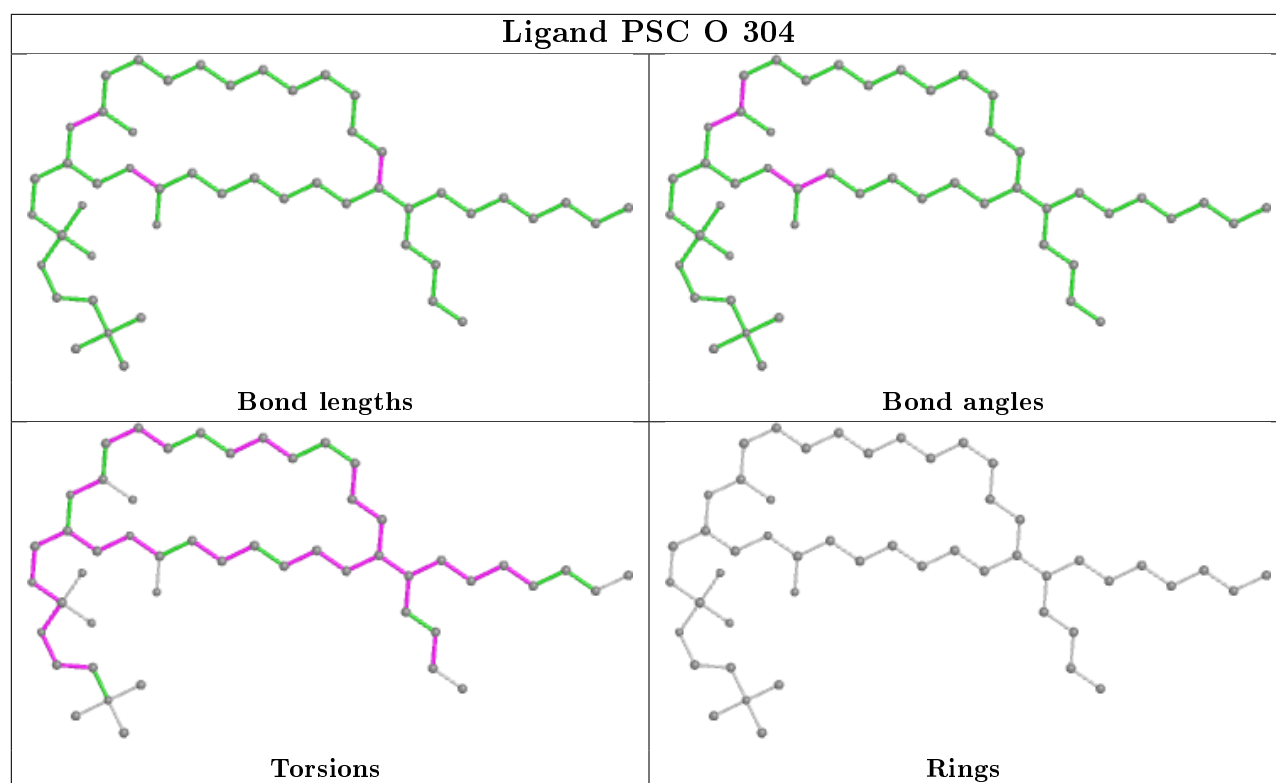


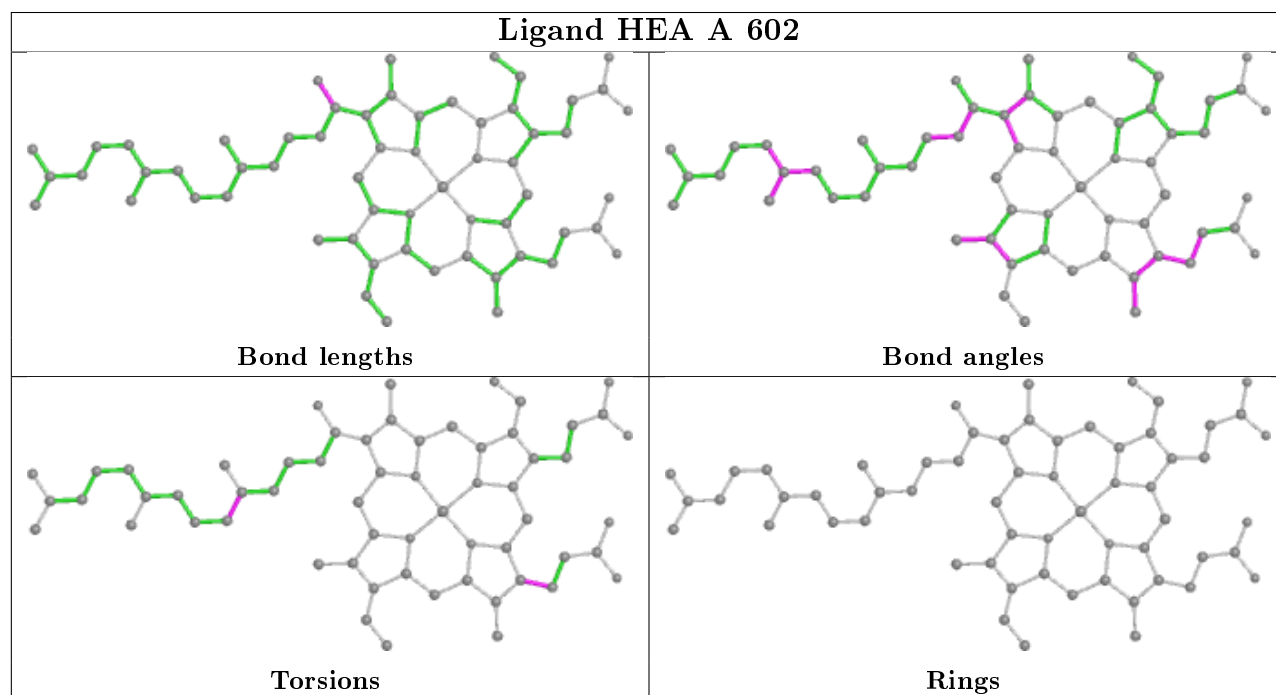
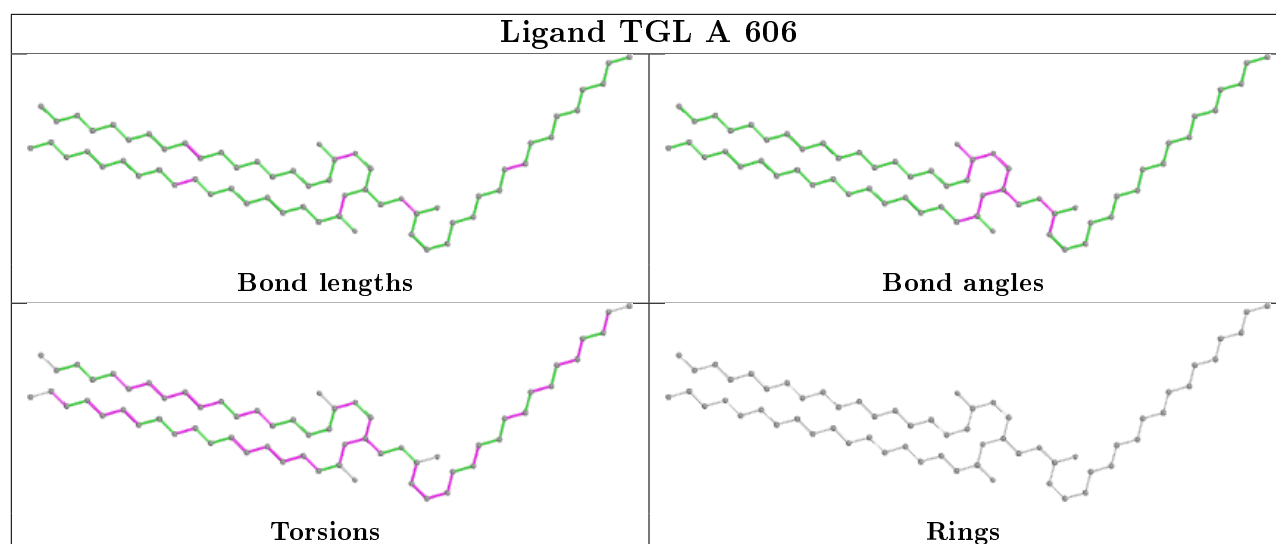
Ligand PGV M 101

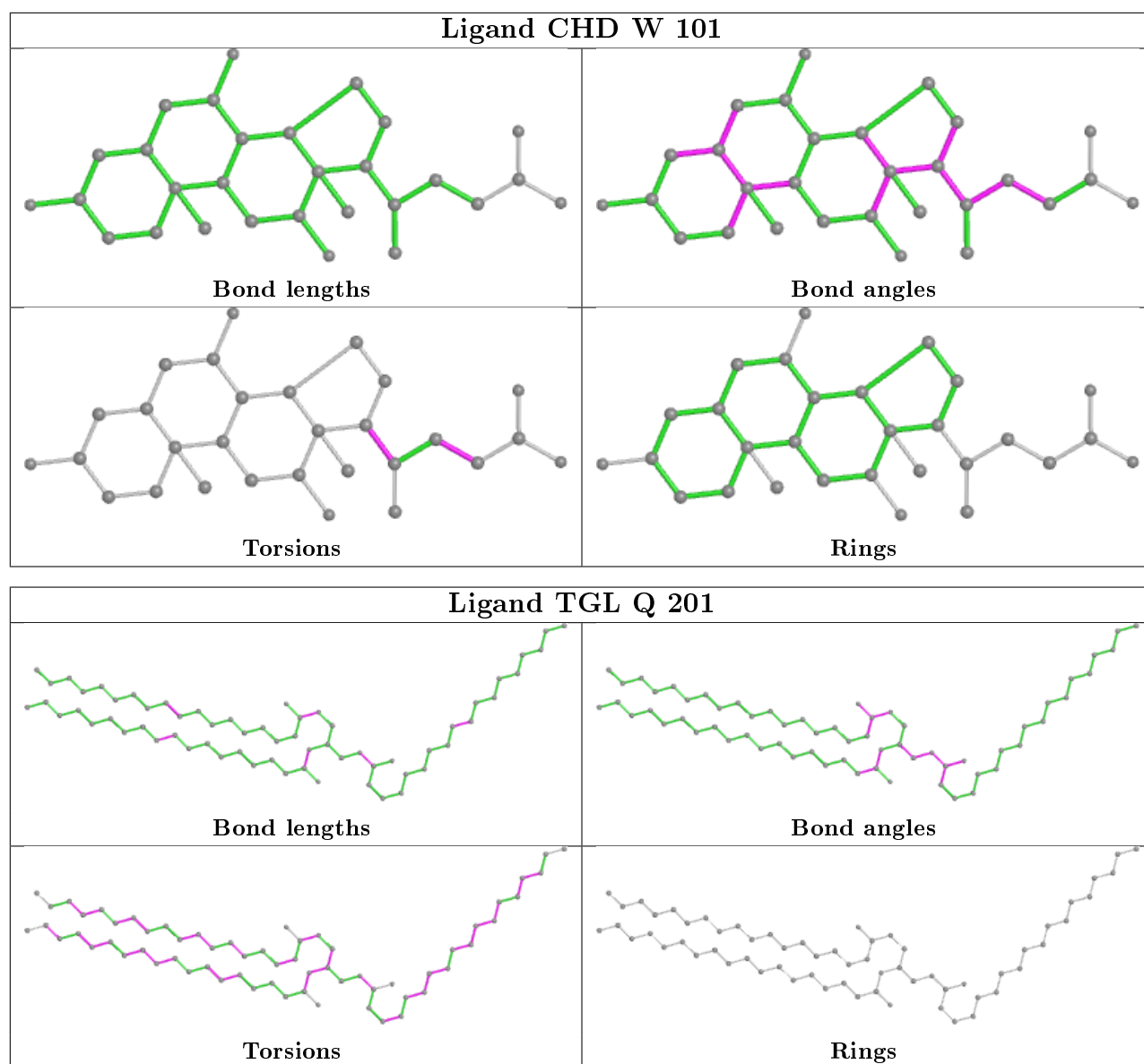


Ligand PGV C 306









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	513/514 (99%)	0.56	6 (1%) 79 78	25, 31, 39, 76	0
1	N	513/514 (99%)	0.44	7 (1%) 75 74	32, 41, 53, 77	0
2	B	226/227 (99%)	0.50	9 (3%) 38 37	28, 38, 58, 75	0
2	O	226/227 (99%)	0.74	16 (7%) 16 15	40, 52, 72, 99	0
3	C	259/260 (99%)	0.68	8 (3%) 49 48	27, 36, 48, 82	0
3	P	259/260 (99%)	0.55	6 (2%) 60 59	33, 40, 51, 78	0
4	D	144/147 (97%)	0.32	2 (1%) 75 74	34, 43, 55, 75	0
4	Q	144/147 (97%)	1.26	21 (14%) 2 2	48, 63, 88, 137	0
5	E	105/109 (96%)	0.36	2 (1%) 66 65	35, 44, 63, 109	0
5	R	105/109 (96%)	0.63	6 (5%) 23 23	45, 55, 71, 115	0
6	F	94/94 (100%)	0.71	3 (3%) 47 46	32, 46, 64, 119	0
6	S	94/94 (100%)	0.66	5 (5%) 26 25	37, 51, 75, 117	0
7	G	83/85 (97%)	1.47	18 (21%) 0 0	34, 44, 114, 145	0
7	T	83/85 (97%)	1.82	22 (26%) 0 0	35, 48, 108, 150	0
8	H	79/85 (92%)	0.69	9 (11%) 5 4	36, 47, 80, 99	0
8	U	79/85 (92%)	0.99	11 (13%) 2 2	43, 54, 93, 109	0
9	I	72/73 (98%)	1.18	12 (16%) 1 1	36, 50, 82, 91	0
9	V	72/73 (98%)	1.44	13 (18%) 1 1	42, 67, 95, 103	0
10	J	58/59 (98%)	0.83	5 (8%) 10 9	38, 47, 73, 117	0
10	W	58/59 (98%)	1.04	8 (13%) 2 2	44, 55, 81, 128	0
11	K	49/56 (87%)	0.38	1 (2%) 65 63	38, 45, 58, 66	0
11	X	49/56 (87%)	0.87	6 (12%) 4 3	56, 66, 83, 90	0
12	L	46/47 (97%)	0.44	1 (2%) 62 60	31, 38, 53, 84	0
12	Y	46/47 (97%)	0.82	5 (10%) 5 5	46, 54, 70, 108	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	M	43/46 (93%)	0.56	3 (6%) 16 15	35, 38, 66, 104	0
13	Z	43/46 (93%)	0.97	8 (18%) 1 1	52, 58, 85, 118	0
All	All	3542/3604 (98%)	0.70	213 (6%) 21 20	25, 43, 75, 150	0

The worst 5 of 213 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	Q	6	VAL	23.2
4	Q	5	VAL	17.0
7	T	8	HIS	13.0
7	G	3	ALA	12.7
6	F	1	ALA	12.5

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	TPO	T	11	11/12	0.48	0.35	93,108,122,125	0
7	TPO	G	11	11/12	0.58	0.30	88,102,120,124	0
9	SAC	V	1	9/10	0.73	0.60	118,123,127,131	0
9	SAC	I	1	9/10	0.82	0.20	74,75,78,79	0
1	FME	A	1	10/11	0.93	0.16	46,53,70,78	0
1	FME	N	1	10/11	0.94	0.21	58,63,83,86	0
2	FME	O	1	10/11	0.96	0.21	52,53,59,60	0
2	FME	B	1	10/11	0.97	0.15	35,36,44,47	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
20	EDO	C	310	4/4	0.37	0.22	71,76,80,84	0
25	DMU	P	308	33/33	0.43	0.28	71,126,142,143	0
23	PEK	C	304	53/53	0.47	0.27	60,91,135,141	0
17	NA	C	302	1/1	0.52	0.20	51,51,51,51	0
23	PEK	T	103	53/53	0.54	0.26	59,85,124,134	0
23	PEK	T	101	53/53	0.59	0.33	56,94,139,144	0
24	CDL	G	101	100/100	0.60	0.40	65,97,136,145	0
24	CDL	T	104	100/100	0.61	0.38	63,92,121,138	0
18	TGL	Y	101	63/63	0.61	0.31	63,83,104,107	0
18	TGL	Q	201	63/63	0.63	0.25	67,86,96,100	0
24	CDL	P	305	100/100	0.63	0.33	65,90,111,114	0
25	DMU	C	313	33/33	0.64	0.53	49,130,146,147	0
19	PGV	N	606	51/51	0.65	0.40	70,87,126,128	0
23	PEK	G	102	53/53	0.66	0.29	55,101,139,150	0
22	CHD	W	101	29/29	0.66	0.36	89,96,106,107	0
26	PSC	E	201	52/52	0.67	0.36	56,87,141,150	0
19	PGV	P	304	51/51	0.68	0.33	66,92,122,129	0
25	DMU	P	309	33/33	0.69	0.46	64,128,146,149	0
24	CDL	C	307	100/100	0.69	0.34	56,85,110,116	0
19	PGV	M	101	51/51	0.70	0.31	49,77,115,124	0
20	EDO	C	309	4/4	0.70	0.27	82,86,90,92	0
25	DMU	C	312	33/33	0.70	0.25	65,100,138,139	0
19	PGV	C	306	51/51	0.72	0.30	62,82,121,124	0
18	TGL	D	201	63/63	0.73	0.21	56,71,89,92	0
18	TGL	L	101	63/63	0.73	0.29	46,71,85,86	0
22	CHD	J	101	29/29	0.73	0.27	73,78,100,104	0
17	NA	P	302	1/1	0.73	0.12	51,51,51,51	0
26	PSC	O	304	52/52	0.73	0.31	60,92,142,154	0
18	TGL	O	302	63/63	0.75	0.32	64,87,99,105	0
18	TGL	A	606	63/63	0.80	0.28	53,78,95,98	0
20	EDO	N	611	4/4	0.81	0.17	60,62,64,66	0
20	EDO	F	104	4/4	0.83	0.29	56,58,60,62	0
20	EDO	A	615	4/4	0.84	0.37	49,50,56,58	0
20	EDO	D	203	4/4	0.84	0.23	61,65,70,71	0
25	DMU	Z	101	33/33	0.84	0.26	57,80,91,94	0
16	MG	N	604	1/1	0.85	0.17	44,44,44,44	0
20	EDO	B	303	4/4	0.85	0.52	62,72,75,79	0
20	EDO	A	611	4/4	0.85	0.75	62,67,67,69	0
20	EDO	D	202	4/4	0.86	0.32	54,63,65,68	0
20	EDO	P	307	4/4	0.86	0.29	82,88,90,97	0
22	CHD	C	308	29/29	0.87	0.25	72,74,79,80	0
20	EDO	S	103	4/4	0.87	0.21	57,66,69,72	0

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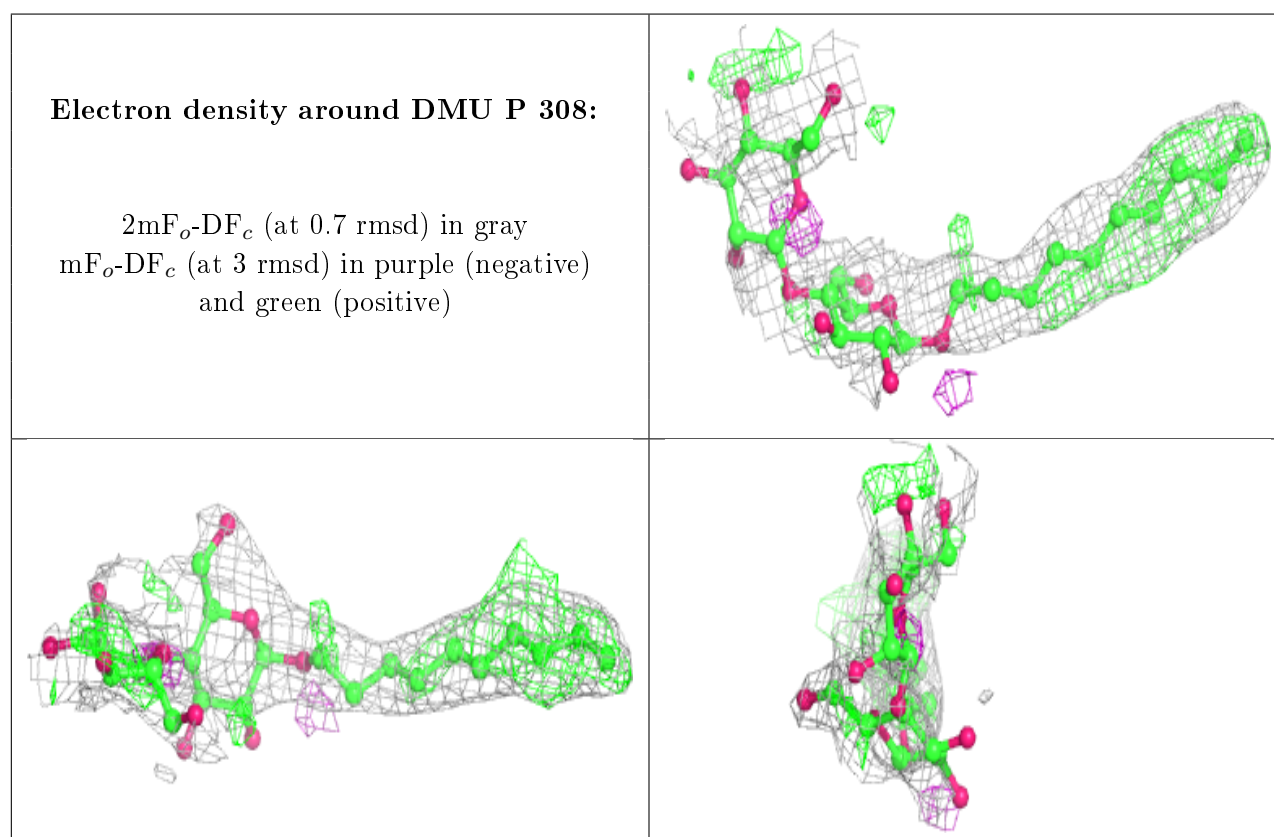
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
22	CHD	P	306	29/29	0.87	0.19	71,76,81,81	0
20	EDO	A	608	4/4	0.88	0.12	55,59,59,65	0
20	EDO	N	610	4/4	0.89	0.21	62,64,65,66	0
25	DMU	M	102	33/33	0.90	0.15	48,58,74,77	0
20	EDO	R	201	4/4	0.91	0.34	52,53,53,53	0
20	EDO	A	613	4/4	0.92	0.15	46,49,51,52	0
20	EDO	N	608	4/4	0.93	0.14	37,39,40,40	0
20	EDO	N	615	4/4	0.93	0.39	54,58,59,65	0
20	EDO	A	614	4/4	0.94	0.24	50,53,54,59	0
20	EDO	N	614	4/4	0.94	0.19	46,48,49,50	0
20	EDO	A	616	4/4	0.94	0.23	43,45,49,49	0
20	EDO	A	612	4/4	0.94	0.22	38,39,43,43	0
20	EDO	N	609	4/4	0.94	0.28	63,69,69,74	0
20	EDO	N	617	4/4	0.94	0.21	58,70,75,86	0
20	EDO	A	610	4/4	0.94	0.21	54,56,56,57	0
20	EDO	E	202	4/4	0.95	0.13	51,52,52,53	0
22	CHD	P	301	29/29	0.95	0.11	37,40,42,45	0
16	MG	A	604	1/1	0.95	0.16	30,30,30,30	0
23	PEK	T	102	53/53	0.95	0.20	40,58,95,98	0
22	CHD	C	301	29/29	0.96	0.11	34,37,39,42	0
23	PEK	C	303	53/53	0.96	0.18	35,52,84,86	0
20	EDO	F	102	4/4	0.96	0.16	41,44,45,47	0
20	EDO	N	612	4/4	0.96	0.17	45,46,46,47	0
20	EDO	N	613	4/4	0.96	0.35	49,51,53,55	0
22	CHD	B	302	29/29	0.96	0.14	36,38,40,44	0
19	PGV	P	303	51/51	0.96	0.19	38,46,76,79	0
19	PGV	A	607	51/51	0.97	0.17	33,40,61,66	0
22	CHD	O	301	29/29	0.97	0.14	40,42,44,47	0
20	EDO	C	311	4/4	0.97	0.11	40,43,45,48	0
20	EDO	N	616	4/4	0.97	0.23	59,59,61,61	0
14	HEA	N	602	60/60	0.97	0.14	35,37,42,44	0
19	PGV	N	607	51/51	0.97	0.15	38,46,68,70	0
19	PGV	C	305	51/51	0.97	0.18	35,41,90,101	0
14	HEA	N	601	60/60	0.97	0.14	36,42,49,51	0
17	NA	A	605	1/1	0.98	0.08	30,30,30,30	0
20	EDO	S	102	4/4	0.98	0.13	37,39,40,42	0
20	EDO	F	103	4/4	0.98	0.14	31,33,33,33	0
21	CUA	B	301	2/2	0.98	0.17	30,30,30,30	0
14	HEA	A	602	60/60	0.98	0.13	26,28,34,38	0
20	EDO	A	609	4/4	0.98	0.14	30,32,32,33	0
14	HEA	A	601	60/60	0.98	0.14	25,28,38,40	0
27	ZN	S	101	1/1	0.99	0.15	46,46,46,46	0

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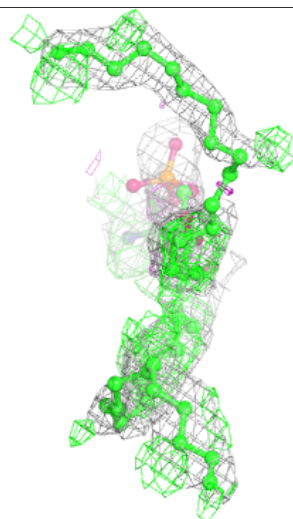
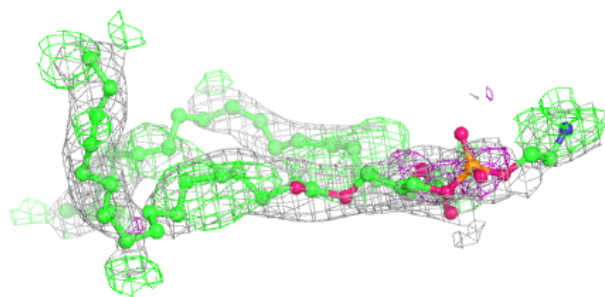
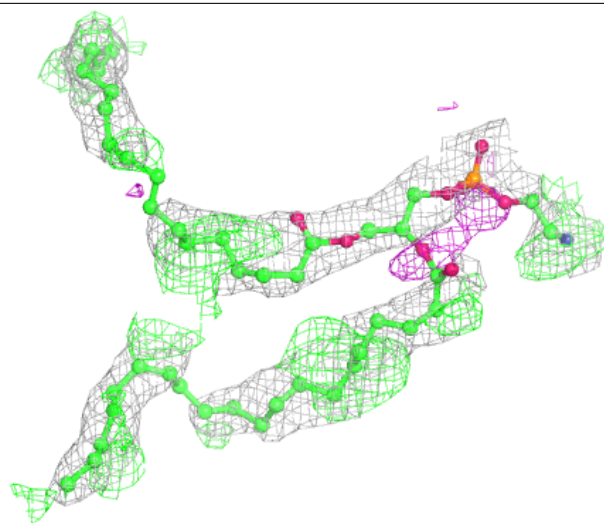
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
21	CUA	O	303	2/2	0.99	0.14	42,42,42,43	0
17	NA	N	605	1/1	0.99	0.09	49,49,49,49	0
15	CU	N	603	1/1	1.00	0.20	37,37,37,37	0
27	ZN	F	101	1/1	1.00	0.18	39,39,39,39	0
15	CU	A	603	1/1	1.00	0.19	28,28,28,28	0

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



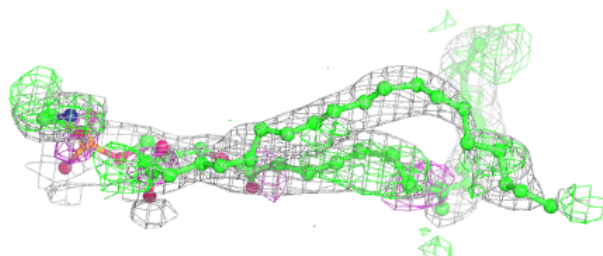
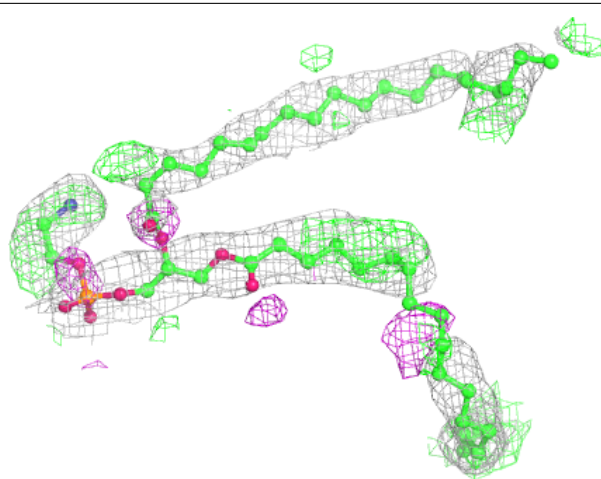
Electron density around PEK C 304:

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



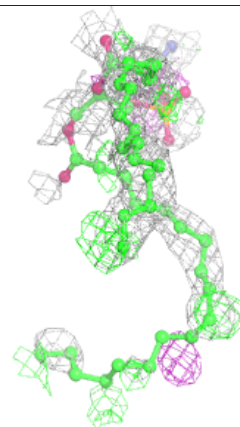
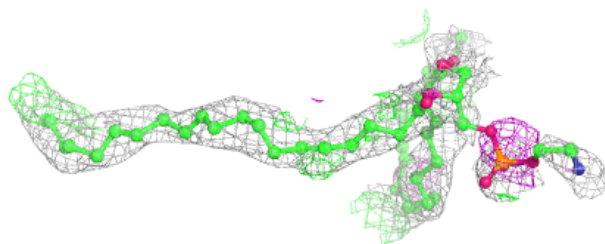
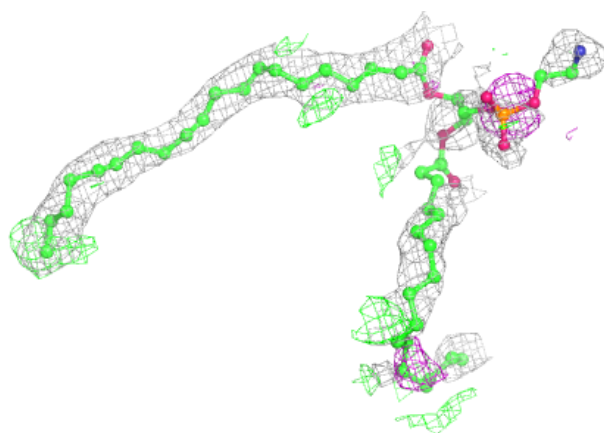
Electron density around PEK T 103:

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

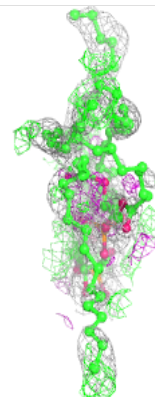
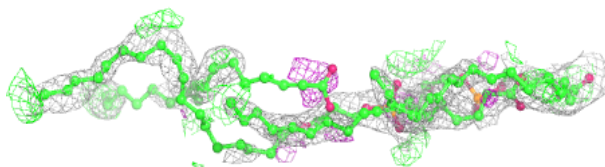
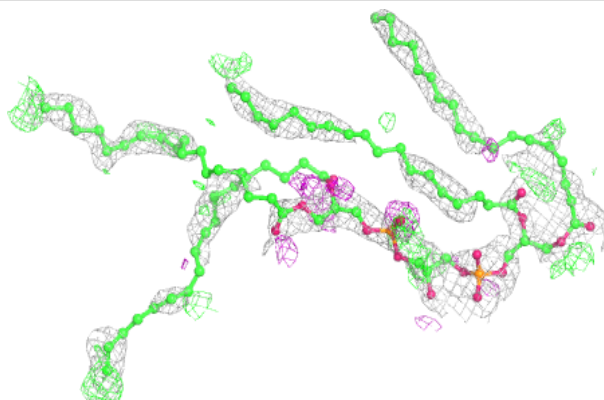


Electron density around PEK T 101:

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

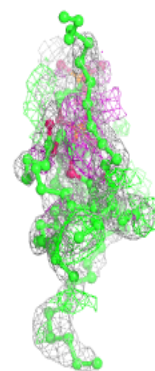
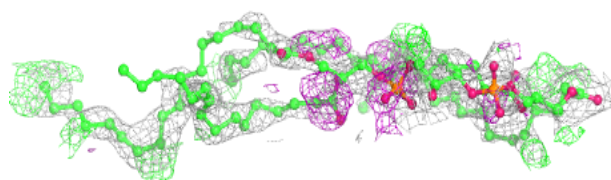
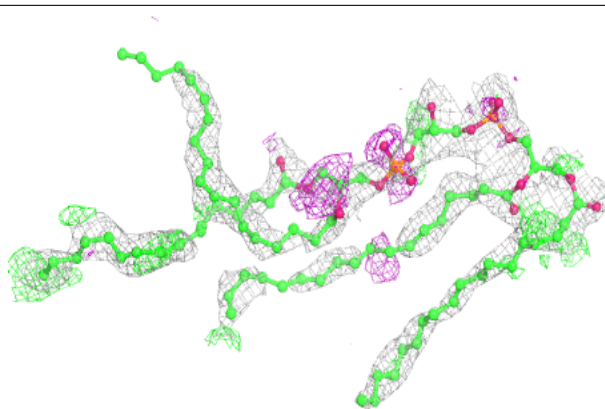
**Electron density around CDL G 101:**

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



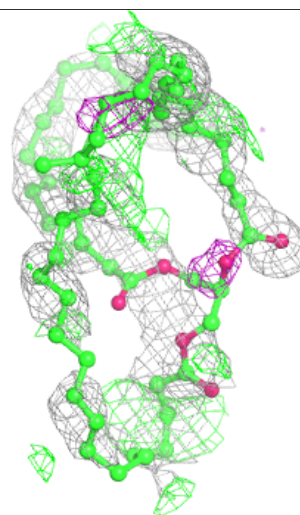
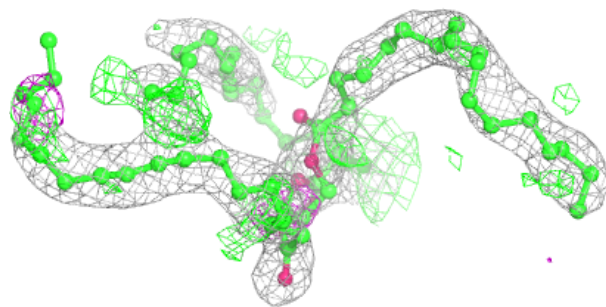
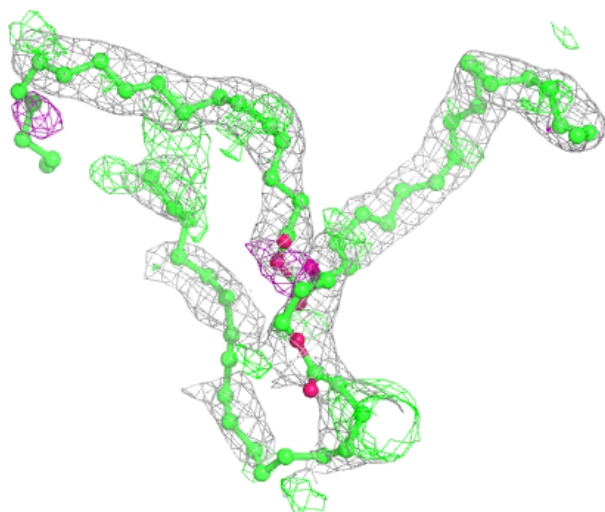
Electron density around CDL T 104:

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



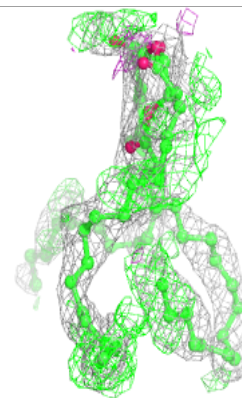
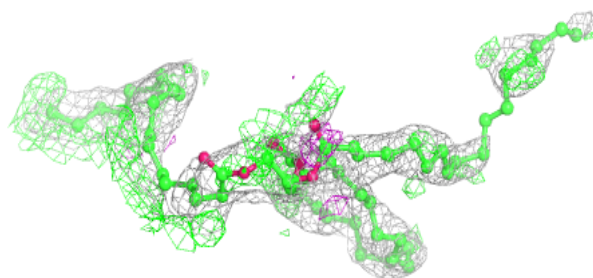
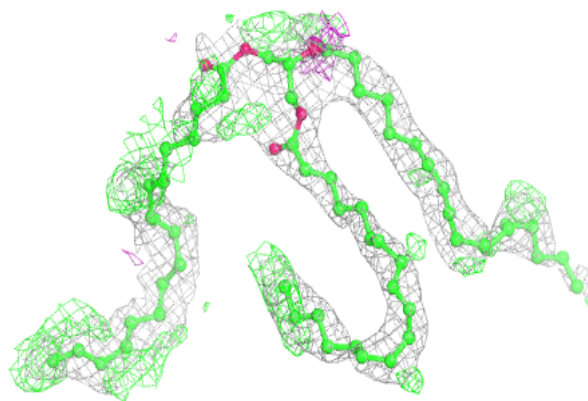
Electron density around TGL Y 101:

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

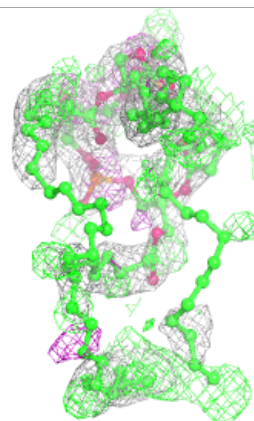
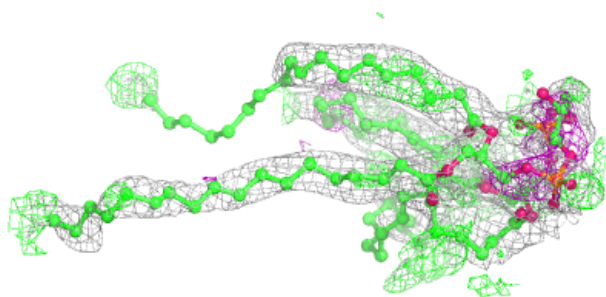
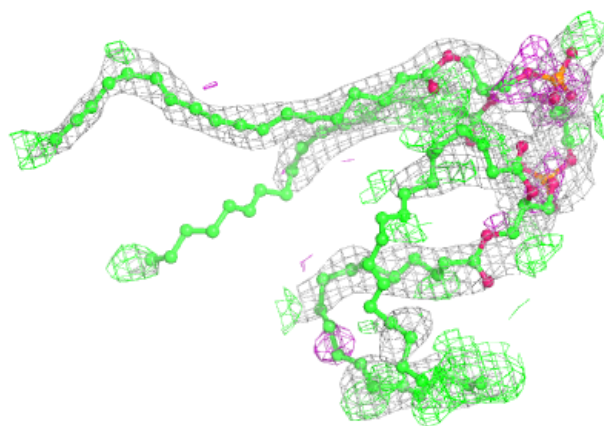


Electron density around TGL Q 201:

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

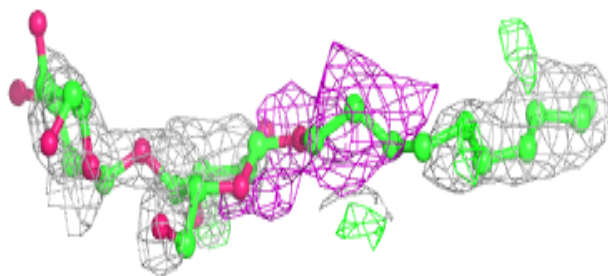
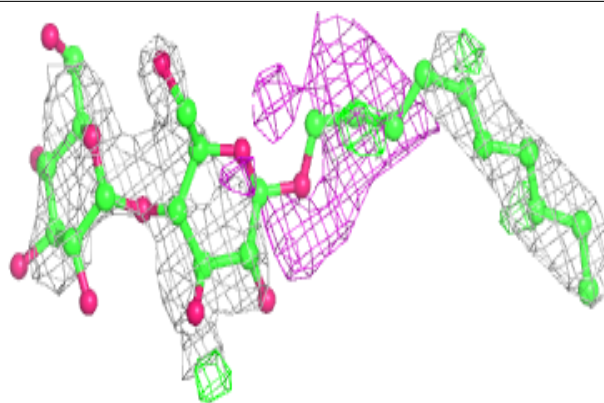
**Electron density around CDL P 305:**

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

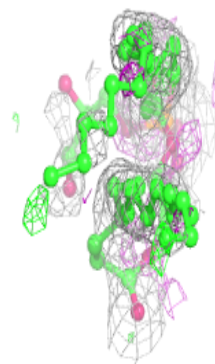
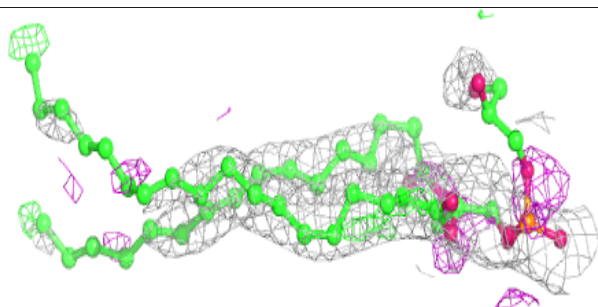
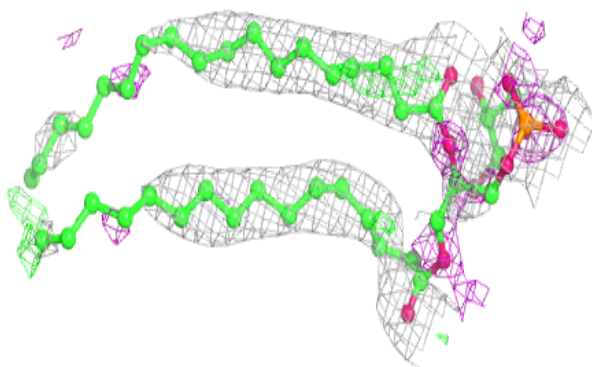


Electron density around DMU C 313:

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

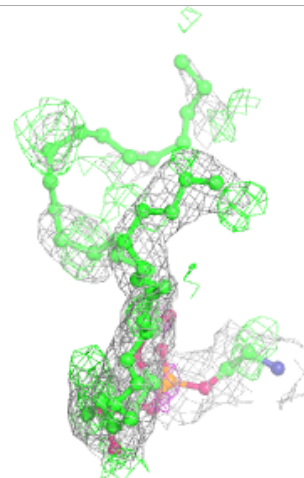
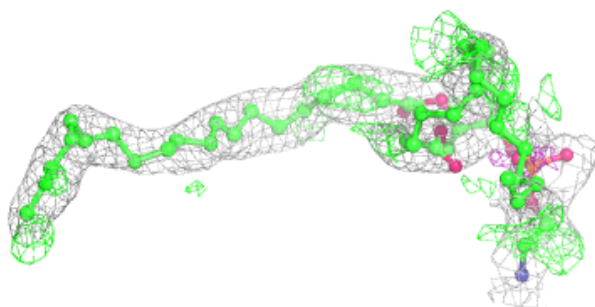
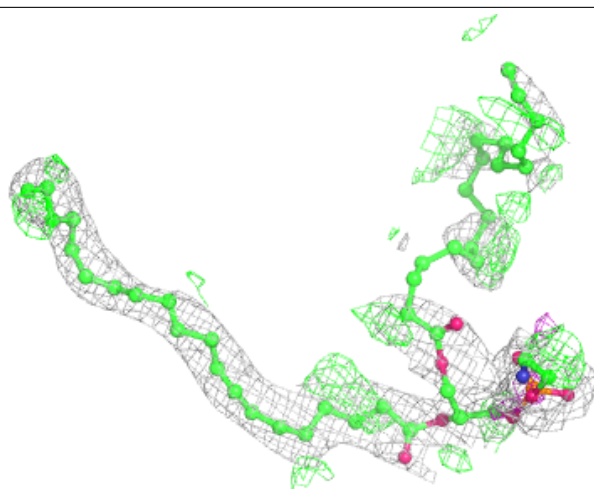
**Electron density around PGV N 606:**

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



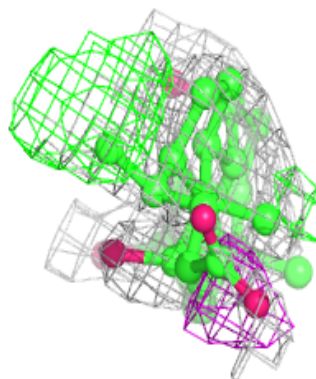
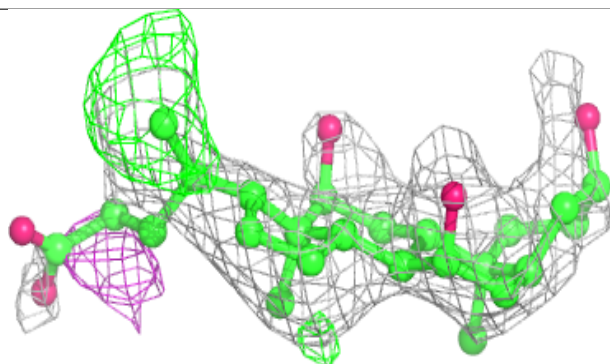
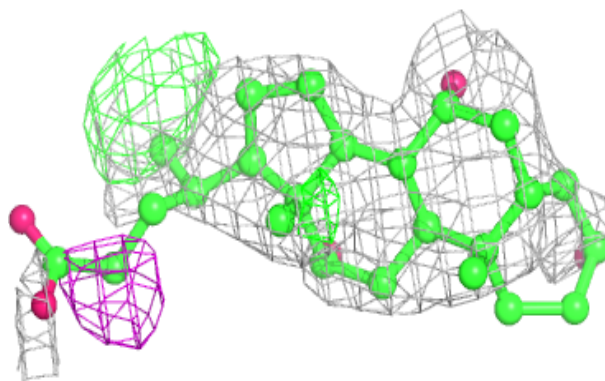
Electron density around PEK G 102:

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

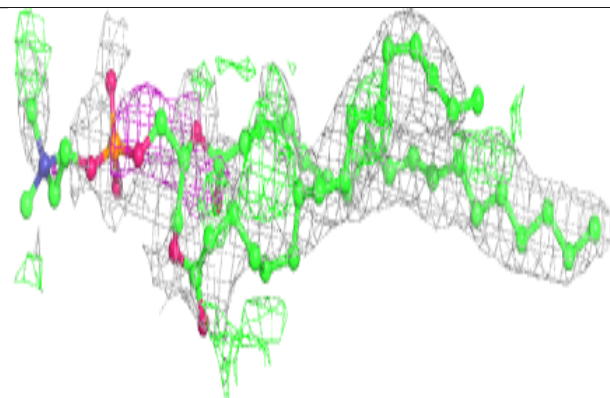
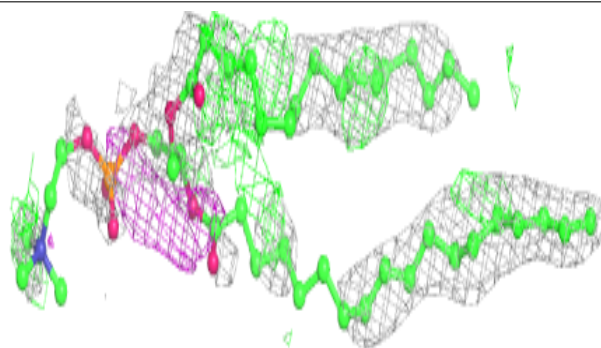


Electron density around CHD W 101:

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

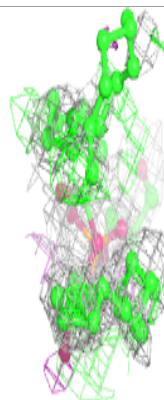
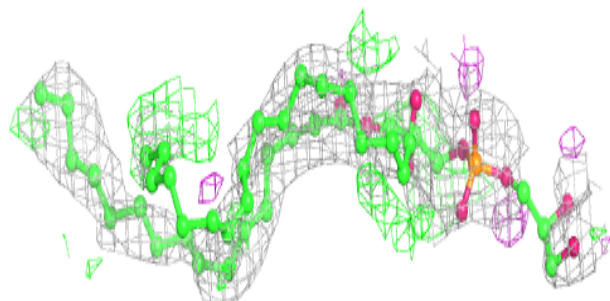
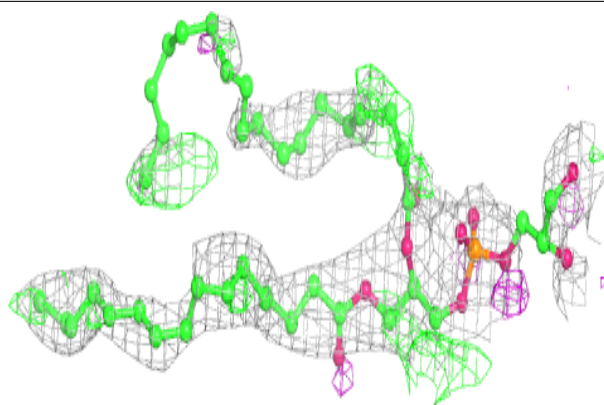
**Electron density around PSC E 201:**

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

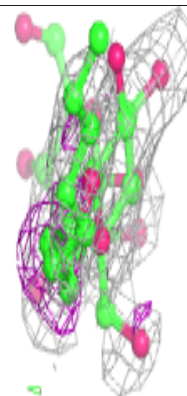
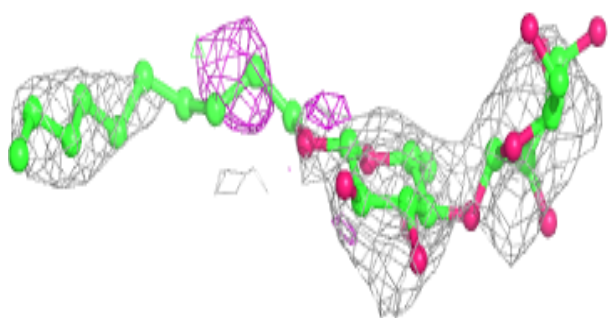
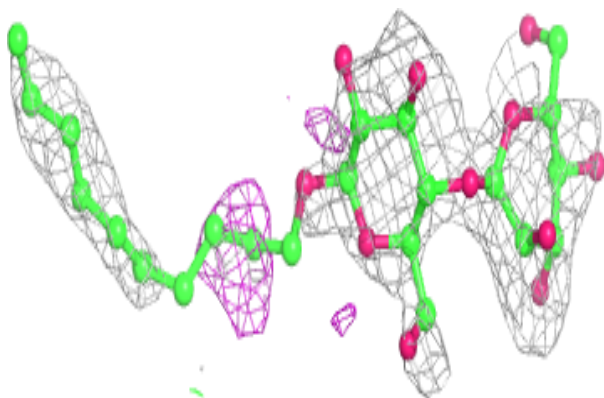


Electron density around PGV P 304:

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

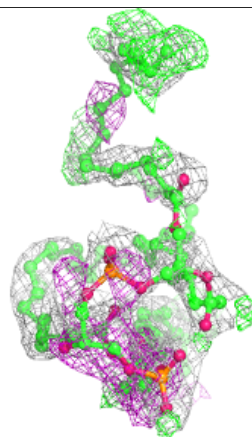
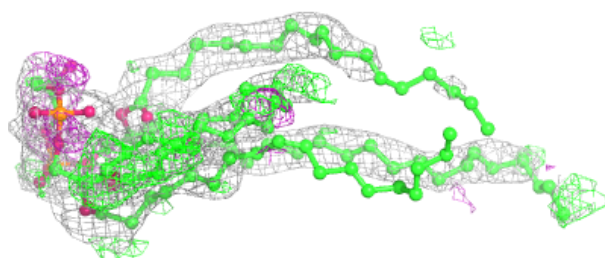
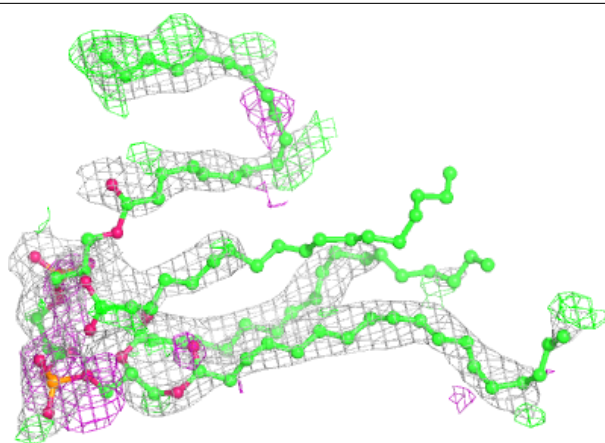
**Electron density around DMU P 309:**

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

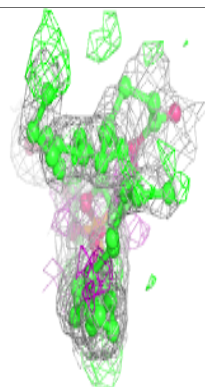
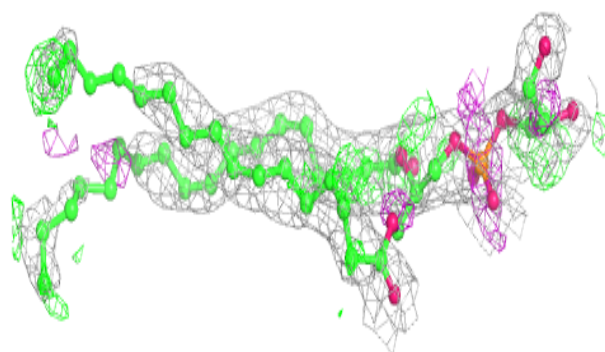
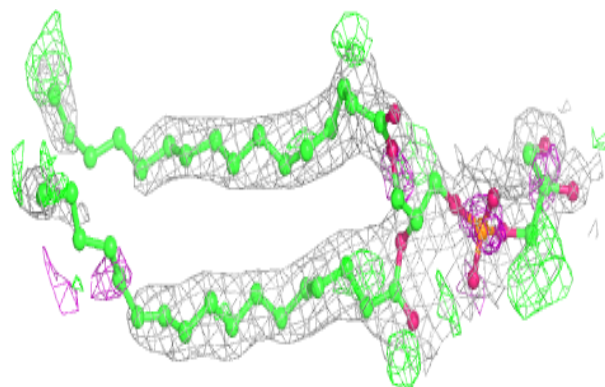


Electron density around CDL C 307:

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

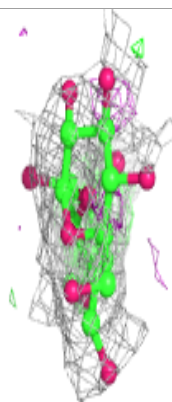
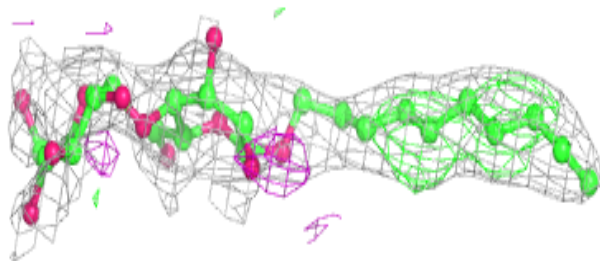
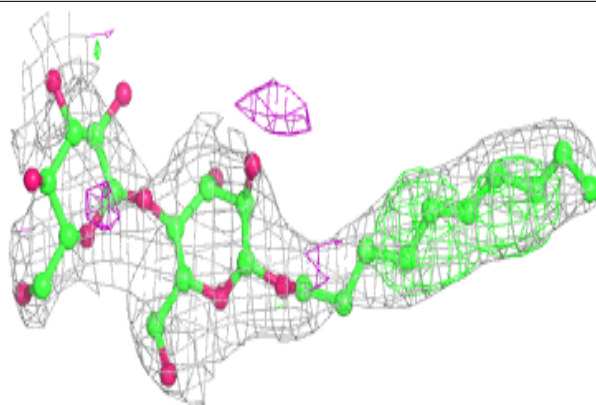
**Electron density around PGV M 101:**

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

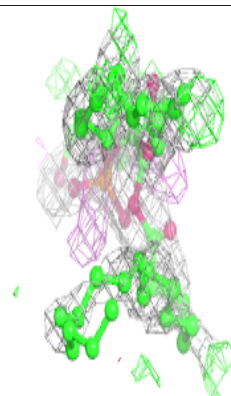
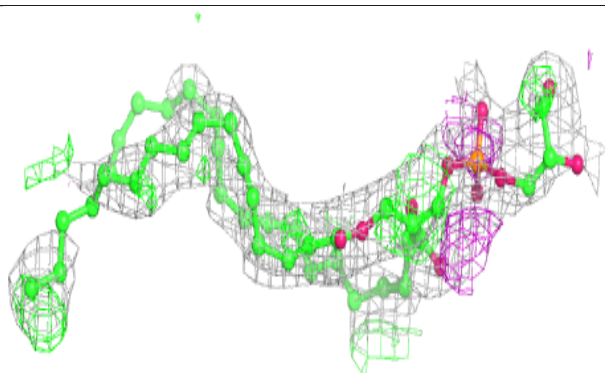
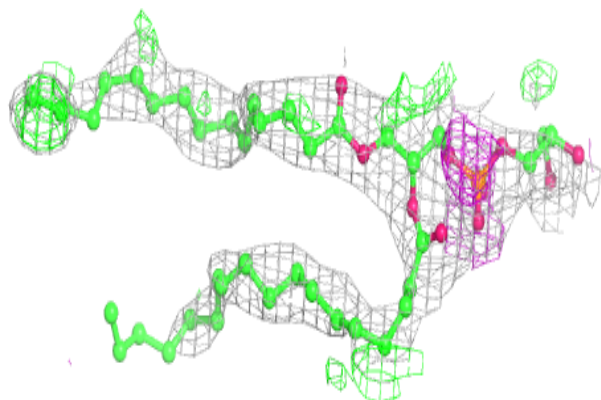


Electron density around DMU C 312:

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

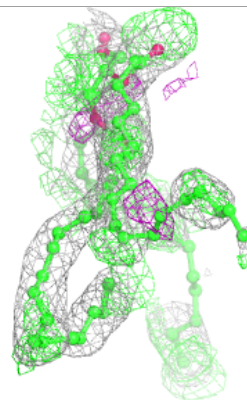
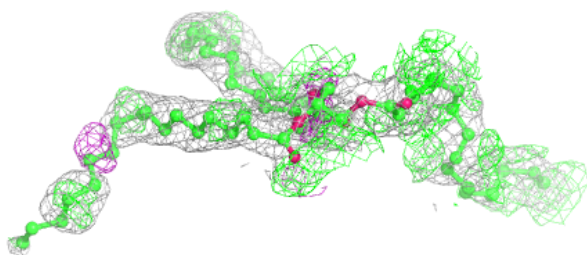
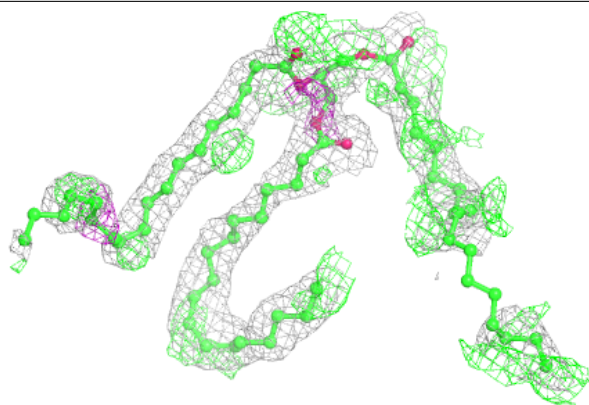
**Electron density around PGV C 306:**

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

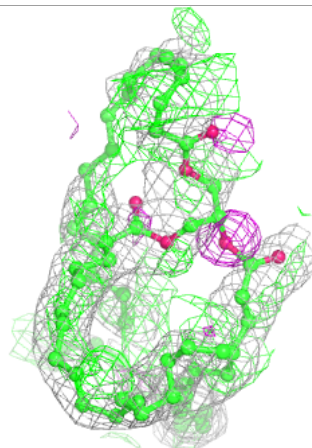
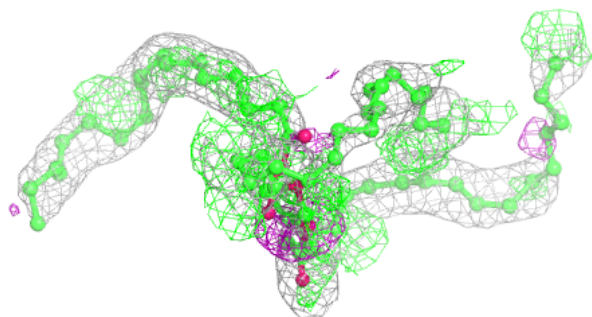
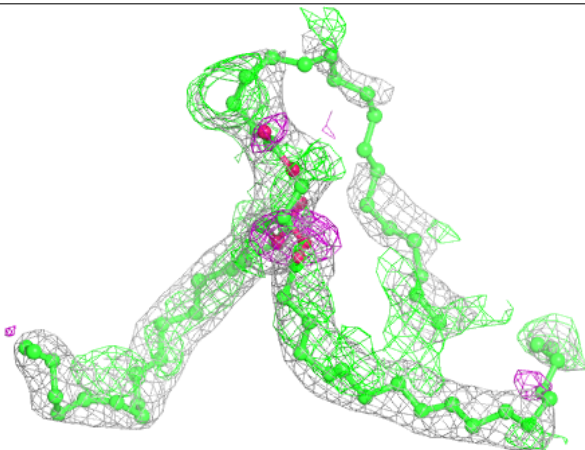


Electron density around TGL D 201:

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

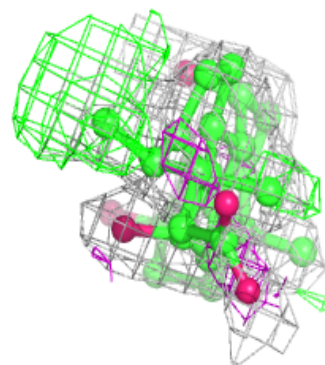
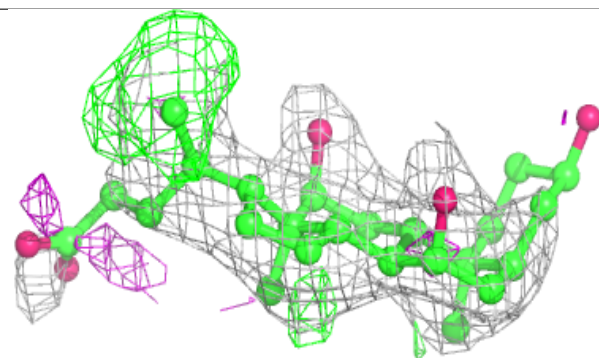
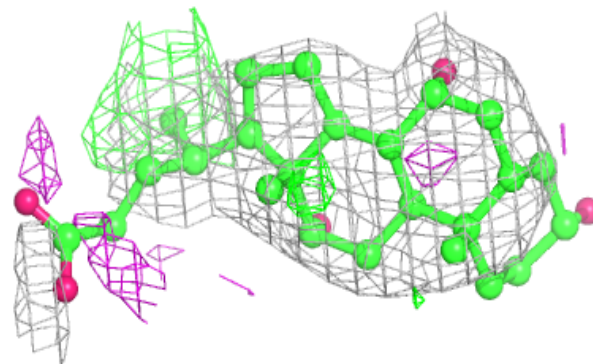
**Electron density around TGL L 101:**

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

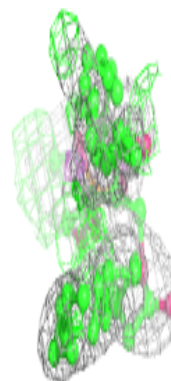
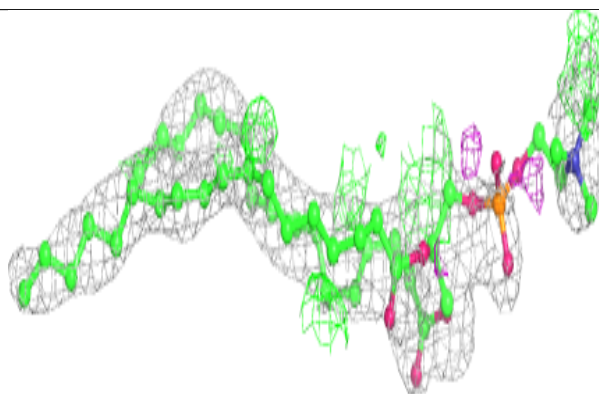
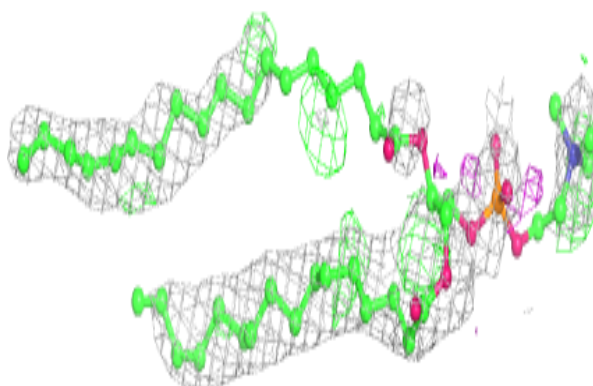


Electron density around CHD J 101:

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

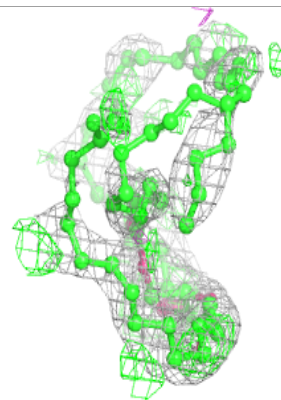
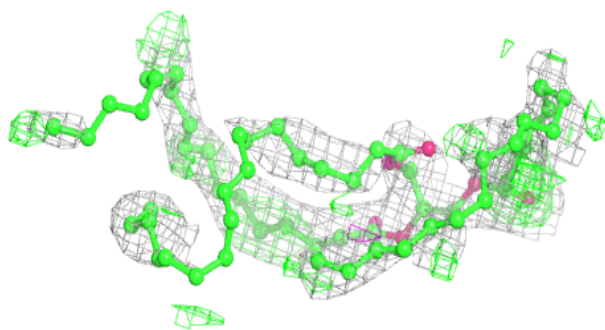
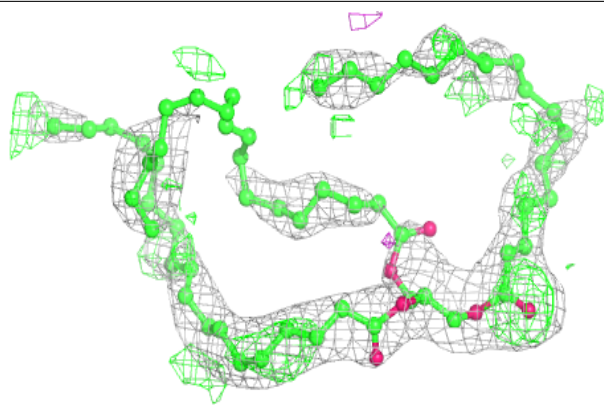
**Electron density around PSC O 304:**

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

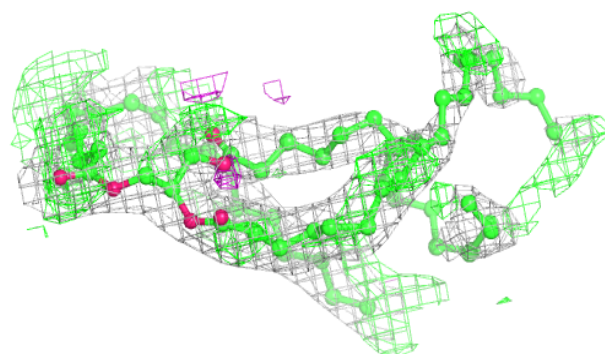
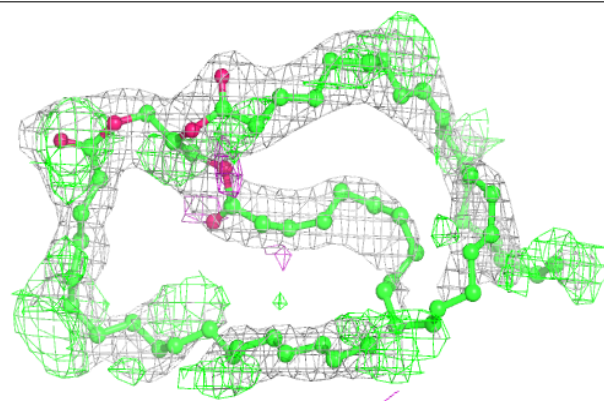


Electron density around TGL O 302:

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

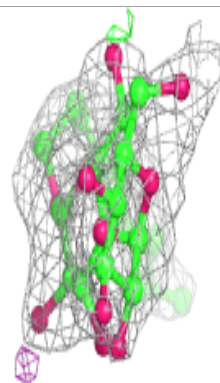
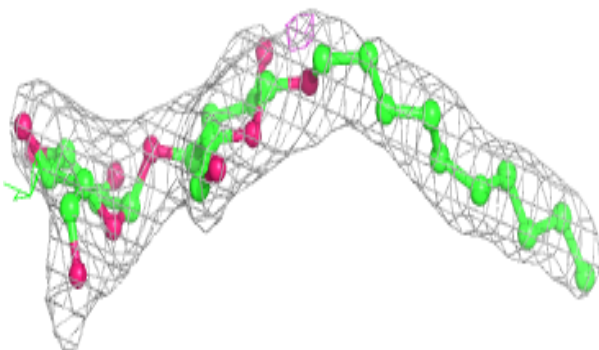
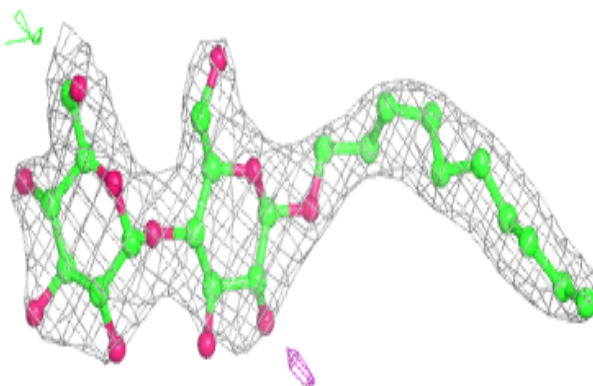
**Electron density around TGL A 606:**

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

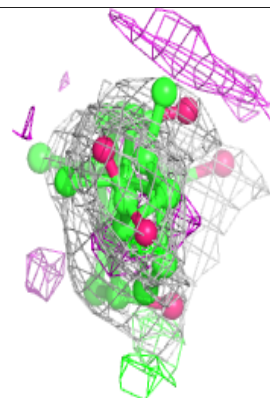
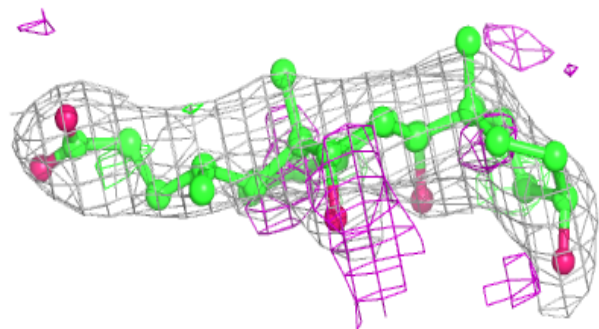
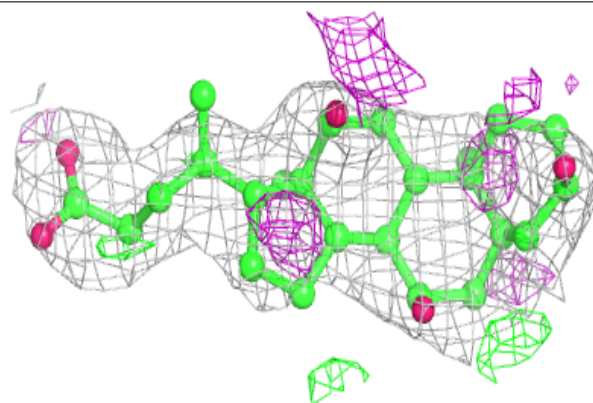


Electron density around DMU Z 101:

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

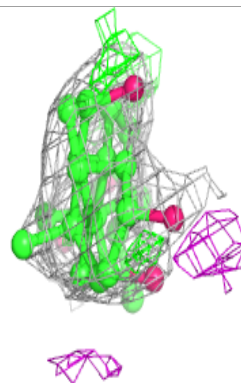
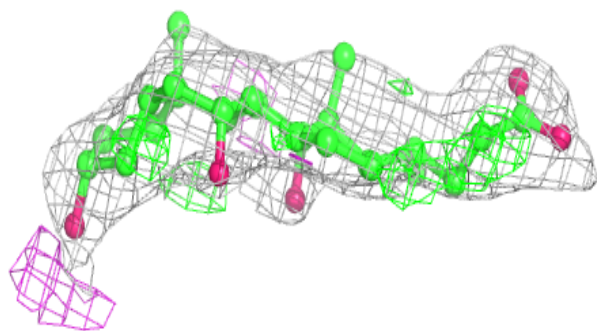
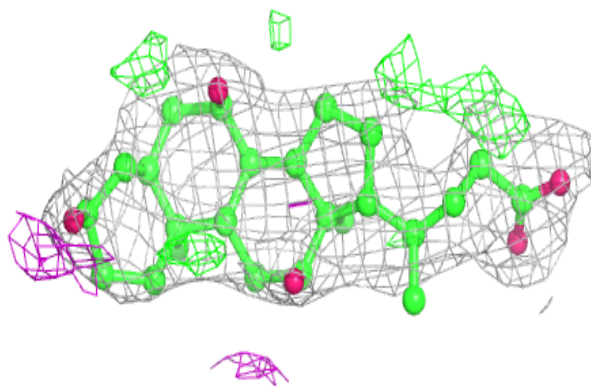
**Electron density around CHD C 308:**

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

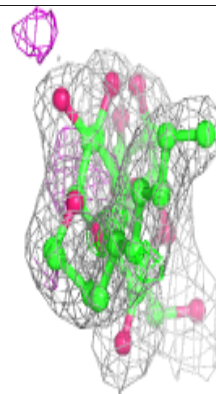
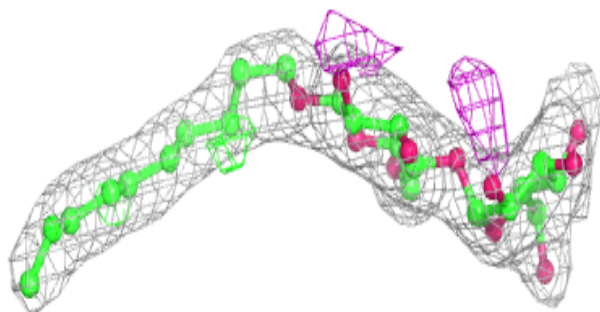
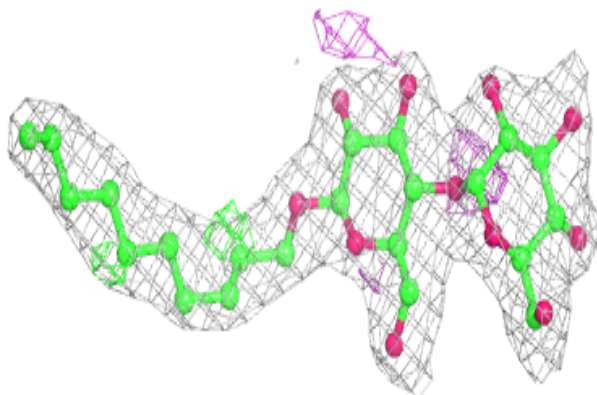


Electron density around CHD P 306:

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

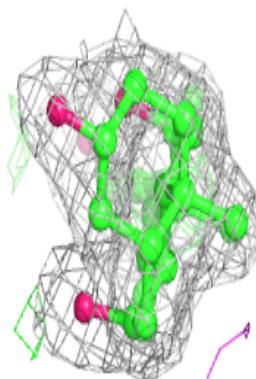
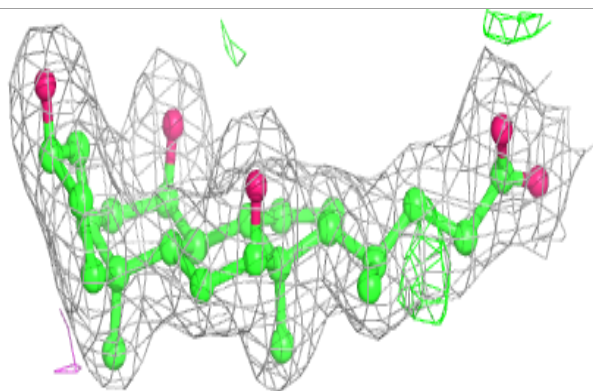
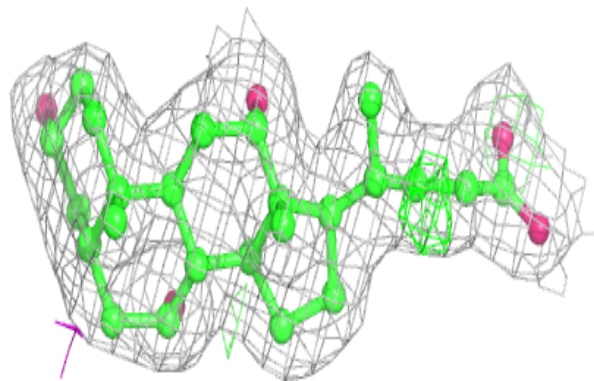
**Electron density around DMU M 102:**

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

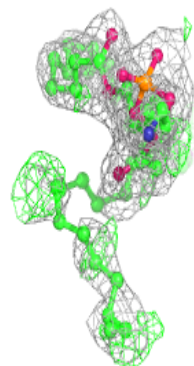
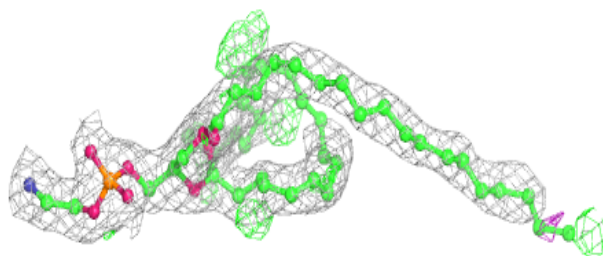
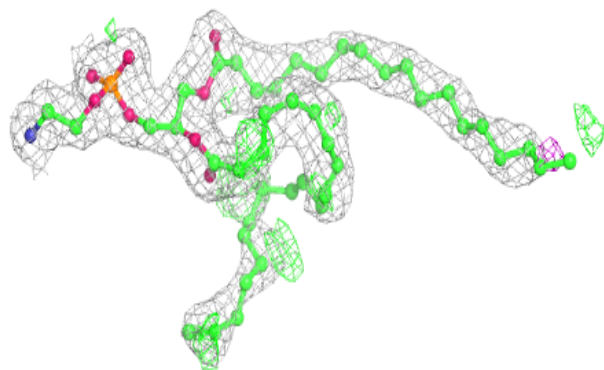


Electron density around CHD P 301:

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

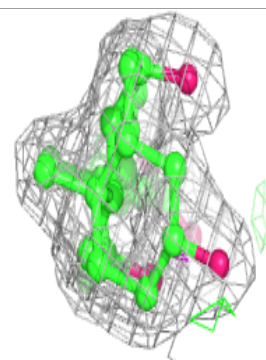
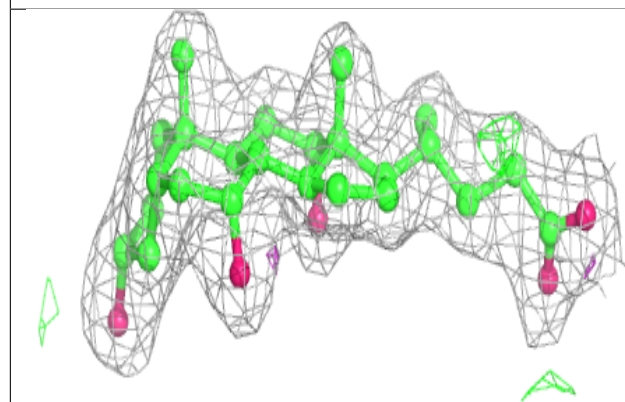
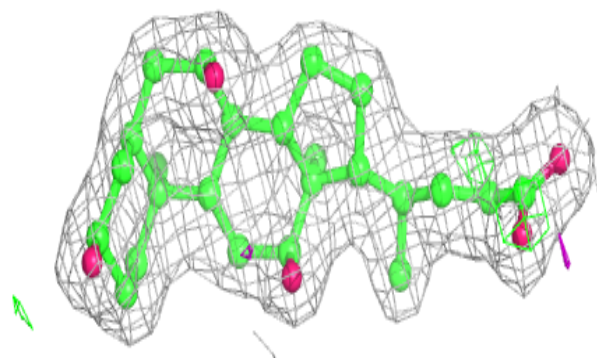
**Electron density around PEK T 102:**

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

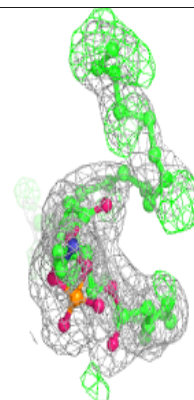
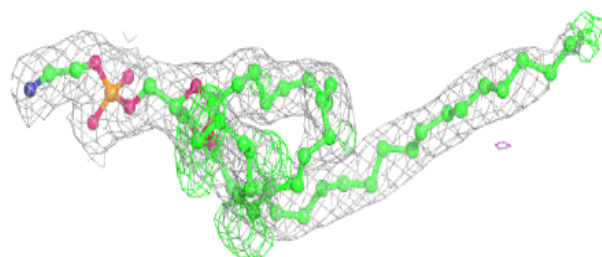
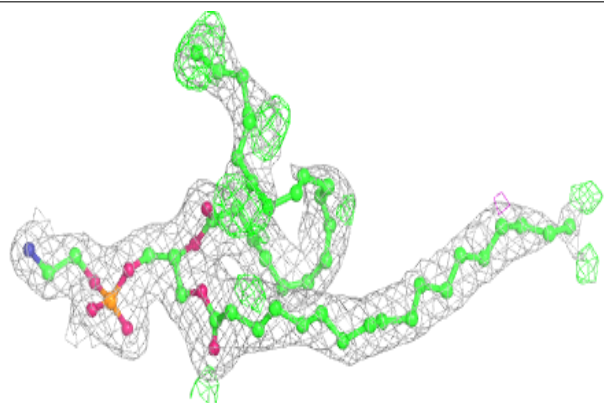


Electron density around CHD C 301:

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

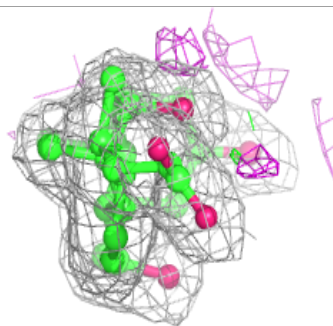
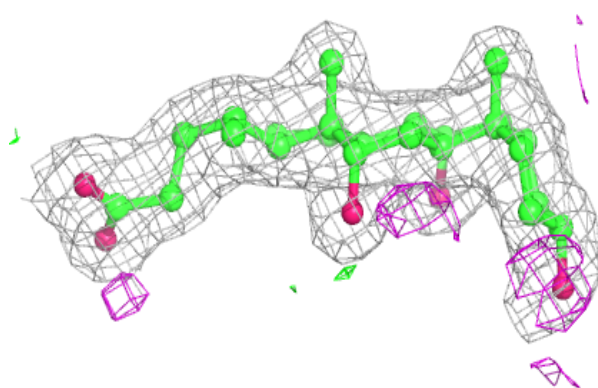
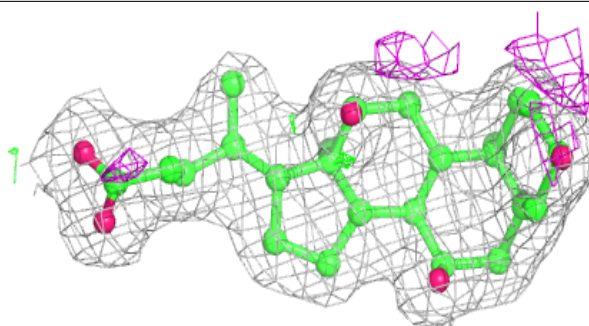
**Electron density around PEK C 303:**

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

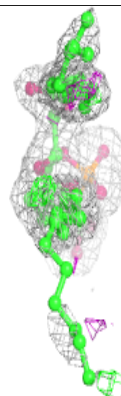
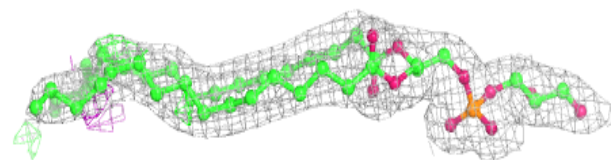
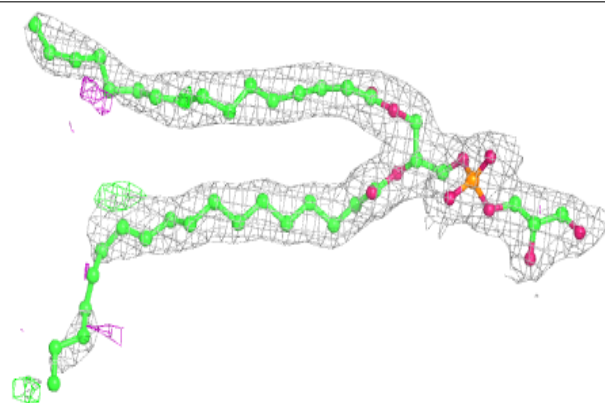


Electron density around CHD B 302:

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

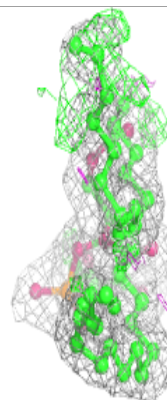
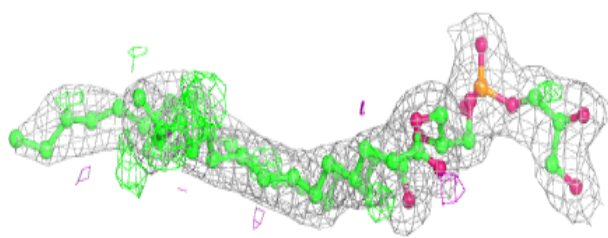
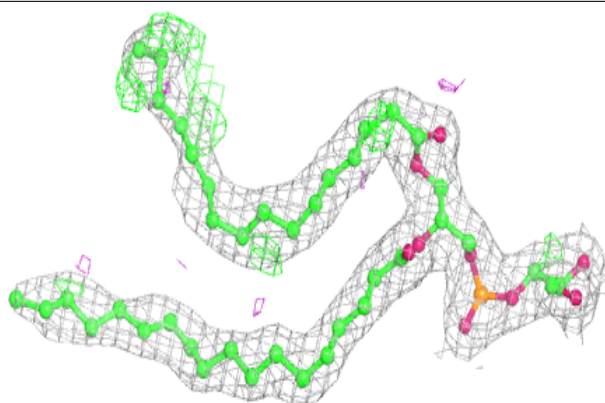
**Electron density around PGV P 303:**

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

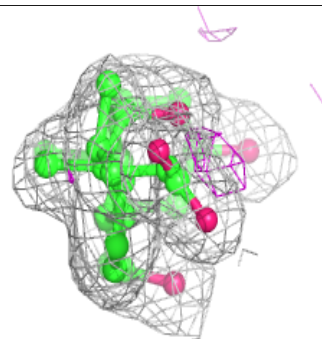
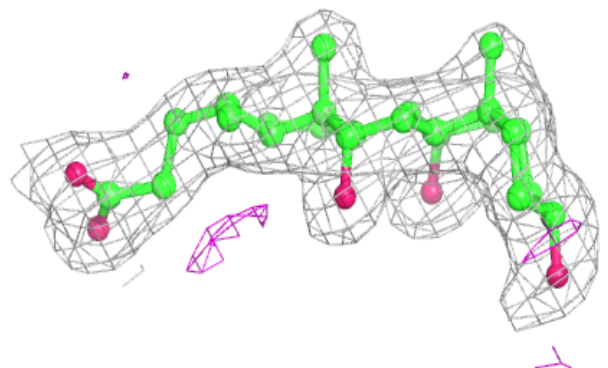
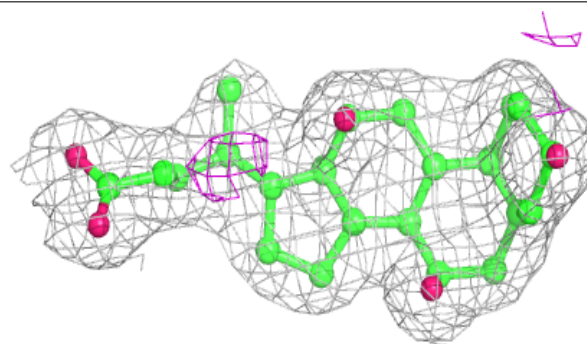


Electron density around PGV A 607:

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

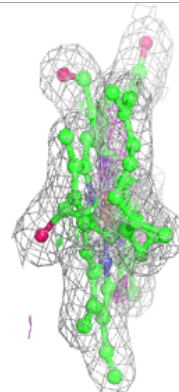
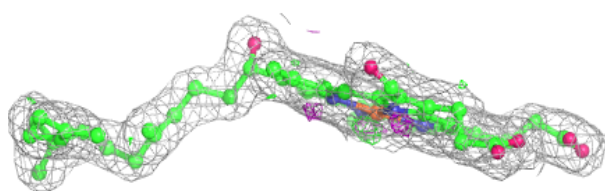
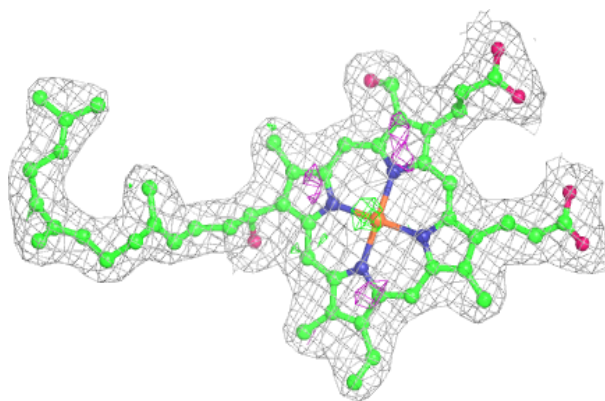
**Electron density around CHD O 301:**

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

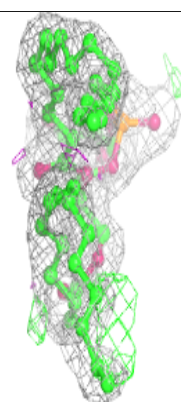
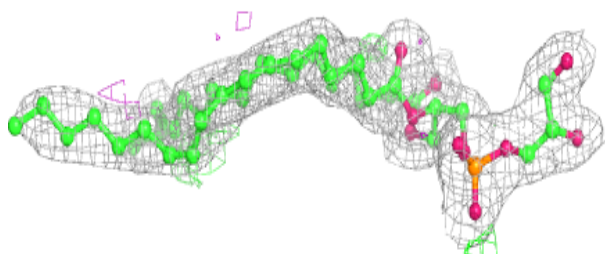
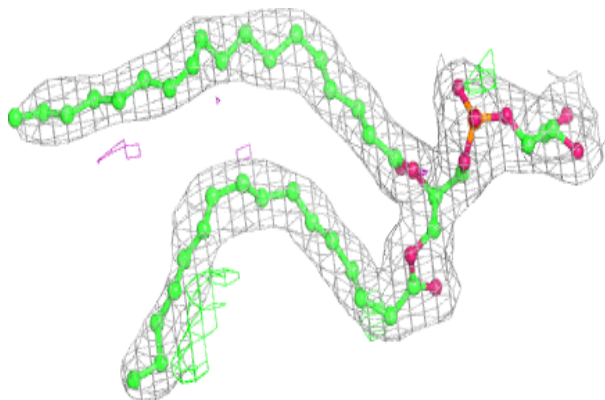


Electron density around HEA N 602:

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

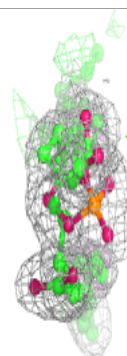
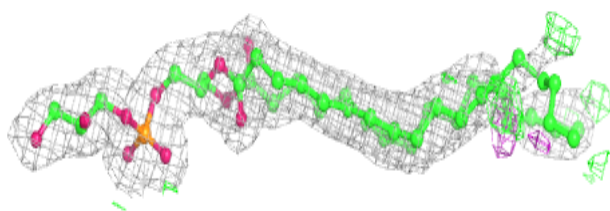
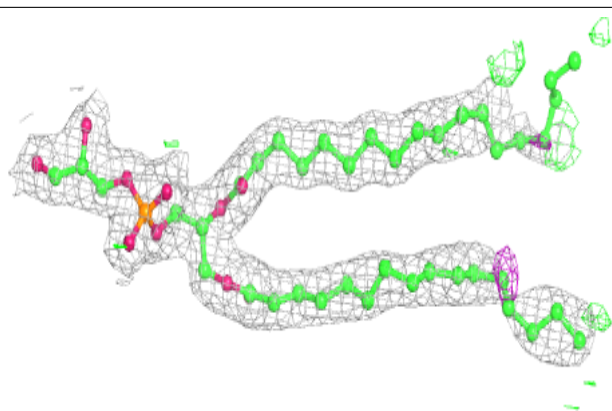
**Electron density around PGV N 607:**

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

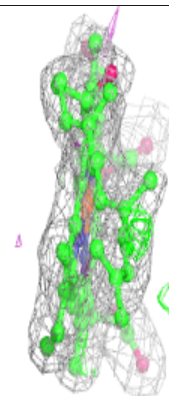
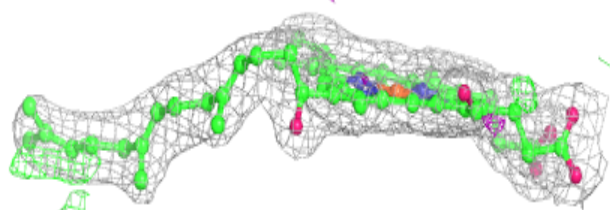
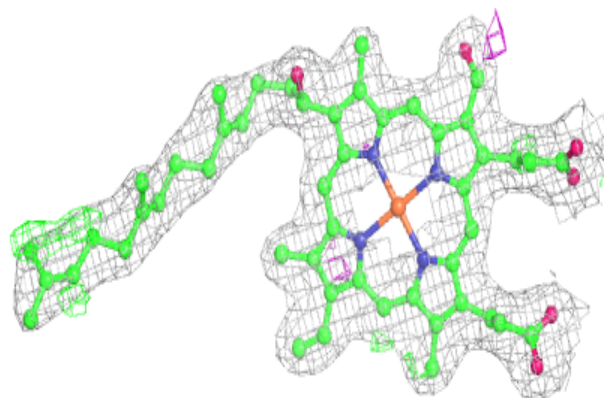


Electron density around PGV C 305:

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

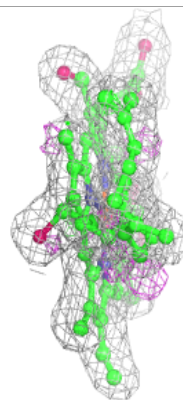
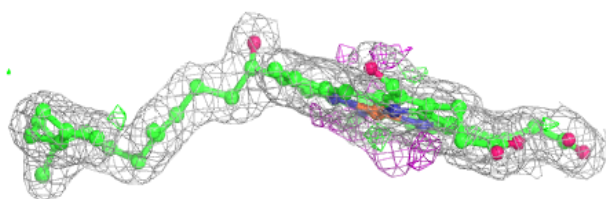
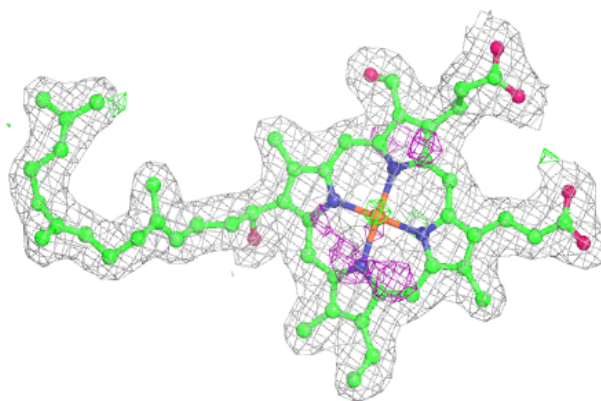
**Electron density around HEA N 601:**

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

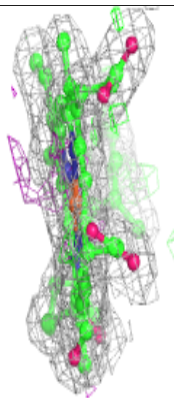
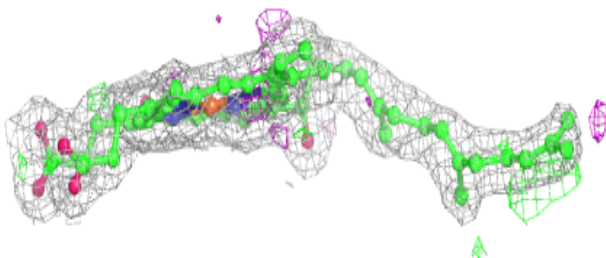
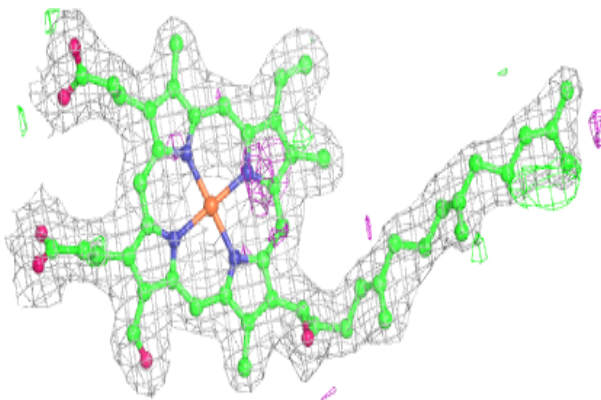


Electron density around HEA A 602:

$2mF_o-DF_c$ (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)



6.5 Other polymers [i](#)

There are no such residues in this entry.