



Full wwPDB EM Validation Report ⓘ

Mar 10, 2026 – 12:14 AM UTC

PDB ID : 5XNO / pdb_00005xno
EMDB ID : EMD-6744
Title : Structure of M-LHCII and CP24 complexes in the unstacked C2S2M2-type PSII-LHCII supercomplex from *Pisum sativum*
Authors : Su, X.D.; Ma, J.; Wei, X.P.; Cao, P.; Zhu, D.J.; Chang, W.R.; Liu, Z.F.; Zhang, X.Z.; Li, M.
Deposited on : 2017-05-23
Resolution : 3.50 Å(reported)

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

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A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

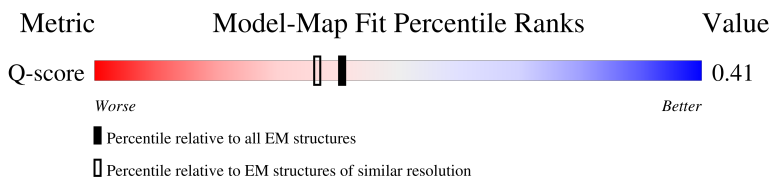
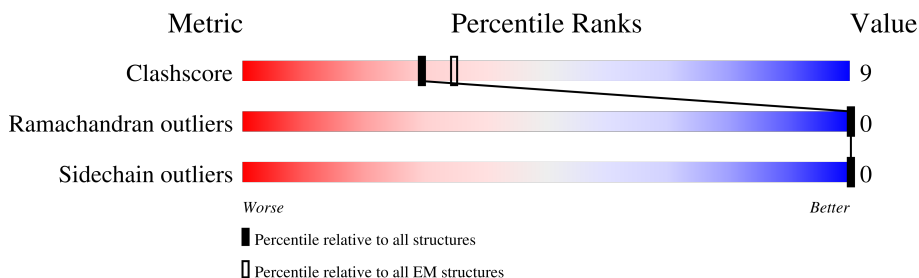
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	13950 (3.00 - 4.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	1	232	31% 78% 16% 6%
1	2	232	44% 77% 17% 6%
2	3	243	26% 75% 15% 9%
3	4	210	47% 76% 18% 6%

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
4	CHL	1	601	X	-	-	-
4	CHL	1	605	X	-	-	-
4	CHL	1	606	X	-	-	-
4	CHL	1	607	X	-	-	-
4	CHL	1	608	X	-	-	-
4	CHL	1	609	X	-	-	-
4	CHL	2	601	X	-	-	-
4	CHL	2	605	X	-	-	-
4	CHL	2	606	X	-	-	-
4	CHL	2	607	X	-	-	-
4	CHL	2	608	X	-	-	-
4	CHL	2	609	X	-	-	-
4	CHL	3	601	X	-	-	-
4	CHL	3	605	X	-	-	-
4	CHL	3	606	X	-	-	-
4	CHL	3	607	X	-	-	-
4	CHL	3	608	X	-	-	-
4	CHL	3	609	X	-	-	-
4	CHL	4	601	X	-	-	-
4	CHL	4	606	X	-	-	-
4	CHL	4	607	X	-	-	-
4	CHL	4	608	X	-	-	-
4	CHL	4	609	X	-	-	-
5	CLA	1	602	X	-	-	-
5	CLA	1	603	X	-	-	-
5	CLA	1	604	X	-	-	-
5	CLA	1	610	X	-	-	-
5	CLA	1	611	X	-	-	-
5	CLA	1	612	X	-	-	-
5	CLA	1	614	X	-	-	-
5	CLA	2	602	X	-	-	-
5	CLA	2	603	X	-	-	-
5	CLA	2	610	X	-	-	-
5	CLA	2	611	X	-	-	-
5	CLA	2	612	X	-	-	-
5	CLA	3	602	X	-	-	-
5	CLA	3	603	X	-	-	-
5	CLA	3	604	X	-	-	-
5	CLA	3	610	X	-	-	-
5	CLA	3	611	X	-	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	CLA	3	612	X	-	-	-
5	CLA	3	613	X	-	-	-
5	CLA	3	614	X	-	-	-
5	CLA	4	602	X	-	-	-
5	CLA	4	603	X	-	-	-
5	CLA	4	610	X	-	-	-
5	CLA	4	611	X	-	-	-
5	CLA	4	612	X	-	-	-

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 10020 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Chlorophyll a-b binding protein 8, chloroplastic.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	1	219	Total	C	N	O	S	0	0
			1668	1081	270	312	5		
1	2	218	Total	C	N	O	S	0	0
			1664	1079	269	311	5		

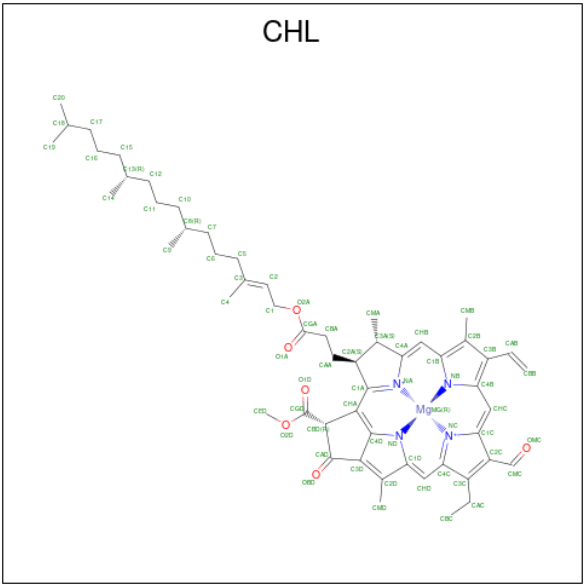
- Molecule 2 is a protein called Chlorophyll a-b binding protein, chloroplastic.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	3	220	Total	C	N	O	S	0	0
			1707	1116	277	309	5		

- Molecule 3 is a protein called Light harvesting chlorophyll a/b-binding protein Lhcb6, CP24.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	4	197	Total	C	N	O	S	0	0
			1534	1009	247	274	4		

- Molecule 4 is CHLOROPHYLL B (CCD ID: CHL) (formula: C₅₅H₇₀MgN₄O₆).



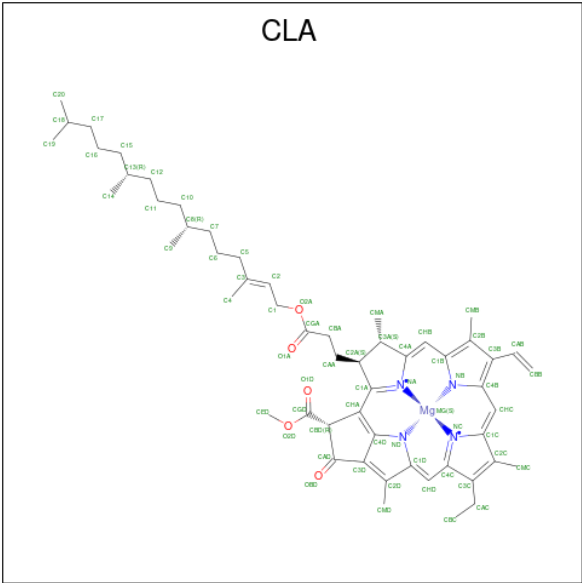
Mol	Chain	Residues	Atoms					AltConf
4	1	1	Total 46	C 35	Mg 1	N 4	O 6	0
4	1	1	Total 46	C 35	Mg 1	N 4	O 6	0
4	1	1	Total 46	C 35	Mg 1	N 4	O 6	0
4	1	1	Total 63	C 52	Mg 1	N 4	O 6	0
4	1	1	Total 46	C 35	Mg 1	N 4	O 6	0
4	1	1	Total 62	C 51	Mg 1	N 4	O 6	0
4	2	1	Total 46	C 35	Mg 1	N 4	O 6	0
4	2	1	Total 46	C 35	Mg 1	N 4	O 6	0
4	2	1	Total 46	C 35	Mg 1	N 4	O 6	0
4	2	1	Total 61	C 50	Mg 1	N 4	O 6	0
4	2	1	Total 46	C 35	Mg 1	N 4	O 6	0
4	2	1	Total 61	C 50	Mg 1	N 4	O 6	0
4	3	1	Total 64	C 53	Mg 1	N 4	O 6	0
4	3	1	Total 46	C 35	Mg 1	N 4	O 6	0
4	3	1	Total 46	C 35	Mg 1	N 4	O 6	0
4	3	1	Total 53	C 42	Mg 1	N 4	O 6	0
4	3	1	Total 46	C 35	Mg 1	N 4	O 6	0
4	3	1	Total 61	C 50	Mg 1	N 4	O 6	0
4	4	1	Total 45	C 34	Mg 1	N 4	O 6	0
4	4	1	Total 46	C 35	Mg 1	N 4	O 6	0
4	4	1	Total 46	C 35	Mg 1	N 4	O 6	0
4	4	1	Total 46	C 35	Mg 1	N 4	O 6	0

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Mol	Chain	Residues	Atoms					AltConf
4	4	1	Total	C	Mg	N	O	0
			46	35	1	4	6	

- Molecule 5 is CHLOROPHYLL A (CCD ID: CLA) (formula: C₅₅H₇₂MgN₄O₅).



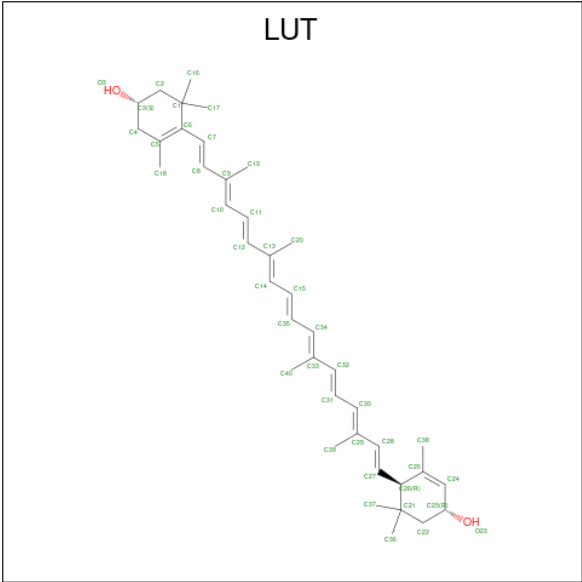
Mol	Chain	Residues	Atoms					AltConf
5	1	1	Total	C	Mg	N	O	0
			61	51	1	4	5	
5	1	1	Total	C	Mg	N	O	0
			55	45	1	4	5	
5	1	1	Total	C	Mg	N	O	0
			50	40	1	4	5	
5	1	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
5	1	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
5	1	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
5	1	1	Total	C	Mg	N	O	0
			55	45	1	4	5	
5	1	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
5	2	1	Total	C	Mg	N	O	0
			61	51	1	4	5	
5	2	1	Total	C	Mg	N	O	0
			55	45	1	4	5	

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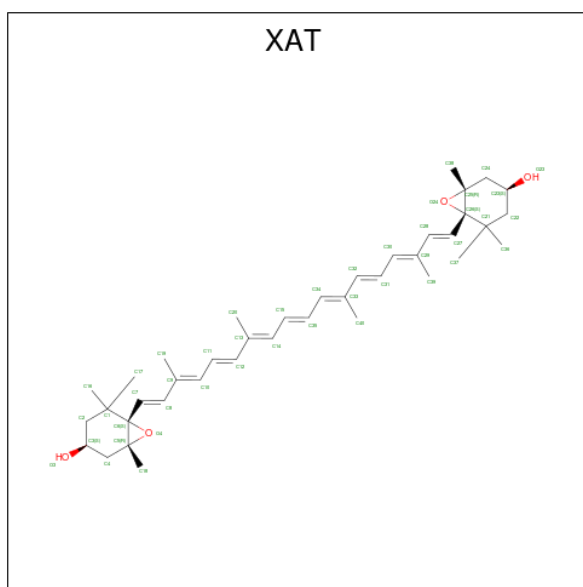
Mol	Chain	Residues	Atoms					AltConf
5	2	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
5	2	1	Total	C	Mg	N	O	0
			50	40	1	4	5	
5	2	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
5	2	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
5	2	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
5	2	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
5	3	1	Total	C	Mg	N	O	0
			60	50	1	4	5	
5	3	1	Total	C	Mg	N	O	0
			55	45	1	4	5	
5	3	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
5	3	1	Total	C	Mg	N	O	0
			60	50	1	4	5	
5	3	1	Total	C	Mg	N	O	0
			55	45	1	4	5	
5	3	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
5	3	1	Total	C	Mg	N	O	0
			58	48	1	4	5	
5	3	1	Total	C	Mg	N	O	0
			48	38	1	4	5	
5	4	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
5	4	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
5	4	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
5	4	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
5	4	1	Total	C	Mg	N	O	0
			45	35	1	4	5	

- Molecule 6 is (3R,3'R,6S)-4,5-DIDEHYDRO-5,6-DIHYDRO-BETA,BETA-CAROTENE-3, 3'-DIOL (CCD ID: LUT) (formula: C₄₀H₅₆O₂).



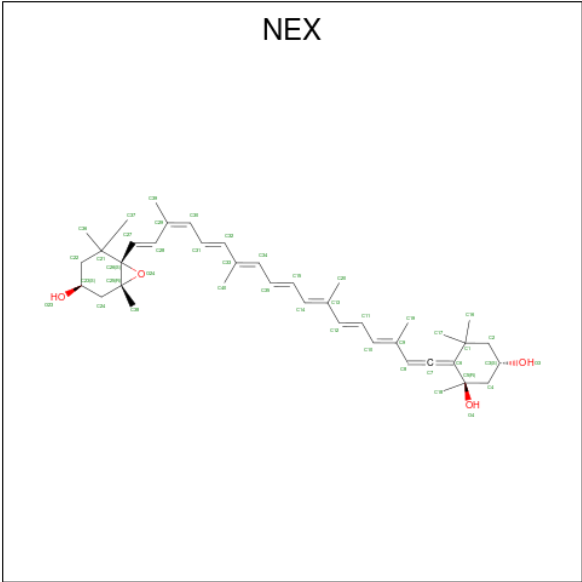
Mol	Chain	Residues	Atoms			AltConf
6	1	1	Total	C	O	0
			42	40	2	
6	1	1	Total	C	O	0
			42	40	2	
6	2	1	Total	C	O	0
			42	40	2	
6	2	1	Total	C	O	0
			42	40	2	
6	3	1	Total	C	O	0
			42	40	2	
6	3	1	Total	C	O	0
			42	40	2	
6	4	1	Total	C	O	0
			42	40	2	

- Molecule 7 is (3S,5R,6S,3'S,5'R,6'S)-5,6,5',6'-DIEPOXY-5,6,5',6'- TETRAHYDRO-BETA, BETA-CAROTENE-3,3'-DIOL (CCD ID: XAT) (formula: C₄₀H₅₆O₄).



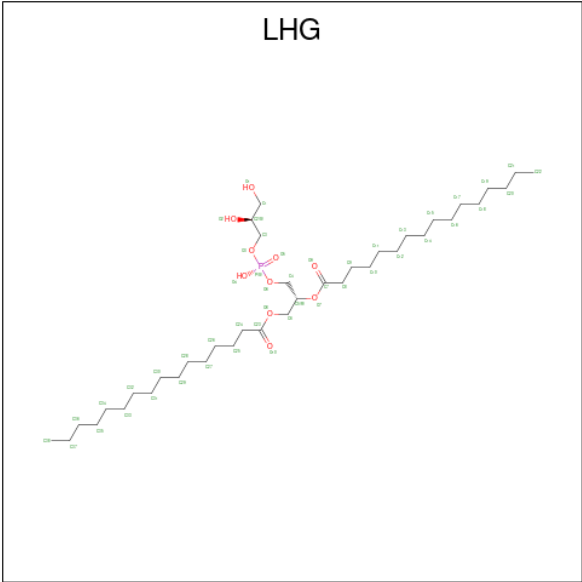
Mol	Chain	Residues	Atoms			AltConf
7	1	1	Total	C	O	0
			44	40	4	
7	2	1	Total	C	O	0
			44	40	4	
7	3	1	Total	C	O	0
			44	40	4	
7	4	1	Total	C	O	0
			44	40	4	

- Molecule 8 is (1R,3R)-6-[(3E,5E,7E,9E,11E,13E,15E,17E)-18-[(1S,4R,6R)-4-HYDROXY-2,2,6-TRIMETHYL-7-OXABICYCLO[4.1.0]HEPT-1-YL]-3,7,12,16-TETRAMETHYLOCTA DECA-1,3,5,7,9,11,13,15,17-NONAENYLIDENE}-1,5,5-TRIMETHYLCYCLOHEXANE-1,3-DIOL (CCD ID: NEX) (formula: C₄₀H₅₆O₄).



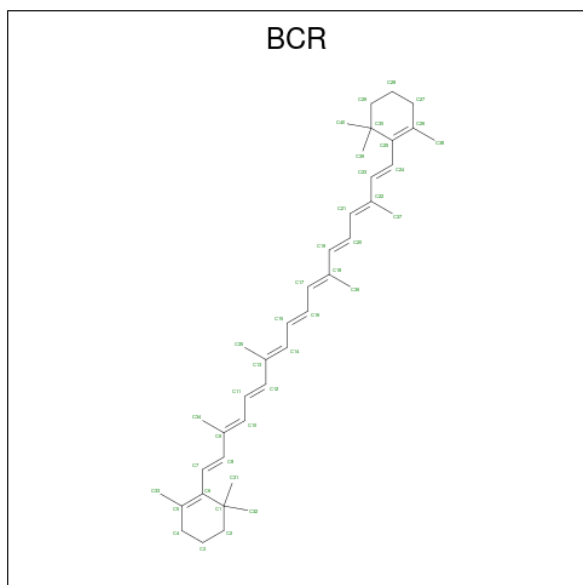
Mol	Chain	Residues	Atoms			AltConf
8	1	1	Total	C	O	0
			44	40	4	
8	2	1	Total	C	O	0
			44	40	4	
8	3	1	Total	C	O	0
			44	40	4	

- Molecule 9 is 1,2-DIPALMITOYL-PHOSPHATIDYL-GLYCEROLE (CCD ID: LHG) (formula: C₃₈H₇₅O₁₀P).



Mol	Chain	Residues	Atoms				AltConf
9	1	1	Total	C	O	P	0
			41	30	10	1	
9	2	1	Total	C	O	P	0
			37	26	10	1	
9	3	1	Total	C	O	P	0
			47	36	10	1	
9	4	1	Total	C	O	P	0
			21	10	10	1	

- Molecule 10 is BETA-CAROTENE (CCD ID: BCR) (formula: $C_{40}H_{56}$).

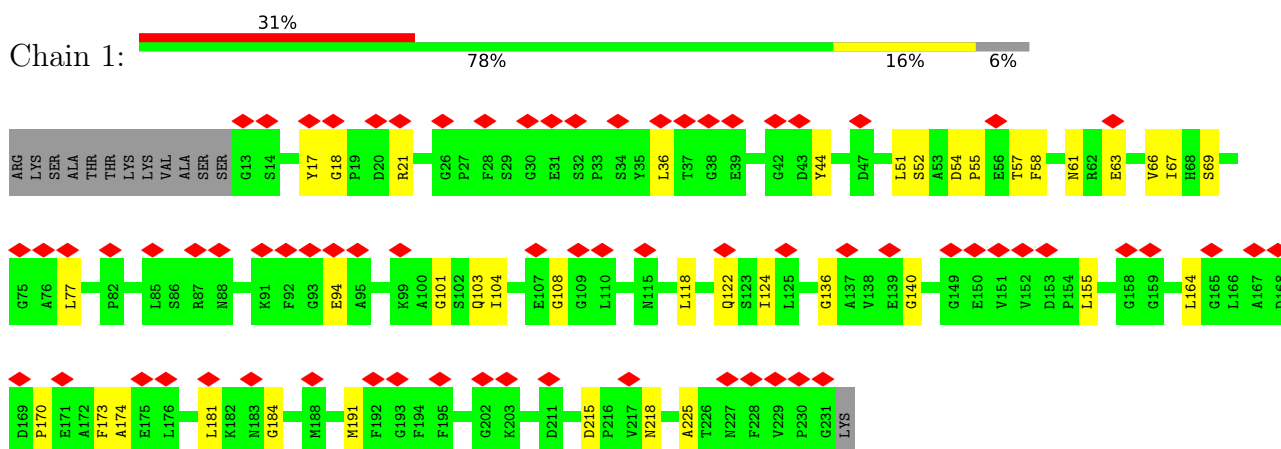


Mol	Chain	Residues	Atoms		AltConf
10	4	1	Total	C	0
			40	40	

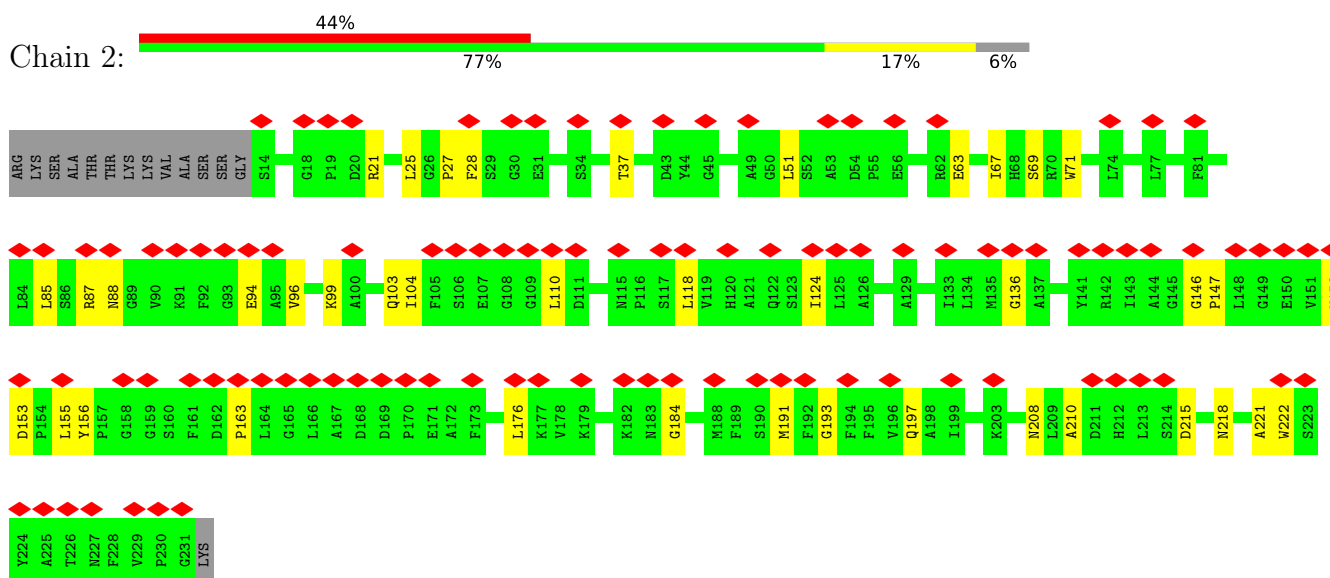
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

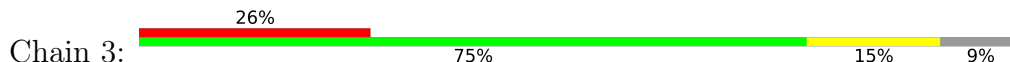
- Molecule 1: Chlorophyll a-b binding protein 8, chloroplastic

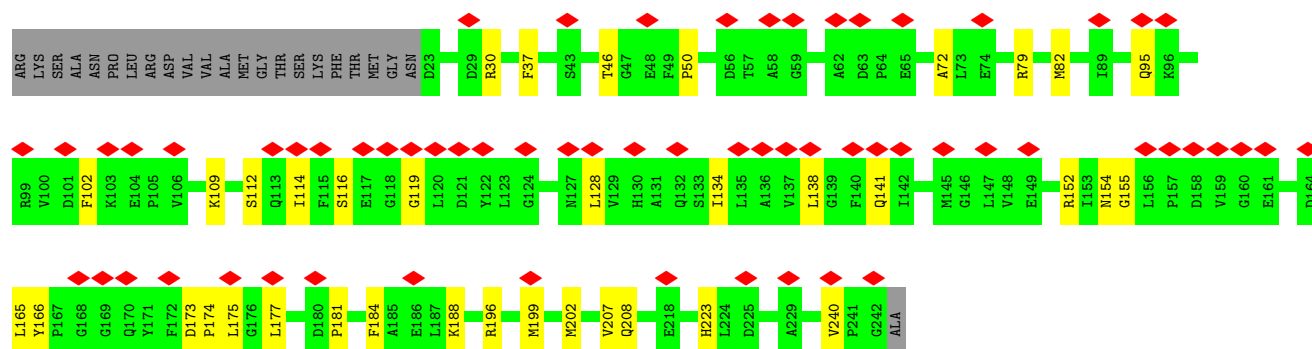


- Molecule 1: Chlorophyll a-b binding protein 8, chloroplastic

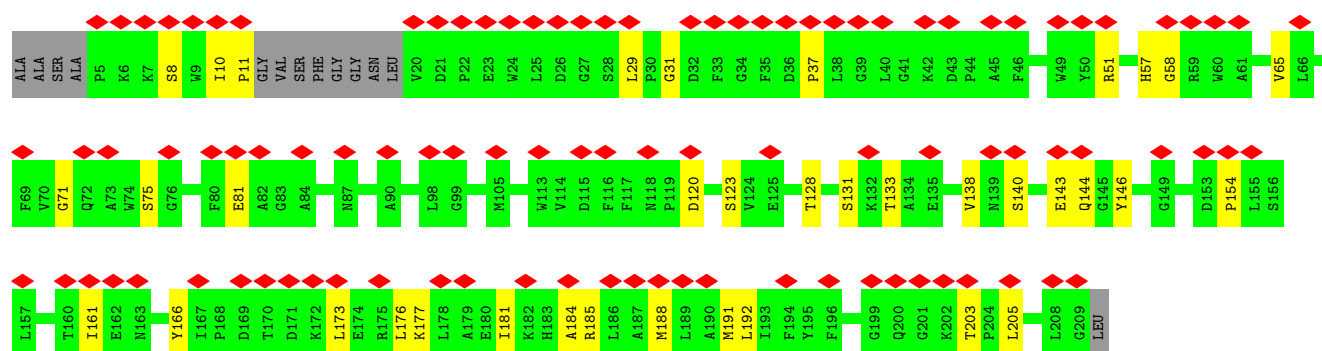
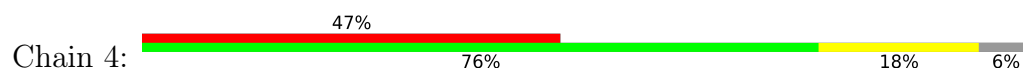


- Molecule 2: Chlorophyll a-b binding protein, chloroplastic





- Molecule 3: Light harvesting chlorophyll a/b-binding protein Lhcb6, CP24



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	50237	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.088	Depositor
Minimum map value	-0.039	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.007	Depositor
Recommended contour level	0.028	Depositor
Map size (Å)	156.0, 156.0, 156.0	wwPDB
Map dimensions	150, 150, 150	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.04, 1.04, 1.04	Depositor

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: CLA, LUT, LHG, CHL, XAT, BCR, NEX

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	1	0.42	0/1720	0.55	0/2342
1	2	0.36	0/1716	0.57	1/2337 (0.0%)
2	3	0.42	0/1759	0.58	0/2396
3	4	0.38	0/1586	0.65	0/2158
All	All	0.40	0/6781	0.59	1/9233 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	2	152	VAL	N-CA-C	-5.02	106.76	111.48

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	1	1668	0	1596	27	0
1	2	1664	0	1593	29	0
2	3	1707	0	1659	30	0
3	4	1534	0	1486	25	0
4	1	309	0	244	11	0
4	2	306	0	238	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	3	316	0	254	17	0
4	4	229	0	152	4	0
5	1	412	0	348	12	0
5	2	391	0	314	12	0
5	3	426	0	373	12	0
5	4	270	0	198	4	0
6	1	84	0	112	9	0
6	2	84	0	112	8	0
6	3	84	0	112	10	0
6	4	42	0	56	4	0
7	1	44	0	56	4	0
7	2	44	0	56	6	0
7	3	44	0	56	5	0
7	4	44	0	56	4	0
8	1	44	0	56	2	0
8	2	44	0	56	1	0
8	3	44	0	56	3	0
9	1	41	0	55	0	0
9	2	37	0	44	2	0
9	3	47	0	67	3	0
9	4	21	0	12	1	0
10	4	40	0	56	3	0
All	All	10020	0	9473	180	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 9.

All (180) 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)
4:3:606:CHL:HBB2	4:3:607:CHL:HBB1	1.57	0.86
5:1:602:CLA:HAB	6:1:1621:LUT:H32	1.66	0.77
1:2:191:MET:HE2	6:2:1621:LUT:H10	1.69	0.74
2:3:202:MET:HE2	6:3:1621:LUT:H10	1.69	0.73
5:2:610:CLA:H2	6:2:1620:LUT:H28	1.72	0.72
2:3:95:GLN:HE22	2:3:102:PHE:H	1.39	0.70
2:3:208:GLN:HE22	6:3:1620:LUT:H41	1.63	0.64
3:4:138:VAL:HG22	3:4:140:SER:H	1.64	0.61
5:1:610:CLA:H2	6:1:1620:LUT:H28	1.81	0.61
5:1:602:CLA:HBA1	6:1:1621:LUT:H382	1.83	0.61
1:2:163:PRO:HD2	6:2:1620:LUT:H23	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:103:GLN:HE22	5:1:604:CLA:HED3	1.67	0.59
1:2:104:ILE:HG21	1:2:124:ILE:HD13	1.83	0.59
7:3:1622:XAT:H12	9:3:2630:LHG:H191	1.83	0.59
2:3:50:PRO:HG3	2:3:188:LYS:HD3	1.83	0.59
1:2:21:ARG:NH1	1:2:37:THR:O	2.35	0.59
1:1:66:VAL:HG22	1:1:181:LEU:HD21	1.85	0.59
2:3:79:ARG:NH1	4:3:608:CHL:OBD	2.36	0.58
5:3:611:CLA:HAB	7:3:1622:XAT:H221	1.86	0.57
2:3:141:GLN:HE22	4:3:607:CHL:HMC	1.69	0.56
2:3:109:LYS:HA	4:3:607:CHL:HED3	1.87	0.56
1:1:21:ARG:NH2	1:1:36:LEU:O	2.39	0.55
1:1:94:GLU:N	1:1:103:GLN:OE1	2.38	0.55
1:2:27:PRO:O	2:3:154:ASN:ND2	2.40	0.55
1:2:193:GLY:O	1:2:197:GLN:HG2	2.07	0.55
1:2:215:ASP:OD2	1:2:218:ASN:ND2	2.41	0.55
2:3:175:LEU:HD13	6:3:1620:LUT:H222	1.88	0.54
4:1:607:CHL:HBB1	4:3:601:CHL:H141	1.89	0.54
3:4:161:ILE:HA	3:4:166:TYR:HA	1.90	0.54
1:1:164:LEU:HD13	6:1:1620:LUT:H222	1.90	0.54
1:2:85:LEU:HD23	1:2:88:ASN:HD22	1.73	0.53
3:4:31:GLY:HA3	3:4:181:ILE:HG21	1.90	0.53
1:1:191:MET:HE2	6:1:1621:LUT:H10	1.89	0.53
5:3:611:CLA:H3A	3:4:131:SER:HA	1.90	0.53
5:2:603:CLA:HED2	4:2:609:CHL:H93	1.90	0.53
7:2:1622:XAT:H14	9:2:2630:LHG:H171	1.91	0.53
2:3:174:PRO:HD2	6:3:1620:LUT:H23	1.91	0.52
4:1:608:CHL:HBB1	4:1:608:CHL:HHC	1.92	0.52
1:1:52:SER:OG	1:1:61:ASN:ND2	2.43	0.52
1:2:221:ALA:N	5:2:613:CLA:O1A	2.43	0.52
2:3:112:SER:HB3	4:3:607:CHL:HED2	1.90	0.52
1:2:25:LEU:HB3	1:2:28:PHE:HB2	1.91	0.52
1:2:63:GLU:HA	1:2:155:LEU:HD21	1.91	0.52
1:1:69:SER:HB3	1:1:184:GLY:HA3	1.92	0.52
3:4:65:VAL:HG21	6:4:620:LUT:H12	1.91	0.51
1:2:94:GLU:N	1:2:103:GLN:OE1	2.43	0.51
5:3:602:CLA:H72	6:3:1621:LUT:H30	1.93	0.51
3:4:120:ASP:O	3:4:123:SER:OG	2.28	0.51
2:3:114:ILE:HG21	2:3:134:ILE:HD13	1.91	0.51
1:1:51:LEU:HD13	5:1:602:CLA:H42	1.93	0.51
2:3:196:ARG:HA	2:3:199:MET:HE3	1.92	0.51
3:4:10:ILE:HG23	3:4:11:PRO:HD3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:4:58:GLY:HA3	3:4:184:ALA:HB1	1.93	0.51
5:2:602:CLA:HBA1	6:2:1621:LUT:H382	1.93	0.50
1:1:77:LEU:HD13	5:1:612:CLA:HBB2	1.92	0.50
7:2:1622:XAT:H41	4:3:607:CHL:HAA2	1.93	0.50
2:3:175:LEU:HB3	2:3:177:LEU:HD13	1.93	0.50
3:4:57:HIS:HD2	7:4:622:XAT:H15	1.77	0.50
5:4:612:CLA:HED2	5:4:612:CLA:H2A	1.94	0.50
2:3:152:ARG:NH2	4:3:609:CHL:O1D	2.45	0.50
3:4:154:PRO:HD2	6:4:620:LUT:H23	1.94	0.50
4:4:608:CHL:HHC	4:4:608:CHL:HBB1	1.93	0.50
1:2:208:ASN:ND2	5:2:613:CLA:O1D	2.45	0.49
4:1:607:CHL:HAA2	7:3:1622:XAT:H41	1.95	0.49
1:2:222:TRP:HZ2	2:3:138:LEU:HD22	1.78	0.49
3:4:71:GLY:O	3:4:75:SER:OG	2.30	0.49
1:2:103:GLN:HG3	1:2:110:LEU:HD13	1.96	0.48
1:2:176:LEU:HB3	5:2:610:CLA:H3A	1.95	0.48
3:4:191:MET:HE2	7:4:622:XAT:H10	1.94	0.48
7:3:1622:XAT:H393	9:3:2630:LHG:H101	1.94	0.48
3:4:143:GLU:H	3:4:146:TYR:HB2	1.79	0.48
1:2:96:VAL:HG12	1:2:99:LYS:H	1.78	0.48
1:2:51:LEU:HD13	5:2:602:CLA:H42	1.95	0.48
1:2:87:ARG:HH12	1:2:210:ALA:HB2	1.79	0.48
9:4:2630:LHG:O3	9:4:2630:LHG:O1	2.24	0.48
7:1:1622:XAT:H41	4:2:607:CHL:HAA2	1.95	0.48
4:3:608:CHL:H2A	4:3:608:CHL:HED3	1.96	0.47
3:4:37:PRO:HD2	7:4:622:XAT:H242	1.95	0.47
1:2:118:LEU:HD23	4:2:605:CHL:HED2	1.96	0.47
2:3:116:SER:OG	2:3:119:GLY:O	2.33	0.47
2:3:30:ARG:NH1	2:3:46:THR:O	2.48	0.47
5:1:613:CLA:H61	5:1:613:CLA:H2	1.60	0.47
1:2:69:SER:HB3	1:2:184:GLY:HA3	1.96	0.47
5:2:610:CLA:H43	5:2:612:CLA:HBA1	1.97	0.47
2:3:128:LEU:HD23	4:3:605:CHL:HED2	1.96	0.47
3:4:29:LEU:HD11	3:4:51:ARG:HD3	1.97	0.47
1:2:153:ASP:OD2	1:2:156:TYR:N	2.43	0.46
1:1:17:TYR:HE2	1:1:174:ALA:HB1	1.81	0.46
10:4:623:BCR:H351	10:4:623:BCR:H15C	1.76	0.46
7:1:1622:XAT:H31	7:1:1622:XAT:H391	1.74	0.46
8:1:1623:NEX:H11	8:1:1623:NEX:H191	1.78	0.46
7:2:1622:XAT:H31	7:2:1622:XAT:H391	1.71	0.46
2:3:72:ALA:HA	2:3:165:LEU:HD11	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:3:602:CLA:HBA1	6:3:1621:LUT:H382	1.98	0.46
5:3:603:CLA:HMD3	4:3:609:CHL:HBA2	1.97	0.46
5:4:611:CLA:HBA2	5:4:612:CLA:HMD3	1.98	0.46
1:1:108:GLY:O	1:1:122:GLN:NE2	2.39	0.46
1:2:147:PRO:HB2	4:2:608:CHL:HBB2	1.98	0.46
4:3:606:CHL:HBC2	4:3:607:CHL:HHD	1.97	0.46
4:3:609:CHL:H91	4:3:609:CHL:H112	1.83	0.46
8:2:1623:NEX:H15	8:2:1623:NEX:H201	1.71	0.46
4:3:606:CHL:HBB1	4:3:606:CHL:HHC	1.98	0.46
3:4:131:SER:O	3:4:133:THR:OG1	2.24	0.46
5:1:603:CLA:HMD3	4:1:609:CHL:HBA2	1.98	0.45
1:1:18:GLY:O	1:1:21:ARG:NH1	2.40	0.45
1:1:104:ILE:HG21	1:1:124:ILE:HD13	1.98	0.45
5:1:610:CLA:CBB	6:1:1620:LUT:H32	2.46	0.45
3:4:173:LEU:HG	3:4:177:LYS:HE3	1.98	0.45
8:3:1623:NEX:H201	8:3:1623:NEX:H15	1.67	0.45
3:4:185:ARG:HA	3:4:188:MET:HE3	1.98	0.45
4:3:606:CHL:HMB2	4:3:609:CHL:HMC	2.00	0.44
2:3:207:VAL:HG11	5:3:613:CLA:HAC2	1.99	0.44
4:2:606:CHL:HBC2	4:2:607:CHL:HHD	1.98	0.44
4:3:601:CHL:HAC1	9:3:2630:LHG:HC2	1.99	0.44
6:4:620:LUT:H15	6:4:620:LUT:H201	1.79	0.44
2:3:173:ASP:OD1	6:3:1620:LUT:O23	2.24	0.43
3:4:51:ARG:HH21	3:4:144:GLN:HB2	1.82	0.43
4:1:607:CHL:H91	4:1:607:CHL:H112	1.79	0.43
6:1:1620:LUT:H15	6:1:1620:LUT:H201	1.86	0.43
4:4:601:CHL:HBA1	4:4:601:CHL:H3A	1.76	0.43
1:2:136:GLY:HA2	4:2:609:CHL:HAB	2.00	0.43
1:2:146:GLY:HA3	1:2:147:PRO:HD3	1.85	0.43
4:4:606:CHL:HBA2	10:4:623:BCR:H19C	2.00	0.43
1:1:215:ASP:OD2	1:1:218:ASN:ND2	2.46	0.43
7:2:1622:XAT:H363	9:2:2630:LHG:HC41	2.01	0.43
5:2:603:CLA:H3A	5:2:603:CLA:HBA1	1.80	0.43
6:1:1621:LUT:H201	6:1:1621:LUT:H15	1.80	0.43
2:3:166:TYR:HB3	5:3:610:CLA:HED2	2.01	0.43
3:4:192:LEU:HD13	5:4:603:CLA:HBB2	2.01	0.43
4:1:606:CHL:HMB1	4:1:609:CHL:HMC	2.01	0.42
1:2:71:TRP:CE2	4:2:608:CHL:HED2	2.54	0.42
1:2:96:VAL:HB	1:2:99:LYS:HB2	2.01	0.42
7:2:1622:XAT:H35	7:2:1622:XAT:H401	1.88	0.42
3:4:8:SER:HB2	3:4:11:PRO:HD2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:225:ALA:HA	7:1:1622:XAT:H42	2.02	0.42
5:2:604:CLA:HBA1	4:2:606:CHL:C1D	2.49	0.42
10:4:623:BCR:H11C	10:4:623:BCR:H341	1.85	0.42
7:2:1622:XAT:H202	2:3:138:LEU:HD21	2.01	0.42
6:4:620:LUT:H401	6:4:620:LUT:H35	1.80	0.42
2:3:181:PRO:HA	2:3:184:PHE:HB3	2.02	0.42
1:1:63:GLU:HA	1:1:155:LEU:HD11	2.01	0.42
6:3:1620:LUT:H35	6:3:1620:LUT:H401	1.86	0.41
1:1:170:PRO:HA	1:1:173:PHE:HB3	2.00	0.41
7:4:622:XAT:H15	7:4:622:XAT:H201	1.82	0.41
1:1:57:THR:HG22	1:1:61:ASN:HD21	1.85	0.41
8:1:1623:NEX:H35	8:1:1623:NEX:H401	1.93	0.41
5:2:612:CLA:HBB1	6:2:1620:LUT:C13	2.51	0.41
1:1:44:TYR:N	5:1:602:CLA:OBD	2.40	0.41
1:1:54:ASP:HA	1:1:55:PRO:HD3	1.95	0.41
5:3:610:CLA:CBB	6:3:1620:LUT:H32	2.51	0.41
1:1:52:SER:HB3	1:1:58:PHE:HD1	1.85	0.41
5:2:602:CLA:CBB	6:2:1621:LUT:H32	2.50	0.41
5:3:613:CLA:H61	5:3:613:CLA:H2	1.81	0.41
8:3:1623:NEX:H11	8:3:1623:NEX:H191	1.80	0.41
8:3:1623:NEX:H35	8:3:1623:NEX:H401	1.87	0.41
5:1:612:CLA:HBB1	6:1:1620:LUT:H35	2.03	0.41
1:2:67:ILE:HG22	4:2:609:CHL:HMD2	2.02	0.41
2:3:79:ARG:HA	2:3:82:MET:HE3	2.03	0.41
2:3:223:HIS:CG	5:3:613:CLA:HAA2	2.56	0.41
3:4:128:THR:OG1	3:4:133:THR:N	2.53	0.41
1:1:118:LEU:HD23	1:1:118:LEU:HA	1.86	0.41
1:1:101:GLY:HA2	4:1:606:CHL:HAC2	2.03	0.41
5:1:603:CLA:H3A	5:1:603:CLA:HBA1	1.78	0.41
4:1:607:CHL:HED1	2:3:240:VAL:HG13	2.02	0.41
4:1:609:CHL:H42	5:3:602:CLA:H143	2.02	0.41
1:2:155:LEU:HA	1:2:155:LEU:HD23	1.89	0.41
6:2:1620:LUT:H15	6:2:1620:LUT:H201	1.85	0.41
3:4:176:LEU:HB3	5:4:610:CLA:H3A	2.02	0.41
4:2:606:CHL:HMB2	4:2:609:CHL:HMC	2.03	0.41
1:1:67:ILE:HG22	4:1:609:CHL:HMD2	2.03	0.40
3:4:203:THR:HG22	3:4:205:LEU:H	1.86	0.40
7:1:1622:XAT:H201	7:1:1622:XAT:H15	1.76	0.40
5:3:602:CLA:H93	5:3:602:CLA:H111	1.86	0.40
7:3:1622:XAT:H15	7:3:1622:XAT:H201	1.78	0.40
6:2:1621:LUT:H11	6:2:1621:LUT:H191	1.98	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:3:1620:LUT:H31	6:3:1620:LUT:H391	1.94	0.40
1:1:136:GLY:HA2	4:1:609:CHL:HAB	2.03	0.40
1:1:140:GLY:HA3	2:3:37:PHE:CD2	2.57	0.40
2:3:155:GLY:HA2	4:3:608:CHL:HAC1	2.04	0.40
3:4:81:GLU:HA	4:4:607:CHL:HED2	2.03	0.40

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 PDB entries followed by that with respect to all EM entries.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	1	217/232 (94%)	211 (97%)	6 (3%)	0	100	100
1	2	216/232 (93%)	211 (98%)	5 (2%)	0	100	100
2	3	218/243 (90%)	208 (95%)	10 (5%)	0	100	100
3	4	193/210 (92%)	179 (93%)	14 (7%)	0	100	100
All	All	844/917 (92%)	809 (96%)	35 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	1	171/182 (94%)	171 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	2	171/182 (94%)	171 (100%)	0	100	100
2	3	175/193 (91%)	175 (100%)	0	100	100
3	4	154/162 (95%)	154 (100%)	0	100	100
All	All	671/719 (93%)	671 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	1	61	ASN
1	2	61	ASN
1	2	88	ASN
2	3	95	GLN
2	3	219	ASN
3	4	72	GLN
3	4	122	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

72 ligands 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	LHG	3	2630	5	46,46,48	0.82	1 (2%)	49,52,54	1.33	4 (8%)
6	LUT	2	1621	-	42,43,43	0.80	1 (2%)	51,60,60	1.87	18 (35%)
9	LHG	1	2630	5	40,40,48	0.77	1 (2%)	43,46,54	1.35	6 (13%)
4	CHL	4	609	3	40,54,74	3.93	19 (47%)	34,90,114	2.93	14 (41%)
5	CLA	2	613	1	49,53,73	1.36	7 (14%)	58,89,113	1.27	6 (10%)
9	LHG	4	2630	5	20,20,48	0.88	1 (5%)	23,26,54	1.36	2 (8%)
4	CHL	3	606	-	40,54,74	4.06	20 (50%)	34,90,114	3.02	13 (38%)
4	CHL	1	606	-	40,54,74	3.87	20 (50%)	34,90,114	3.00	11 (32%)
4	CHL	4	601	3	39,53,74	4.32	18 (46%)	33,89,114	2.10	12 (36%)
5	CLA	2	614	1	49,53,73	1.35	9 (18%)	58,89,113	1.29	7 (12%)
5	CLA	1	613	1	59,63,73	1.25	9 (15%)	70,101,113	1.29	8 (11%)
6	LUT	1	1620	-	42,43,43	0.80	0	51,60,60	1.99	14 (27%)
5	CLA	2	612	1	49,53,73	1.37	9 (18%)	58,89,113	1.36	8 (13%)
7	XAT	3	1622	-	41,47,47	1.02	2 (4%)	54,74,74	2.96	26 (48%)
5	CLA	2	602	1	65,69,73	1.20	9 (13%)	77,108,113	1.21	8 (10%)
4	CHL	2	608	-	40,54,74	4.02	20 (50%)	34,90,114	2.80	10 (29%)
5	CLA	2	611	9	49,53,73	1.34	8 (16%)	58,89,113	1.42	6 (10%)
4	CHL	2	607	-	55,69,74	3.47	20 (36%)	52,108,114	1.95	14 (26%)
5	CLA	3	602	2	64,68,73	1.20	9 (14%)	76,107,113	1.20	7 (9%)
6	LUT	2	1620	-	42,43,43	0.73	0	51,60,60	1.88	14 (27%)
5	CLA	1	604	-	54,58,73	1.31	9 (16%)	64,95,113	1.36	7 (10%)
5	CLA	2	610	1	54,58,73	1.28	7 (12%)	64,95,113	1.14	4 (6%)
5	CLA	4	604	-	49,53,73	1.37	9 (18%)	58,89,113	1.37	6 (10%)
5	CLA	1	611	9	49,53,73	1.35	10 (20%)	58,89,113	1.43	4 (6%)
5	CLA	4	602	3	49,53,73	1.30	9 (18%)	58,89,113	1.46	8 (13%)
7	XAT	4	622	-	41,47,47	0.93	2 (4%)	54,74,74	2.63	21 (38%)
4	CHL	3	605	2	40,54,74	4.05	19 (47%)	34,90,114	2.84	13 (38%)
6	LUT	1	1621	-	42,43,43	0.92	2 (4%)	51,60,60	1.98	15 (29%)
5	CLA	3	612	2	49,53,73	1.39	10 (20%)	58,89,113	1.39	6 (10%)
6	LUT	4	620	-	42,43,43	0.87	2 (4%)	51,60,60	2.28	17 (33%)
5	CLA	4	603	3	49,53,73	1.37	9 (18%)	58,89,113	1.38	5 (8%)
8	NEX	3	1623	-	40,46,46	0.99	2 (5%)	50,70,70	2.32	16 (32%)
5	CLA	3	614	2	52,56,73	1.34	7 (13%)	61,92,113	1.31	6 (9%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	XAT	2	1622	-	41,47,47	0.97	2 (4%)	54,74,74	2.78	22 (40%)
4	CHL	1	605	1	40,54,74	3.98	21 (52%)	34,90,114	2.67	10 (29%)
8	NEX	2	1623	-	40,46,46	1.09	3 (7%)	50,70,70	2.27	14 (28%)
4	CHL	1	609	1	56,70,74	3.37	20 (35%)	53,109,114	2.29	14 (26%)
4	CHL	3	601	2	58,72,74	3.21	20 (34%)	55,111,114	2.72	18 (32%)
4	CHL	1	601	1	40,54,74	3.94	20 (50%)	34,90,114	2.84	14 (41%)
5	CLA	1	614	1	49,53,73	1.36	9 (18%)	58,89,113	1.31	7 (12%)
5	CLA	4	612	3	49,53,73	1.36	9 (18%)	58,89,113	1.39	7 (12%)
8	NEX	1	1623	-	40,46,46	1.24	4 (10%)	50,70,70	2.49	18 (36%)
4	CHL	3	608	-	40,54,74	3.94	20 (50%)	34,90,114	2.85	13 (38%)
5	CLA	3	613	2	62,66,73	1.22	9 (14%)	73,104,113	1.28	4 (5%)
4	CHL	3	607	-	47,61,74	3.71	19 (40%)	41,98,114	2.43	15 (36%)
4	CHL	2	609	1	55,69,74	3.37	20 (36%)	52,108,114	2.51	17 (32%)
4	CHL	4	607	-	40,54,74	4.04	19 (47%)	34,90,114	2.58	14 (41%)
9	LHG	2	2630	5	36,36,48	0.75	1 (2%)	39,42,54	1.29	4 (10%)
5	CLA	3	611	9	59,63,73	1.30	10 (16%)	70,101,113	1.34	8 (11%)
10	BCR	4	623	-	41,41,41	0.72	0	56,56,56	2.20	15 (26%)
4	CHL	1	608	-	40,54,74	3.95	20 (50%)	34,90,114	2.93	12 (35%)
5	CLA	1	612	1	49,53,73	1.39	9 (18%)	58,89,113	1.49	9 (15%)
4	CHL	1	607	-	57,71,74	3.37	20 (35%)	54,110,114	2.54	19 (35%)
5	CLA	3	603	2	59,63,73	1.24	11 (18%)	70,101,113	1.44	8 (11%)
5	CLA	3	604	-	49,53,73	1.36	8 (16%)	58,89,113	1.28	4 (6%)
5	CLA	1	610	1	60,64,73	1.19	6 (10%)	71,102,113	1.19	9 (12%)
4	CHL	4	608	-	40,54,74	3.96	20 (50%)	34,90,114	3.06	15 (44%)
4	CHL	4	606	-	40,54,74	4.07	19 (47%)	34,90,114	2.95	13 (38%)
4	CHL	2	606	-	40,54,74	3.95	19 (47%)	34,90,114	2.89	13 (38%)
5	CLA	1	603	1	59,63,73	1.27	10 (16%)	70,101,113	1.42	7 (10%)
5	CLA	4	610	3	49,53,73	1.36	9 (18%)	58,89,113	1.24	7 (12%)
4	CHL	3	609	2	55,69,74	3.46	20 (36%)	52,108,114	2.30	15 (28%)
5	CLA	3	610	2	64,68,73	1.21	7 (10%)	76,107,113	1.15	7 (9%)
6	LUT	3	1620	-	42,43,43	0.86	2 (4%)	51,60,60	2.20	21 (41%)
4	CHL	2	601	1	40,54,74	4.06	20 (50%)	34,90,114	2.85	16 (47%)
5	CLA	2	603	1	59,63,73	1.26	9 (15%)	70,101,113	1.37	6 (8%)
5	CLA	1	602	1	65,69,73	1.23	10 (15%)	77,108,113	1.18	7 (9%)
5	CLA	2	604	-	49,53,73	1.36	9 (18%)	58,89,113	1.41	10 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	XAT	1	1622	-	41,47,47	0.93	2 (4%)	54,74,74	2.64	20 (37%)
6	LUT	3	1621	-	42,43,43	0.86	1 (2%)	51,60,60	1.70	12 (23%)
4	CHL	2	605	1	40,54,74	4.11	19 (47%)	34,90,114	2.59	12 (35%)
5	CLA	4	611	9	49,53,73	1.34	8 (16%)	58,89,113	1.26	6 (10%)

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	LHG	3	2630	5	-	23/51/51/53	-
6	LUT	2	1621	-	-	0/29/67/67	0/2/2/2
9	LHG	1	2630	5	-	9/45/45/53	-
4	CHL	4	609	3	3/3/16/26	6/15/113/137	-
5	CLA	2	613	1	-	5/15/91/115	-
9	LHG	4	2630	5	-	8/23/23/53	-
4	CHL	3	606	-	3/3/16/26	5/15/113/137	-
4	CHL	1	606	-	3/3/16/26	7/15/113/137	-
4	CHL	4	601	3	3/3/16/26	3/13/111/137	-
5	CLA	2	614	1	-	8/15/91/115	-
5	CLA	1	613	1	-	7/27/103/115	-
6	LUT	1	1620	-	-	2/29/67/67	0/2/2/2
5	CLA	2	612	1	1/1/11/20	8/15/91/115	-
7	XAT	3	1622	-	-	1/31/93/93	0/4/4/4
5	CLA	2	602	1	1/1/14/20	13/35/111/115	-
4	CHL	2	608	-	3/3/16/26	7/15/113/137	-
5	CLA	2	611	9	1/1/11/20	4/15/91/115	-
4	CHL	2	607	-	4/4/19/26	20/33/131/137	-
5	CLA	3	602	2	1/1/14/20	15/33/109/115	-
6	LUT	2	1620	-	-	2/29/67/67	0/2/2/2
5	CLA	1	604	-	1/1/12/20	10/21/97/115	-
5	CLA	2	610	1	1/1/12/20	5/21/97/115	-
5	CLA	4	604	-	-	9/15/91/115	-
5	CLA	1	611	9	1/1/11/20	7/15/91/115	-
5	CLA	4	602	3	1/1/11/20	4/15/91/115	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	XAT	4	622	-	-	2/31/93/93	0/4/4/4
4	CHL	3	605	2	3/3/16/26	9/15/113/137	-
6	LUT	1	1621	-	-	5/29/67/67	0/2/2/2
5	CLA	3	612	2	1/1/11/20	6/15/91/115	-
6	LUT	4	620	-	-	5/29/67/67	0/2/2/2
5	CLA	4	603	3	1/1/11/20	6/15/91/115	-
8	NEX	3	1623	-	-	4/27/83/83	0/3/3/3
5	CLA	3	614	2	1/1/11/20	5/19/95/115	-
7	XAT	2	1622	-	-	4/31/93/93	0/4/4/4
4	CHL	1	605	1	3/3/16/26	8/15/113/137	-
8	NEX	2	1623	-	-	4/27/83/83	0/3/3/3
4	CHL	1	609	1	4/4/19/26	13/35/133/137	-
4	CHL	3	601	2	4/4/19/26	12/37/135/137	-
4	CHL	1	601	1	3/3/16/26	6/15/113/137	-
5	CLA	1	614	1	1/1/11/20	5/15/91/115	-
5	CLA	4	612	3	1/1/11/20	8/15/91/115	-
8	NEX	1	1623	-	-	8/27/83/83	0/3/3/3
4	CHL	3	608	-	3/3/16/26	8/15/113/137	-
5	CLA	3	613	2	1/1/13/20	8/31/107/115	-
4	CHL	3	607	-	3/3/17/26	13/24/122/137	-
4	CHL	2	609	1	4/4/19/26	13/33/131/137	-
4	CHL	4	607	-	3/3/16/26	8/15/113/137	-
9	LHG	2	2630	5	-	14/41/41/53	-
5	CLA	3	611	9	1/1/13/20	11/27/103/115	-
10	BCR	4	623	-	-	12/29/63/63	0/2/2/2
4	CHL	1	608	-	3/3/16/26	3/15/113/137	-
5	CLA	1	612	1	1/1/11/20	6/15/91/115	-
4	CHL	1	607	-	4/4/19/26	20/36/134/137	-
5	CLA	3	603	2	1/1/13/20	10/27/103/115	-
5	CLA	3	604	-	1/1/11/20	10/15/91/115	-
5	CLA	1	610	1	1/1/13/20	4/29/105/115	-
4	CHL	4	608	-	3/3/16/26	10/15/113/137	-
4	CHL	4	606	-	3/3/16/26	6/15/113/137	-
4	CHL	2	606	-	3/3/16/26	5/15/113/137	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	CLA	1	603	1	1/1/13/20	11/27/103/115	-
5	CLA	4	610	3	1/1/11/20	6/15/91/115	-
4	CHL	3	609	2	4/4/19/26	14/33/131/137	-
5	CLA	3	610	2	1/1/14/20	6/33/109/115	-
6	LUT	3	1620	-	-	1/29/67/67	0/2/2/2
4	CHL	2	601	1	3/3/16/26	6/15/113/137	-
5	CLA	2	603	1	1/1/13/20	14/27/103/115	-
5	CLA	1	602	1	1/1/14/20	10/35/111/115	-
5	CLA	2	604	-	-	7/15/91/115	-
7	XAT	1	1622	-	-	2/31/93/93	0/4/4/4
6	LUT	3	1621	-	-	3/29/67/67	0/2/2/2
4	CHL	2	605	1	3/3/16/26	8/15/113/137	-
5	CLA	4	611	9	1/1/11/20	8/15/91/115	-

All (744) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	4	601	CHL	C2C-C3C	10.79	1.46	1.36
4	2	605	CHL	C2C-C3C	10.53	1.46	1.36
4	3	608	CHL	C2C-C3C	10.39	1.46	1.36
4	2	608	CHL	C2C-C3C	10.39	1.46	1.36
4	2	606	CHL	C2C-C3C	10.31	1.46	1.36
4	3	609	CHL	C2C-C3C	10.31	1.46	1.36
4	4	606	CHL	C2C-C3C	10.31	1.46	1.36
4	2	607	CHL	C2C-C3C	10.23	1.46	1.36
4	3	605	CHL	C2C-C3C	10.17	1.46	1.36
4	1	605	CHL	C2C-C3C	10.01	1.46	1.36
4	1	609	CHL	C2C-C3C	10.00	1.46	1.36
4	1	606	CHL	C2C-C3C	9.95	1.46	1.36
4	2	601	CHL	C2C-C3C	9.81	1.45	1.36
4	4	601	CHL	C1D-C2D	9.78	1.50	1.39
4	4	607	CHL	C2C-C3C	9.76	1.45	1.36
4	1	608	CHL	C2C-C3C	9.70	1.45	1.36
4	1	607	CHL	C2C-C3C	9.68	1.45	1.36
4	3	606	CHL	C2C-C3C	9.61	1.45	1.36
4	1	607	CHL	C1D-C2D	9.60	1.50	1.39
4	2	601	CHL	C1D-C2D	9.56	1.50	1.39
4	1	609	CHL	C1D-C2D	9.50	1.50	1.39
4	4	609	CHL	C2C-C3C	9.48	1.45	1.36
4	4	607	CHL	C1D-C2D	9.46	1.50	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	2	609	CHL	C1D-C2D	9.41	1.50	1.39
4	3	605	CHL	C1D-C2D	9.39	1.50	1.39
4	3	609	CHL	C1D-C2D	9.39	1.50	1.39
4	1	601	CHL	C2C-C3C	9.35	1.45	1.36
4	1	601	CHL	C1D-C2D	9.32	1.50	1.39
4	2	609	CHL	C2C-C3C	9.30	1.45	1.36
4	4	606	CHL	C1D-C2D	9.28	1.50	1.39
4	4	609	CHL	C1D-C2D	9.28	1.50	1.39
4	3	607	CHL	C1D-C2D	9.24	1.50	1.39
4	2	608	CHL	C1D-C2D	9.22	1.50	1.39
4	4	601	CHL	C1A-CHA	9.21	1.50	1.40
4	2	605	CHL	C1D-C2D	9.20	1.50	1.39
4	3	606	CHL	C1D-C2D	9.19	1.50	1.39
4	1	605	CHL	C1D-C2D	9.14	1.50	1.39
4	3	607	CHL	C2C-C3C	9.12	1.45	1.36
4	3	601	CHL	C1D-C2D	9.06	1.49	1.39
4	2	607	CHL	C1D-C2D	8.98	1.49	1.39
4	4	608	CHL	C2C-C3C	8.87	1.45	1.36
4	4	608	CHL	C1D-C2D	8.86	1.49	1.39
4	2	606	CHL	C1D-C2D	8.85	1.49	1.39
4	3	608	CHL	C1D-C2D	8.84	1.49	1.39
4	1	608	CHL	C1D-C2D	8.82	1.49	1.39
4	3	601	CHL	C2C-C3C	8.79	1.44	1.36
4	1	606	CHL	C1D-C2D	8.52	1.49	1.39
4	4	601	CHL	C1B-C2B	8.23	1.48	1.39
4	2	605	CHL	C1B-C2B	8.19	1.48	1.39
4	2	608	CHL	C1B-C2B	8.14	1.48	1.39
4	3	607	CHL	C1A-CHA	8.12	1.49	1.40
4	2	607	CHL	C1B-C2B	8.11	1.48	1.39
4	1	605	CHL	C1B-C2B	8.05	1.48	1.39
4	4	608	CHL	C1B-C2B	7.99	1.48	1.39
4	3	605	CHL	C1B-C2B	7.94	1.48	1.39
4	1	601	CHL	C1B-C2B	7.89	1.48	1.39
4	3	606	CHL	C3B-C4B	7.82	1.49	1.41
4	2	601	CHL	C1B-C2B	7.77	1.48	1.39
4	1	607	CHL	C1A-CHA	7.76	1.48	1.40
4	2	607	CHL	C1A-CHA	7.76	1.48	1.40
4	2	608	CHL	C1A-CHA	7.71	1.48	1.40
4	3	607	CHL	C1B-C2B	7.70	1.48	1.39
4	4	607	CHL	C1B-C2B	7.67	1.48	1.39
4	4	609	CHL	C1A-CHA	7.61	1.48	1.40
4	2	605	CHL	C1A-CHA	7.58	1.48	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	4	609	CHL	C1B-C2B	7.56	1.48	1.39
4	3	607	CHL	C3B-C4B	7.55	1.48	1.41
4	1	608	CHL	C3B-C4B	7.55	1.48	1.41
4	4	601	CHL	C3B-C4B	7.52	1.48	1.41
4	3	601	CHL	C1B-C2B	7.52	1.48	1.39
4	2	601	CHL	C3B-C4B	7.48	1.48	1.41
4	2	606	CHL	C1A-CHA	7.46	1.48	1.40
4	4	606	CHL	C3B-C4B	7.45	1.48	1.41
4	4	608	CHL	C3B-C4B	7.43	1.48	1.41
4	1	606	CHL	C1A-CHA	7.42	1.48	1.40
4	3	609	CHL	C1B-C2B	7.42	1.48	1.39
4	3	605	CHL	C1A-CHA	7.35	1.48	1.40
4	2	609	CHL	C1A-CHA	7.32	1.48	1.40
4	4	607	CHL	C1A-CHA	7.23	1.48	1.40
4	1	607	CHL	C3B-C4B	7.18	1.48	1.41
4	4	606	CHL	C1B-C2B	7.17	1.47	1.39
4	3	608	CHL	C1B-C2B	7.16	1.47	1.39
4	4	606	CHL	C1A-CHA	7.16	1.48	1.40
4	1	607	CHL	C1B-C2B	7.08	1.47	1.39
4	2	606	CHL	C1B-C2B	7.06	1.47	1.39
4	3	608	CHL	C1A-CHA	7.06	1.48	1.40
4	4	607	CHL	C3B-C4B	7.04	1.48	1.41
4	3	606	CHL	C1B-C2B	7.03	1.47	1.39
4	4	608	CHL	C1A-CHA	7.01	1.48	1.40
4	3	609	CHL	C1A-CHA	6.98	1.47	1.40
4	1	605	CHL	C1A-CHA	6.94	1.47	1.40
4	1	609	CHL	C1A-CHA	6.90	1.47	1.40
4	3	605	CHL	C3B-C4B	6.89	1.48	1.41
4	2	605	CHL	C3B-C4B	6.88	1.48	1.41
4	1	608	CHL	C1A-CHA	6.88	1.47	1.40
4	3	609	CHL	C3B-C4B	6.86	1.48	1.41
4	2	609	CHL	C1B-C2B	6.84	1.47	1.39
4	1	606	CHL	C1B-C2B	6.78	1.47	1.39
4	2	601	CHL	C1A-CHA	6.70	1.47	1.40
4	3	601	CHL	C1A-CHA	6.65	1.47	1.40
4	2	601	CHL	C3A-C2A	-6.63	1.49	1.54
4	3	606	CHL	C3D-C4D	-6.62	1.31	1.41
4	3	601	CHL	C3D-C4D	-6.60	1.31	1.41
4	1	605	CHL	C3B-C4B	6.59	1.47	1.41
4	1	601	CHL	C1A-CHA	6.58	1.47	1.40
4	1	601	CHL	C3D-C4D	-6.56	1.31	1.41
4	3	608	CHL	C3B-C4B	6.55	1.47	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	1	608	CHL	C3D-C4D	-6.51	1.31	1.41
4	1	609	CHL	C3D-C4D	-6.44	1.31	1.41
4	1	607	CHL	C3A-C2A	-6.43	1.49	1.54
4	3	606	CHL	C3A-C2A	-6.37	1.49	1.54
4	4	608	CHL	C3D-C4D	-6.37	1.31	1.41
4	4	608	CHL	C3A-C2A	-6.36	1.49	1.54
4	3	606	CHL	C1A-CHA	6.35	1.47	1.40
4	1	608	CHL	C3A-C2A	-6.34	1.49	1.54
4	1	608	CHL	C1B-C2B	6.27	1.46	1.39
4	1	609	CHL	C1B-C2B	6.24	1.46	1.39
4	2	607	CHL	C3B-C4B	6.24	1.47	1.41
4	2	608	CHL	C3B-C4B	6.23	1.47	1.41
4	2	601	CHL	C3D-C4D	-6.23	1.32	1.41
4	3	601	CHL	C3B-C4B	6.21	1.47	1.41
4	3	609	CHL	C3D-C4D	-6.19	1.32	1.41
4	1	601	CHL	C3B-C4B	6.16	1.47	1.41
4	2	609	CHL	C3A-C2A	-6.16	1.49	1.54
4	4	606	CHL	C3A-C2A	-6.14	1.49	1.54
4	1	606	CHL	C3D-C4D	-6.14	1.32	1.41
4	2	609	CHL	C3B-C4B	6.12	1.47	1.41
4	3	608	CHL	C3D-C4D	-6.12	1.32	1.41
4	4	609	CHL	C3B-C4B	6.10	1.47	1.41
4	2	609	CHL	C3D-C4D	-6.08	1.32	1.41
4	2	606	CHL	C3B-C4B	6.07	1.47	1.41
4	2	607	CHL	C3D-C4D	-6.07	1.32	1.41
4	1	609	CHL	C3A-C2A	-6.06	1.49	1.54
4	1	605	CHL	C3D-C4D	-6.05	1.32	1.41
4	1	609	CHL	C3B-C4B	6.05	1.47	1.41
4	1	606	CHL	C3B-C4B	6.05	1.47	1.41
4	3	607	CHL	C3D-C4D	-5.97	1.32	1.41
4	1	607	CHL	C3D-C4D	-5.97	1.32	1.41
4	4	609	CHL	C3D-C4D	-5.95	1.32	1.41
4	2	606	CHL	C3A-C2A	-5.87	1.49	1.54
4	2	606	CHL	C3D-C4D	-5.87	1.32	1.41
4	4	607	CHL	C3D-C4D	-5.84	1.32	1.41
4	4	606	CHL	C3D-C4D	-5.83	1.32	1.41
4	2	608	CHL	C3A-C2A	-5.81	1.49	1.54
4	2	608	CHL	C3D-C4D	-5.80	1.32	1.41
4	3	606	CHL	CHB-C4A	-5.76	1.31	1.38
4	3	608	CHL	C3A-C2A	-5.75	1.49	1.54
4	1	609	CHL	CHB-C4A	-5.75	1.31	1.38
4	4	609	CHL	C3A-C2A	-5.75	1.49	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	3	605	CHL	C3D-C4D	-5.73	1.32	1.41
4	3	605	CHL	C3A-C2A	-5.71	1.49	1.54
4	2	605	CHL	C3D-C4D	-5.63	1.32	1.41
4	4	607	CHL	C3A-C2A	-5.60	1.49	1.54
4	2	607	CHL	C3A-C2A	-5.58	1.49	1.54
4	1	606	CHL	C3A-C2A	-5.57	1.49	1.54
4	1	601	CHL	C3A-C2A	-5.53	1.49	1.54
4	3	601	CHL	C3A-C2A	-5.53	1.49	1.54
4	2	605	CHL	C3A-C2A	-5.53	1.49	1.54
4	1	607	CHL	CHB-C4A	-5.51	1.32	1.38
4	3	601	CHL	CHB-C4A	-5.50	1.32	1.38
4	1	608	CHL	CHB-C4A	-5.49	1.32	1.38
4	4	601	CHL	C3D-C4D	-5.41	1.33	1.41
4	4	606	CHL	CHB-C4A	-5.35	1.32	1.38
4	3	609	CHL	C3A-C2A	-5.30	1.50	1.54
4	4	601	CHL	C3B-C2B	5.29	1.47	1.40
4	2	606	CHL	CHB-C4A	-5.28	1.32	1.38
4	3	607	CHL	C3A-C2A	-5.26	1.50	1.54
4	3	609	CHL	CHB-C4A	-5.21	1.32	1.38
4	4	609	CHL	CHB-C4A	-5.16	1.32	1.38
4	2	607	CHL	CHB-C4A	-5.12	1.32	1.38
4	2	609	CHL	CHB-C4A	-5.07	1.32	1.38
4	2	607	CHL	O2D-CGD	5.06	1.45	1.33
4	3	607	CHL	CHB-C4A	-5.06	1.32	1.38
4	4	609	CHL	O2D-CGD	5.05	1.45	1.33
4	1	605	CHL	C3A-C2A	-5.04	1.50	1.54
4	3	608	CHL	CHB-C4A	-5.02	1.32	1.38
4	1	606	CHL	CHB-C4A	-5.00	1.32	1.38
4	2	605	CHL	O2D-CGD	4.98	1.45	1.33
4	2	601	CHL	O2D-CGD	4.96	1.45	1.33
4	2	605	CHL	CHB-C4A	-4.94	1.32	1.38
4	3	605	CHL	CHB-C4A	-4.94	1.32	1.38
4	4	607	CHL	CHB-C4A	-4.93	1.32	1.38
4	2	609	CHL	O2D-CGD	4.92	1.45	1.33
4	3	605	CHL	O2D-CGD	4.91	1.45	1.33
4	3	606	CHL	O2D-CGD	4.91	1.45	1.33
4	1	605	CHL	CHB-C4A	-4.90	1.32	1.38
4	4	608	CHL	O2D-CGD	4.90	1.45	1.33
4	2	606	CHL	O2D-CGD	4.90	1.45	1.33
4	4	608	CHL	CHB-C4A	-4.88	1.32	1.38
4	4	601	CHL	OBD-CAD	4.87	1.28	1.22
4	2	608	CHL	CHB-C4A	-4.85	1.32	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	4	606	CHL	O2D-CGD	4.84	1.45	1.33
4	2	605	CHL	OBD-CAD	4.83	1.28	1.22
4	1	605	CHL	O2D-CGD	4.82	1.45	1.33
4	2	601	CHL	CHB-C4A	-4.82	1.32	1.38
4	4	601	CHL	CHC-C4B	4.82	1.47	1.39
4	2	608	CHL	O2D-CGD	4.82	1.45	1.33
4	1	606	CHL	O2D-CGD	4.78	1.45	1.33
4	4	601	CHL	O2D-CGD	4.78	1.45	1.30
4	3	601	CHL	O2D-CGD	4.78	1.45	1.33
4	4	601	CHL	CHD-C4C	4.78	1.48	1.39
4	3	606	CHL	CHC-C4B	4.77	1.47	1.39
4	4	607	CHL	O2D-CGD	4.77	1.45	1.33
4	1	601	CHL	CHB-C4A	-4.77	1.32	1.38
4	4	601	CHL	CHD-C1D	4.77	1.47	1.39
4	3	608	CHL	O2D-CGD	4.73	1.44	1.33
4	2	605	CHL	CHC-C4B	4.72	1.47	1.39
4	4	608	CHL	CHB-C1B	4.72	1.47	1.39
4	3	609	CHL	C3B-C2B	4.71	1.46	1.40
4	4	607	CHL	OBD-CAD	4.68	1.28	1.22
4	1	601	CHL	O2D-CGD	4.68	1.44	1.33
4	2	605	CHL	C3B-C2B	4.67	1.46	1.40
4	3	609	CHL	O2D-CGD	4.67	1.44	1.33
4	1	609	CHL	CHD-C1D	4.66	1.47	1.39
4	3	605	CHL	OBD-CAD	4.66	1.28	1.22
4	1	607	CHL	O2D-CGD	4.66	1.44	1.33
4	1	608	CHL	O2D-CGD	4.65	1.44	1.33
4	3	607	CHL	O2D-CGD	4.64	1.44	1.33
4	3	605	CHL	O2A-CGA	4.62	1.46	1.30
4	4	601	CHL	CHB-C1B	4.60	1.47	1.39
4	4	609	CHL	OBD-CAD	4.60	1.28	1.22
4	1	609	CHL	O2D-CGD	4.60	1.44	1.33
4	4	606	CHL	OBD-CAD	4.59	1.28	1.22
4	2	608	CHL	O2A-CGA	4.58	1.46	1.30
4	1	605	CHL	OBD-CAD	4.57	1.28	1.22
4	4	607	CHL	C3B-C2B	4.57	1.46	1.40
4	3	609	CHL	OBD-CAD	4.57	1.28	1.22
4	4	601	CHL	CHC-C1C	4.57	1.48	1.39
4	4	601	CHL	O2A-CGA	4.56	1.45	1.30
4	2	607	CHL	OBD-CAD	4.56	1.28	1.22
4	3	608	CHL	CHC-C4B	4.55	1.47	1.39
4	2	601	CHL	C3B-C2B	4.55	1.46	1.40
4	4	606	CHL	CHC-C4B	4.55	1.47	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	1	608	CHL	CHC-C4B	4.55	1.47	1.39
4	1	601	CHL	O2A-CGA	4.54	1.45	1.30
4	4	601	CHL	CHB-C4A	-4.54	1.33	1.38
4	2	605	CHL	O2A-CGA	4.53	1.45	1.30
4	2	601	CHL	O2A-CGA	4.53	1.45	1.30
4	2	606	CHL	O2A-CGA	4.53	1.45	1.30
4	3	607	CHL	OBD-CAD	4.53	1.28	1.22
4	3	606	CHL	O2A-CGA	4.52	1.45	1.30
4	1	606	CHL	O2A-CGA	4.52	1.45	1.30
4	2	609	CHL	OBD-CAD	4.52	1.28	1.22
4	2	607	CHL	C3B-C2B	4.52	1.46	1.40
4	4	607	CHL	O2A-CGA	4.51	1.45	1.30
4	4	609	CHL	O2A-CGA	4.51	1.45	1.30
4	1	605	CHL	C3B-C2B	4.50	1.46	1.40
4	3	609	CHL	CHD-C1D	4.50	1.46	1.39
4	1	608	CHL	O2A-CGA	4.50	1.45	1.30
4	3	605	CHL	C3B-C2B	4.49	1.46	1.40
4	1	607	CHL	OBD-CAD	4.49	1.28	1.22
4	3	608	CHL	O2A-CGA	4.49	1.45	1.30
4	1	609	CHL	CHD-C4C	4.49	1.48	1.39
4	2	606	CHL	OBD-CAD	4.49	1.28	1.22
4	2	601	CHL	OBD-CAD	4.48	1.28	1.22
4	2	607	CHL	CHD-C1D	4.48	1.46	1.39
4	1	605	CHL	O2A-CGA	4.47	1.45	1.30
4	1	609	CHL	OBD-CAD	4.47	1.28	1.22
4	3	606	CHL	OBD-CAD	4.45	1.28	1.22
4	1	606	CHL	OBD-CAD	4.45	1.28	1.22
4	1	606	CHL	CHC-C4B	4.45	1.46	1.39
4	3	609	CHL	CHD-C4C	4.44	1.47	1.39
4	1	605	CHL	CHB-C1B	4.43	1.46	1.39
4	3	607	CHL	CHC-C4B	4.43	1.46	1.39
4	3	609	CHL	O2A-CGA	4.43	1.46	1.33
4	4	606	CHL	O2A-CGA	4.43	1.45	1.30
4	4	607	CHL	CHB-C1B	4.42	1.46	1.39
4	3	608	CHL	CHC-C1C	4.41	1.47	1.39
4	4	607	CHL	CHD-C4C	4.41	1.47	1.39
4	2	601	CHL	CHB-C1B	4.39	1.46	1.39
4	1	601	CHL	CHB-C1B	4.38	1.46	1.39
4	2	607	CHL	CHD-C4C	4.37	1.47	1.39
4	4	608	CHL	O2A-CGA	4.37	1.45	1.30
4	4	606	CHL	C3B-C2B	4.36	1.46	1.40
4	3	609	CHL	CHB-C1B	4.34	1.46	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	2	605	CHL	CHB-C1B	4.34	1.46	1.39
4	1	601	CHL	OBD-CAD	4.33	1.28	1.22
4	3	607	CHL	O2A-CGA	4.33	1.46	1.33
4	2	601	CHL	CHC-C4B	4.33	1.46	1.39
4	4	607	CHL	CHC-C4B	4.31	1.46	1.39
4	3	607	CHL	CHC-C1C	4.31	1.47	1.39
4	2	608	CHL	CHC-C4B	4.31	1.46	1.39
4	1	601	CHL	CHD-C4C	4.31	1.47	1.39
4	3	605	CHL	CHC-C4B	4.29	1.46	1.39
4	2	608	CHL	OBD-CAD	4.29	1.27	1.22
4	2	609	CHL	CHB-C1B	4.28	1.46	1.39
4	2	609	CHL	C3B-C2B	4.28	1.46	1.40
4	1	607	CHL	O2A-CGA	4.25	1.45	1.33
4	2	605	CHL	CHC-C1C	4.25	1.47	1.39
4	2	607	CHL	CHB-C1B	4.25	1.46	1.39
4	1	609	CHL	O2A-CGA	4.24	1.45	1.33
4	2	601	CHL	CHD-C4C	4.24	1.47	1.39
4	4	601	CHL	C3D-C2D	4.23	1.46	1.39
4	2	605	CHL	CHD-C4C	4.23	1.47	1.39
4	4	608	CHL	C3B-C2B	4.23	1.46	1.40
4	1	601	CHL	C3B-C2B	4.22	1.46	1.40
8	1	1623	NEX	C7-C6	-4.21	1.25	1.30
4	4	609	CHL	CHC-C4B	4.21	1.46	1.39
4	2	609	CHL	O2A-CGA	4.21	1.45	1.33
4	3	605	CHL	CHB-C1B	4.21	1.46	1.39
4	2	608	CHL	C3B-C2B	4.20	1.46	1.40
4	4	607	CHL	CHD-C1D	4.20	1.46	1.39
4	3	607	CHL	C3B-C2B	4.20	1.46	1.40
4	2	606	CHL	CHD-C1D	4.19	1.46	1.39
4	3	608	CHL	CHB-C1B	4.19	1.46	1.39
4	3	607	CHL	CHB-C1B	4.19	1.46	1.39
4	3	601	CHL	O2A-CGA	4.18	1.45	1.33
4	2	607	CHL	CHC-C1C	4.17	1.47	1.39
4	2	609	CHL	CHD-C4C	4.17	1.47	1.39
4	1	608	CHL	CHC-C1C	4.16	1.47	1.39
4	3	606	CHL	CHD-C1D	4.16	1.46	1.39
4	2	607	CHL	O2A-CGA	4.14	1.45	1.33
4	3	606	CHL	CHC-C1C	4.14	1.47	1.39
4	3	606	CHL	CHD-C4C	4.12	1.47	1.39
4	2	609	CHL	CHC-C4B	4.12	1.46	1.39
4	2	606	CHL	CHB-C1B	4.11	1.46	1.39
4	3	606	CHL	C3B-C2B	4.11	1.45	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	3	609	CHL	CHC-C4B	4.10	1.46	1.39
4	2	608	CHL	CHB-C1B	4.10	1.46	1.39
4	3	601	CHL	CHC-C1C	4.09	1.47	1.39
4	1	605	CHL	CHD-C4C	4.08	1.47	1.39
4	4	608	CHL	CHC-C4B	4.08	1.46	1.39
4	1	607	CHL	CHD-C1D	4.08	1.46	1.39
4	4	606	CHL	CHD-C4C	4.07	1.47	1.39
4	1	607	CHL	CHD-C4C	4.06	1.47	1.39
4	3	605	CHL	CHD-C4C	4.06	1.47	1.39
4	1	601	CHL	CHC-C4B	4.05	1.46	1.39
4	3	601	CHL	OBD-CAD	4.05	1.27	1.22
4	4	606	CHL	CHC-C1C	4.05	1.47	1.39
4	1	609	CHL	CHB-C1B	4.04	1.46	1.39
4	4	606	CHL	CHD-C1D	4.04	1.46	1.39
4	2	606	CHL	CHD-C4C	4.03	1.47	1.39
4	3	608	CHL	CHD-C4C	4.03	1.47	1.39
4	4	606	CHL	CHB-C1B	4.01	1.46	1.39
4	1	606	CHL	C3B-C2B	4.01	1.45	1.40
4	2	608	CHL	CHD-C4C	4.00	1.47	1.39
4	3	601	CHL	CHC-C4B	4.00	1.46	1.39
4	4	607	CHL	CHC-C1C	3.99	1.47	1.39
4	1	605	CHL	CHC-C4B	3.99	1.46	1.39
4	2	606	CHL	CHC-C4B	3.99	1.46	1.39
4	3	608	CHL	C3B-C2B	3.99	1.45	1.40
4	4	609	CHL	CHD-C4C	3.98	1.47	1.39
4	2	609	CHL	CHD-C1D	3.98	1.46	1.39
4	2	609	CHL	CHC-C1C	3.98	1.47	1.39
4	2	607	CHL	CHC-C4B	3.97	1.46	1.39
4	4	609	CHL	CHB-C1B	3.97	1.46	1.39
4	3	606	CHL	CHB-C1B	3.96	1.46	1.39
4	3	607	CHL	CHD-C1D	3.96	1.46	1.39
4	1	606	CHL	CHC-C1C	3.95	1.46	1.39
4	3	608	CHL	OBD-CAD	3.94	1.27	1.22
4	2	605	CHL	CHD-C1D	3.94	1.46	1.39
4	1	609	CHL	C3B-C2B	3.93	1.45	1.40
4	1	608	CHL	CHB-C1B	3.91	1.46	1.39
4	4	609	CHL	C3B-C2B	3.89	1.45	1.40
4	3	601	CHL	C3B-C2B	3.89	1.45	1.40
4	1	605	CHL	CHC-C1C	3.89	1.46	1.39
4	2	606	CHL	C3B-C2B	3.89	1.45	1.40
4	3	601	CHL	CHD-C4C	3.89	1.46	1.39
4	1	601	CHL	CHD-C1D	3.88	1.45	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	1	605	CHL	CHD-C1D	3.87	1.45	1.39
4	4	609	CHL	CHD-C1D	3.87	1.45	1.39
4	3	609	CHL	CHC-C1C	3.85	1.46	1.39
4	1	606	CHL	CHB-C1B	3.85	1.45	1.39
4	1	608	CHL	OBD-CAD	3.85	1.27	1.22
4	2	601	CHL	CHD-C1D	3.84	1.45	1.39
4	3	605	CHL	CHD-C1D	3.84	1.45	1.39
4	4	608	CHL	CHD-C4C	3.81	1.46	1.39
4	4	608	CHL	CHC-C1C	3.81	1.46	1.39
4	1	601	CHL	CHC-C1C	3.80	1.46	1.39
4	2	608	CHL	CHC-C1C	3.79	1.46	1.39
4	1	608	CHL	CBD-CGD	-3.79	1.47	1.52
4	2	608	CHL	CHD-C1D	3.76	1.45	1.39
4	1	607	CHL	CHC-C4B	3.76	1.45	1.39
4	3	607	CHL	CHD-C4C	3.76	1.46	1.39
4	3	605	CHL	C3D-C2D	3.75	1.46	1.39
4	1	608	CHL	CHD-C4C	3.74	1.46	1.39
4	1	607	CHL	CHC-C1C	3.74	1.46	1.39
4	1	605	CHL	C3D-C2D	3.73	1.46	1.39
4	2	605	CHL	C3D-C2D	3.72	1.46	1.39
4	4	609	CHL	CHC-C1C	3.71	1.46	1.39
4	1	607	CHL	CHB-C1B	3.71	1.45	1.39
4	3	605	CHL	CHC-C1C	3.70	1.46	1.39
4	2	606	CHL	CHC-C1C	3.69	1.46	1.39
4	1	607	CHL	C3B-C2B	3.66	1.45	1.40
4	3	601	CHL	CHB-C1B	3.65	1.45	1.39
4	2	607	CHL	C3D-C2D	3.64	1.45	1.39
4	1	606	CHL	CHD-C4C	3.63	1.46	1.39
4	1	607	CHL	CBD-CGD	-3.62	1.47	1.52
4	1	609	CHL	CHC-C1C	3.61	1.46	1.39
4	2	601	CHL	CHC-C1C	3.61	1.46	1.39
4	4	607	CHL	C3D-C2D	3.61	1.45	1.39
4	1	606	CHL	CHD-C1D	3.60	1.45	1.39
4	1	601	CHL	CBD-CGD	-3.58	1.48	1.52
4	4	608	CHL	OBD-CAD	3.53	1.26	1.22
5	3	614	CLA	C4B-NB	3.52	1.42	1.37
4	1	609	CHL	CBD-CGD	-3.52	1.48	1.52
4	1	608	CHL	CHD-C1D	3.51	1.45	1.39
4	2	609	CHL	C3D-C2D	3.51	1.45	1.39
4	4	606	CHL	C3D-C2D	3.50	1.45	1.39
4	3	601	CHL	CHD-C1D	3.49	1.45	1.39
8	1	1623	NEX	C7-C8	-3.49	1.26	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	3	614	CLA	C1D-ND	3.47	1.42	1.37
4	3	601	CHL	CBD-CGD	-3.44	1.48	1.52
4	4	601	CHL	C3A-C2A	-3.43	1.51	1.54
4	3	609	CHL	C3D-C2D	3.42	1.45	1.39
4	4	608	CHL	CBD-CGD	-3.41	1.48	1.52
4	3	607	CHL	C3D-C2D	3.41	1.45	1.39
4	4	608	CHL	CHD-C1D	3.40	1.45	1.39
4	1	609	CHL	CHC-C4B	3.40	1.45	1.39
4	1	608	CHL	C3B-C2B	3.39	1.45	1.40
4	2	609	CHL	CBD-CGD	-3.35	1.48	1.52
5	3	611	CLA	CMB-C2B	-3.33	1.43	1.50
4	2	606	CHL	C3D-C2D	3.32	1.45	1.39
4	4	606	CHL	CBD-CGD	-3.32	1.48	1.52
4	2	608	CHL	C3D-C2D	3.30	1.45	1.39
4	3	607	CHL	CBD-CGD	-3.30	1.48	1.52
5	2	604	CLA	C1D-ND	3.23	1.42	1.37
8	2	1623	NEX	C7-C8	-3.22	1.26	1.31
4	1	606	CHL	CBD-CGD	-3.21	1.48	1.52
4	1	609	CHL	C3D-C2D	3.20	1.45	1.39
4	3	608	CHL	CBD-CGD	-3.19	1.48	1.52
4	4	609	CHL	C3D-C2D	3.19	1.45	1.39
4	2	608	CHL	CBD-CGD	-3.19	1.48	1.52
4	3	608	CHL	CHD-C1D	3.18	1.44	1.39
9	3	2630	LHG	O7-C5	-3.18	1.39	1.46
4	3	605	CHL	CBD-CGD	-3.18	1.48	1.52
4	4	608	CHL	C3D-C2D	3.17	1.45	1.39
4	3	609	CHL	CBD-CGD	-3.13	1.48	1.52
5	1	612	CLA	C1D-ND	3.12	1.42	1.37
5	1	604	CLA	CMB-C2B	-3.09	1.44	1.50
8	2	1623	NEX	C7-C6	-3.08	1.27	1.30
5	2	612	CLA	C1D-ND	3.06	1.41	1.37
4	1	606	CHL	C3D-C2D	3.04	1.44	1.39
5	4	611	CLA	C1D-ND	3.02	1.41	1.37
5	2	613	CLA	C4B-NB	3.01	1.41	1.37
4	2	601	CHL	CBD-CGD	-2.99	1.48	1.52
5	2	612	CLA	C4B-NB	2.95	1.41	1.37
4	1	607	CHL	C3D-C2D	2.91	1.44	1.39
5	3	604	CLA	C1D-ND	2.91	1.41	1.37
5	4	604	CLA	C1D-ND	2.91	1.41	1.37
5	1	614	CLA	C1D-ND	2.89	1.41	1.37
4	2	605	CHL	CBD-CGD	-2.88	1.48	1.52
4	2	606	CHL	CBD-CGD	-2.87	1.48	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	4	607	CHL	CBD-CGD	-2.87	1.48	1.52
5	2	614	CLA	C1D-ND	2.86	1.41	1.37
5	3	611	CLA	CMD-C2D	-2.84	1.44	1.50
5	2	611	CLA	C1D-ND	2.83	1.41	1.37
4	3	608	CHL	C3D-C2D	2.83	1.44	1.39
5	4	610	CLA	CMC-C2C	-2.82	1.45	1.50
4	3	601	CHL	C3D-C2D	2.82	1.44	1.39
4	3	606	CHL	C3D-C2D	2.81	1.44	1.39
4	1	608	CHL	C3D-C2D	2.80	1.44	1.39
5	4	603	CLA	C1D-ND	2.80	1.41	1.37
5	2	610	CLA	C1D-ND	2.80	1.41	1.37
5	4	610	CLA	CMD-C2D	-2.79	1.45	1.50
5	3	613	CLA	C1D-ND	2.79	1.41	1.37
8	3	1623	NEX	C7-C8	-2.78	1.27	1.31
5	3	604	CLA	C4B-NB	2.78	1.41	1.37
5	2	603	CLA	C1D-ND	2.77	1.41	1.37
5	3	602	CLA	CMC-C2C	-2.77	1.45	1.50
5	3	612	CLA	CMD-C2D	-2.77	1.45	1.50
5	4	612	CLA	C1D-ND	2.77	1.41	1.37
5	1	611	CLA	C1D-ND	2.77	1.41	1.37
5	3	603	CLA	CMD-C2D	-2.76	1.45	1.50
6	4	620	LUT	C22-C21	-2.76	1.51	1.54
5	4	603	CLA	C4B-NB	2.76	1.41	1.37
4	3	606	CHL	CBD-CGD	-2.75	1.49	1.52
4	2	601	CHL	C3D-C2D	2.75	1.44	1.39
5	2	602	CLA	CMC-C2C	-2.74	1.45	1.50
5	4	612	CLA	C4B-NB	2.73	1.41	1.37
6	3	1620	LUT	C22-C21	-2.72	1.51	1.54
5	1	613	CLA	C1D-ND	2.72	1.41	1.37
5	3	611	CLA	MG-NB	-2.72	2.00	2.05
5	4	611	CLA	C4B-NB	2.71	1.41	1.37
5	2	613	CLA	MG-ND	-2.70	2.00	2.05
5	3	613	CLA	C4B-NB	2.70	1.41	1.37
5	1	602	CLA	C1D-ND	2.70	1.41	1.37
5	2	614	CLA	C4B-NB	2.69	1.41	1.37
5	1	610	CLA	C1D-ND	2.69	1.41	1.37
5	1	612	CLA	C4B-NB	2.69	1.41	1.37
6	1	1621	LUT	C1-C6	-2.69	1.50	1.53
5	1	613	CLA	MG-ND	-2.67	2.00	2.05
5	2	613	CLA	CMD-C2D	-2.66	1.45	1.50
5	1	614	CLA	C4B-NB	2.66	1.41	1.37
5	1	613	CLA	C4B-NB	2.66	1.41	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	3	1622	XAT	O4-C5	-2.66	1.42	1.46
5	4	610	CLA	C1D-ND	2.64	1.41	1.37
5	1	603	CLA	CMB-C2B	-2.61	1.45	1.50
4	4	609	CHL	CBD-CGD	-2.61	1.49	1.52
5	3	610	CLA	MG-NB	-2.60	2.00	2.05
5	1	604	CLA	MG-NB	-2.59	2.00	2.05
5	4	603	CLA	CMD-C2D	-2.59	1.45	1.50
5	3	604	CLA	CMB-C2B	-2.59	1.45	1.50
9	1	2630	LHG	O7-C5	-2.59	1.40	1.46
5	3	614	CLA	C3B-C4B	2.59	1.50	1.42
5	4	604	CLA	CMB-C2B	-2.58	1.45	1.50
5	3	602	CLA	C1B-C2B	2.58	1.49	1.43
5	1	602	CLA	CMC-C2C	-2.58	1.45	1.50
5	2	611	CLA	C4B-NB	2.58	1.41	1.37
8	1	1623	NEX	O24-C25	-2.58	1.42	1.46
5	3	602	CLA	CMD-C2D	-2.58	1.45	1.50
4	2	607	CHL	CBD-CGD	-2.57	1.49	1.52
5	2	610	CLA	C4B-NB	2.57	1.41	1.37
5	3	610	CLA	C1D-ND	2.56	1.41	1.37
5	4	602	CLA	C4B-NB	2.56	1.41	1.37
5	3	613	CLA	C1B-C2B	2.56	1.49	1.43
5	3	603	CLA	MG-ND	-2.56	2.00	2.05
5	2	611	CLA	MG-NB	-2.55	2.00	2.05
5	4	611	CLA	C1B-C2B	2.55	1.49	1.43
5	1	603	CLA	C1D-ND	2.55	1.41	1.37
5	2	603	CLA	C4B-NB	2.55	1.41	1.37
5	3	611	CLA	MG-ND	-2.54	2.00	2.05
5	1	603	CLA	MG-ND	-2.54	2.00	2.05
5	2	602	CLA	CMD-C2D	-2.54	1.45	1.50
5	2	613	CLA	C1D-ND	2.52	1.41	1.37
5	4	602	CLA	C1D-ND	2.52	1.41	1.37
5	3	603	CLA	C1D-ND	2.52	1.41	1.37
5	2	602	CLA	C4B-NB	2.51	1.41	1.37
5	3	611	CLA	C4B-NB	2.51	1.41	1.37
6	3	1621	LUT	C22-C21	-2.51	1.51	1.54
4	1	601	CHL	C3D-C2D	2.50	1.43	1.39
5	2	614	CLA	CMD-C2D	-2.50	1.45	1.50
5	3	611	CLA	C1D-ND	2.49	1.41	1.37
5	1	604	CLA	MG-ND	-2.49	2.00	2.05
5	2	604	CLA	C4B-NB	2.49	1.41	1.37
5	3	613	CLA	CMD-C2D	-2.49	1.45	1.50
5	1	614	CLA	CMD-C2D	-2.49	1.45	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	2	603	CLA	MG-ND	-2.48	2.00	2.05
7	4	622	XAT	O4-C5	-2.48	1.43	1.46
5	4	612	CLA	CMD-C2D	-2.48	1.45	1.50
5	1	602	CLA	MG-ND	-2.48	2.00	2.05
5	2	603	CLA	CMD-C2D	-2.47	1.45	1.50
5	3	610	CLA	CMB-C2B	-2.47	1.45	1.50
5	1	602	CLA	CMD-C2D	-2.47	1.45	1.50
5	3	613	CLA	MG-ND	-2.47	2.00	2.05
4	1	605	CHL	CBD-CGD	-2.47	1.49	1.52
5	3	612	CLA	MG-ND	-2.46	2.00	2.05
7	1	1622	XAT	O4-C5	-2.46	1.43	1.46
5	1	604	CLA	CMD-C2D	-2.45	1.45	1.50
5	3	612	CLA	CMC-C2C	-2.45	1.45	1.50
5	3	612	CLA	MG-NB	-2.45	2.00	2.05
5	1	611	CLA	C4B-NB	2.45	1.41	1.37
5	1	611	CLA	MG-NB	-2.45	2.00	2.05
5	4	603	CLA	MG-ND	-2.45	2.00	2.05
5	4	610	CLA	MG-NB	-2.45	2.00	2.05
5	2	602	CLA	MG-NB	-2.44	2.00	2.05
5	1	604	CLA	C4B-NB	2.44	1.41	1.37
6	1	1621	LUT	C22-C21	-2.44	1.51	1.54
5	1	603	CLA	CMD-C2D	-2.44	1.45	1.50
5	2	611	CLA	MG-ND	-2.44	2.01	2.05
5	3	612	CLA	C1D-ND	2.44	1.41	1.37
4	1	608	CHL	CHA-CBD	-2.43	1.48	1.51
8	3	1623	NEX	O24-C25	-2.43	1.43	1.46
5	1	613	CLA	CMD-C2D	-2.42	1.45	1.50
5	2	603	CLA	CMC-C2C	-2.42	1.45	1.50
5	2	610	CLA	CMC-C2C	-2.41	1.45	1.50
5	3	602	CLA	MG-ND	-2.41	2.01	2.05
5	1	604	CLA	C1D-ND	2.40	1.41	1.37
5	4	612	CLA	MG-NB	-2.40	2.01	2.05
5	1	603	CLA	CMC-C2C	-2.40	1.45	1.50
5	1	602	CLA	MG-NB	-2.39	2.01	2.05
5	1	613	CLA	CMB-C2B	-2.38	1.45	1.50
5	1	612	CLA	CMC-C2C	-2.38	1.45	1.50
5	1	612	CLA	MG-ND	-2.38	2.01	2.05
5	2	610	CLA	CMB-C2B	-2.37	1.45	1.50
5	2	603	CLA	C1B-C2B	2.37	1.48	1.43
5	2	614	CLA	MG-NB	-2.37	2.01	2.05
5	4	604	CLA	C4B-NB	2.36	1.41	1.37
5	3	612	CLA	CMB-C2B	-2.36	1.45	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	1	610	CLA	MG-NB	-2.36	2.01	2.05
5	3	613	CLA	CMB-C2B	-2.35	1.45	1.50
5	3	610	CLA	CMD-C2D	-2.35	1.45	1.50
5	2	612	CLA	CMB-C2B	-2.35	1.45	1.50
5	2	604	CLA	CMB-C2B	-2.35	1.46	1.50
9	2	2630	LHG	O7-C5	-2.35	1.41	1.46
5	4	604	CLA	MG-NB	-2.35	2.01	2.05
5	3	610	CLA	C4B-NB	2.35	1.40	1.37
5	3	612	CLA	C1B-NB	-2.35	1.34	1.37
5	4	610	CLA	C4B-NB	2.35	1.40	1.37
5	3	610	CLA	CMC-C2C	-2.35	1.46	1.50
5	4	611	CLA	CMD-C2D	-2.34	1.46	1.50
5	1	612	CLA	CMD-C2D	-2.34	1.46	1.50
5	4	603	CLA	C1B-C2B	2.34	1.48	1.43
5	2	610	CLA	CMD-C2D	-2.34	1.46	1.50
5	1	610	CLA	CMC-C2C	-2.34	1.46	1.50
5	2	611	CLA	CMD-C2D	-2.33	1.46	1.50
5	2	604	CLA	MG-NB	-2.33	2.01	2.05
7	3	1622	XAT	O24-C25	-2.33	1.43	1.46
5	3	611	CLA	C1B-NB	-2.32	1.34	1.37
5	1	611	CLA	CMB-C2B	-2.32	1.46	1.50
4	1	609	CHL	CHA-CBD	-2.32	1.48	1.51
5	1	612	CLA	CMB-C2B	-2.32	1.46	1.50
5	3	613	CLA	CMC-C2C	-2.32	1.46	1.50
5	2	604	CLA	CMC-C2C	-2.32	1.46	1.50
5	3	603	CLA	CMC-C2C	-2.32	1.46	1.50
5	3	614	CLA	C1B-C2B	2.32	1.48	1.43
5	2	604	CLA	CMD-C2D	-2.32	1.46	1.50
5	4	604	CLA	MG-ND	-2.32	2.01	2.05
5	3	604	CLA	C1B-C2B	2.32	1.48	1.43
5	2	604	CLA	C1B-C2B	2.31	1.48	1.43
5	3	611	CLA	CHC-C4B	-2.31	1.34	1.39
5	3	610	CLA	MG-ND	-2.31	2.01	2.05
5	4	604	CLA	CMD-C2D	-2.31	1.46	1.50
5	2	613	CLA	C1B-C2B	2.31	1.48	1.43
5	4	602	CLA	CMC-C2C	-2.30	1.46	1.50
5	3	614	CLA	CMB-C2B	-2.30	1.46	1.50
5	2	602	CLA	C1B-C2B	2.30	1.48	1.43
5	1	603	CLA	CHC-C4B	-2.29	1.34	1.39
5	2	611	CLA	C1B-C2B	2.29	1.48	1.43
5	4	610	CLA	MG-ND	-2.29	2.01	2.05
5	4	602	CLA	CMB-C2B	-2.28	1.46	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	1	612	CLA	C1B-C2B	2.28	1.48	1.43
5	1	602	CLA	C4B-NB	2.28	1.40	1.37
5	3	602	CLA	C1B-NB	-2.28	1.34	1.37
5	2	610	CLA	C1B-C2B	2.28	1.48	1.43
5	1	611	CLA	C1B-C2B	2.28	1.48	1.43
5	1	602	CLA	CMB-C2B	-2.27	1.46	1.50
5	4	612	CLA	C3B-C4B	2.27	1.49	1.42
5	3	603	CLA	CMB-C2B	-2.27	1.46	1.50
5	4	612	CLA	CMC-C2C	-2.27	1.46	1.50
6	4	620	LUT	C1-C6	-2.27	1.50	1.53
5	3	602	CLA	C4B-NB	2.27	1.40	1.37
5	2	612	CLA	C1B-C2B	2.26	1.48	1.43
5	1	613	CLA	C1B-C2B	2.26	1.48	1.43
5	1	614	CLA	CMB-C2B	-2.26	1.46	1.50
5	4	603	CLA	CMC-C2C	-2.26	1.46	1.50
5	3	602	CLA	MG-NB	-2.25	2.01	2.05
5	1	612	CLA	CHC-C4B	-2.25	1.34	1.39
5	1	604	CLA	CMC-C2C	-2.25	1.46	1.50
5	2	602	CLA	MG-ND	-2.25	2.01	2.05
5	1	603	CLA	CAC-C3C	-2.25	1.45	1.51
5	1	603	CLA	C1B-C2B	2.25	1.48	1.43
7	4	622	XAT	O24-C25	-2.25	1.43	1.46
5	2	612	CLA	C3B-C4B	2.25	1.49	1.42
5	4	602	CLA	C3B-C4B	2.25	1.49	1.42
5	2	602	CLA	C1D-ND	2.24	1.40	1.37
5	4	604	CLA	C1B-NB	-2.24	1.34	1.37
5	1	614	CLA	MG-ND	-2.24	2.01	2.05
4	4	608	CHL	CHA-CBD	-2.24	1.49	1.51
5	3	613	CLA	CHC-C4B	-2.24	1.34	1.39
5	3	603	CLA	C1B-C2B	2.24	1.48	1.43
5	4	612	CLA	C1B-C2B	2.23	1.48	1.43
5	4	603	CLA	CMB-C2B	-2.23	1.46	1.50
5	1	611	CLA	CMC-C2C	-2.23	1.46	1.50
5	2	614	CLA	CMB-C2B	-2.23	1.46	1.50
5	1	613	CLA	CMC-C2C	-2.23	1.46	1.50
5	1	610	CLA	CMD-C2D	-2.22	1.46	1.50
5	2	610	CLA	MG-NB	-2.22	2.01	2.05
7	1	1622	XAT	O24-C25	-2.22	1.43	1.46
5	3	602	CLA	C1D-ND	2.22	1.40	1.37
5	3	604	CLA	CMD-C2D	-2.22	1.46	1.50
5	1	611	CLA	MG-ND	-2.21	2.01	2.05
5	2	613	CLA	CMB-C2B	-2.21	1.46	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	1	611	CLA	CMD-C2D	-2.20	1.46	1.50
8	1	1623	NEX	C22-C21	-2.20	1.51	1.54
5	3	604	CLA	CMC-C2C	-2.20	1.46	1.50
5	2	612	CLA	CMD-C2D	-2.20	1.46	1.50
5	2	603	CLA	CMB-C2B	-2.20	1.46	1.50
5	4	602	CLA	MG-NB	-2.20	2.01	2.05
5	3	612	CLA	C4B-NB	2.20	1.40	1.37
5	2	612	CLA	MG-ND	-2.20	2.01	2.05
5	3	602	CLA	CMB-C2B	-2.19	1.46	1.50
5	1	614	CLA	MG-NB	-2.19	2.01	2.05
5	4	604	CLA	CMC-C2C	-2.19	1.46	1.50
5	3	612	CLA	C1B-C2B	2.19	1.48	1.43
5	2	604	CLA	CHC-C4B	-2.19	1.34	1.39
5	2	614	CLA	MG-ND	-2.19	2.01	2.05
4	3	606	CHL	CHA-CBD	-2.19	1.49	1.51
5	1	610	CLA	CMB-C2B	-2.18	1.46	1.50
5	3	603	CLA	CAC-C3C	-2.18	1.45	1.51
5	1	614	CLA	CMC-C2C	-2.18	1.46	1.50
5	4	612	CLA	MG-ND	-2.18	2.01	2.05
5	1	602	CLA	C1B-C2B	2.18	1.48	1.43
4	1	606	CHL	MG-ND	-2.17	2.01	2.05
5	3	603	CLA	MG-NB	-2.17	2.01	2.05
5	2	613	CLA	CMC-C2C	-2.17	1.46	1.50
5	2	611	CLA	CMC-C2C	-2.16	1.46	1.50
4	3	608	CHL	CMC-C2C	2.16	1.49	1.44
5	4	611	CLA	CMB-C2B	-2.16	1.46	1.50
5	3	614	CLA	CMD-C2D	-2.16	1.46	1.50
5	1	613	CLA	CHC-C4B	-2.16	1.34	1.39
5	2	614	CLA	C3B-C4B	2.16	1.49	1.42
5	2	614	CLA	C1B-C2B	2.15	1.48	1.43
5	2	611	CLA	CMB-C2B	-2.15	1.46	1.50
5	2	603	CLA	C3B-C4B	2.15	1.49	1.42
5	4	603	CLA	C3B-C4B	2.14	1.49	1.42
4	1	607	CHL	C4A-C3A	-2.14	1.46	1.51
5	1	603	CLA	MG-NB	-2.14	2.01	2.05
5	2	614	CLA	CMC-C2C	-2.14	1.46	1.50
8	2	1623	NEX	O24-C25	-2.14	1.43	1.46
7	2	1622	XAT	O4-C5	-2.14	1.43	1.46
5	3	603	CLA	C4B-NB	2.14	1.40	1.37
4	2	609	CHL	CHA-CBD	-2.13	1.49	1.51
5	3	604	CLA	MG-NB	-2.13	2.01	2.05
4	2	601	CHL	CMC-C2C	2.13	1.49	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	2	1621	LUT	C22-C21	-2.13	1.52	1.54
5	2	602	CLA	CMB-C2B	-2.13	1.46	1.50
5	2	604	CLA	MG-ND	-2.13	2.01	2.05
4	1	605	CHL	CHA-CBD	-2.13	1.49	1.51
5	1	614	CLA	C1B-C2B	2.13	1.48	1.43
5	3	612	CLA	CHC-C4B	-2.13	1.34	1.39
5	4	610	CLA	CMB-C2B	-2.12	1.46	1.50
5	1	602	CLA	C1B-NB	-2.12	1.35	1.37
5	3	603	CLA	CHC-C4B	-2.11	1.34	1.39
5	1	613	CLA	MG-NB	-2.11	2.01	2.05
7	2	1622	XAT	O24-C25	-2.10	1.43	1.46
5	4	603	CLA	MG-NB	-2.10	2.01	2.05
5	4	611	CLA	MG-NB	-2.10	2.01	2.05
5	3	604	CLA	MG-ND	-2.10	2.01	2.05
5	3	611	CLA	CMC-C2C	-2.10	1.46	1.50
5	1	602	CLA	C3B-C4B	2.10	1.48	1.42
5	4	602	CLA	CMD-C2D	-2.09	1.46	1.50
5	3	611	CLA	C1B-C2B	2.09	1.48	1.43
5	2	603	CLA	MG-NB	-2.09	2.01	2.05
5	1	603	CLA	C1B-NB	-2.09	1.35	1.37
5	1	614	CLA	C3B-C4B	2.09	1.48	1.42
5	3	603	CLA	C1B-NB	-2.09	1.35	1.37
5	2	602	CLA	C3B-C4B	2.08	1.48	1.42
5	1	604	CLA	C1B-NB	-2.08	1.35	1.37
5	4	602	CLA	MG-ND	-2.08	2.01	2.05
4	2	607	CHL	MG-ND	-2.07	2.01	2.05
5	4	610	CLA	C1B-C2B	2.07	1.48	1.43
5	4	612	CLA	CMB-C2B	-2.07	1.46	1.50
5	1	611	CLA	O2A-CGA	2.07	1.37	1.30
5	2	612	CLA	CMC-C2C	-2.05	1.46	1.50
5	4	604	CLA	C1B-C2B	2.05	1.47	1.43
5	1	604	CLA	CHC-C4B	-2.05	1.34	1.39
5	4	611	CLA	C3B-C4B	2.04	1.48	1.42
5	1	611	CLA	C1B-NB	-2.04	1.35	1.37
5	3	613	CLA	CAC-C3C	-2.04	1.45	1.51
5	4	611	CLA	CMC-C2C	-2.04	1.46	1.50
4	3	601	CHL	CAC-C3C	-2.04	1.48	1.51
5	4	610	CLA	C3B-C4B	2.04	1.48	1.42
5	1	612	CLA	C1B-NB	-2.03	1.35	1.37
6	3	1620	LUT	C1-C6	-2.02	1.51	1.53
5	3	614	CLA	CMC-C2C	-2.02	1.46	1.50
4	3	609	CHL	CHA-CBD	-2.02	1.49	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	1	601	CHL	MG-ND	-2.02	2.01	2.05
4	1	605	CHL	MG-ND	-2.02	2.01	2.05
5	2	612	CLA	MG-NB	-2.01	2.01	2.05
4	2	608	CHL	CMC-C2C	2.01	1.49	1.44
5	4	602	CLA	C1B-NB	-2.01	1.35	1.37
5	1	610	CLA	C4B-NB	2.00	1.40	1.37
9	4	2630	LHG	O7-C5	-2.00	1.41	1.46

All (798) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	3	601	CHL	C1B-CHB-C4A	9.98	127.74	121.32
4	3	609	CHL	O2D-CGD-CBD	9.88	121.80	110.95
4	1	606	CHL	C1B-CHB-C4A	9.87	127.67	121.32
4	4	606	CHL	C1B-CHB-C4A	9.69	127.56	121.32
4	1	607	CHL	C1B-CHB-C4A	9.57	127.48	121.32
4	3	606	CHL	O2D-CGD-CBD	9.51	121.39	110.95
4	4	608	CHL	C1B-CHB-C4A	9.43	127.39	121.32
4	2	609	CHL	O2D-CGD-CBD	9.25	121.11	110.95
4	3	601	CHL	O2D-CGD-CBD	9.00	120.84	110.95
4	1	608	CHL	C1B-CHB-C4A	8.80	126.98	121.32
4	2	608	CHL	C1B-CHB-C4A	8.78	126.97	121.32
4	2	606	CHL	C1B-CHB-C4A	8.74	126.94	121.32
4	1	609	CHL	O2D-CGD-CBD	8.73	120.54	110.95
4	2	605	CHL	O2D-CGD-CBD	8.36	120.13	110.95
7	1	1622	XAT	O24-C25-C24	8.29	121.26	113.49
4	4	609	CHL	C1B-CHB-C4A	8.25	126.63	121.32
4	3	608	CHL	C1B-CHB-C4A	8.23	126.62	121.32
4	1	606	CHL	O2D-CGD-CBD	8.19	119.94	110.95
4	4	609	CHL	O2D-CGD-CBD	8.05	119.80	110.95
4	2	601	CHL	C1B-CHB-C4A	8.03	126.49	121.32
4	4	607	CHL	O2D-CGD-CBD	8.02	119.76	110.95
4	1	601	CHL	C1B-CHB-C4A	7.83	126.36	121.32
4	3	605	CHL	O2D-CGD-CBD	7.82	119.53	110.95
7	2	1622	XAT	O4-C5-C4	7.79	120.79	113.49
10	4	623	BCR	C24-C23-C22	-7.77	114.74	126.23
4	1	605	CHL	O2D-CGD-CBD	7.71	119.42	110.95
8	2	1623	NEX	O24-C25-C24	7.66	120.67	113.49
4	3	607	CHL	O2D-CGD-CBD	7.63	119.33	110.95
4	2	606	CHL	O2D-CGD-CBD	7.61	119.31	110.95
4	1	601	CHL	O2D-CGD-CBD	7.49	119.18	110.95
8	1	1623	NEX	O24-C25-C24	7.47	120.49	113.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	2	1622	XAT	O24-C25-C24	7.44	120.47	113.49
4	1	608	CHL	O2D-CGD-CBD	7.38	119.06	110.95
4	3	605	CHL	C1B-CHB-C4A	7.35	126.05	121.32
8	3	1623	NEX	O24-C25-C24	7.32	120.35	113.49
4	3	606	CHL	C1B-CHB-C4A	7.15	125.92	121.32
4	1	607	CHL	O2D-CGD-CBD	7.10	118.74	110.95
4	3	608	CHL	O2D-CGD-CBD	7.02	118.66	110.95
7	3	1622	XAT	O4-C5-C4	6.93	119.99	113.49
4	2	607	CHL	O2D-CGD-CBD	6.91	118.54	110.95
4	4	606	CHL	O2D-CGD-CBD	6.91	118.53	110.95
4	2	608	CHL	O2D-CGD-CBD	6.87	118.49	110.95
4	4	608	CHL	O2D-CGD-CBD	6.82	118.44	110.95
4	2	601	CHL	O2D-CGD-CBD	6.77	118.38	110.95
4	3	608	CHL	C4D-ND-C1D	-6.69	100.14	105.22
4	4	608	CHL	C4D-ND-C1D	-6.59	100.22	105.22
4	3	601	CHL	C4D-ND-C1D	-6.55	100.25	105.22
4	2	608	CHL	C4D-ND-C1D	-6.34	100.41	105.22
5	1	611	CLA	C4A-NA-C1A	6.33	109.57	106.68
4	4	601	CHL	C1B-CHB-C4A	-6.32	117.26	121.32
4	1	608	CHL	C4D-ND-C1D	-6.30	100.44	105.22
6	3	1620	LUT	C7-C8-C9	-6.25	117.00	126.23
4	3	605	CHL	C4D-ND-C1D	-6.24	100.49	105.22
4	2	609	CHL	C1B-CHB-C4A	6.22	125.33	121.32
4	2	601	CHL	C4D-ND-C1D	-6.18	100.53	105.22
7	1	1622	XAT	O4-C5-C4	6.18	119.28	113.49
5	1	603	CLA	C4A-NA-C1A	6.17	109.49	106.68
4	1	607	CHL	C4D-ND-C1D	-6.06	100.62	105.22
4	3	601	CHL	CBC-CAC-C3C	-6.04	104.22	112.87
4	1	601	CHL	C4D-ND-C1D	-5.96	100.69	105.22
4	4	606	CHL	C4D-ND-C1D	-5.95	100.70	105.22
7	3	1622	XAT	C31-C30-C29	-5.92	118.98	127.28
4	1	605	CHL	C1B-CHB-C4A	5.85	125.08	121.32
7	4	622	XAT	C6-C7-C8	-5.84	113.64	125.99
4	2	601	CHL	CHA-C1A-C2A	-5.84	119.59	133.31
4	4	609	CHL	C4D-ND-C1D	-5.83	100.79	105.22
5	3	603	CLA	C4A-NA-C1A	5.83	109.34	106.68
4	2	609	CHL	C4D-ND-C1D	-5.80	100.82	105.22
7	3	1622	XAT	C27-C28-C29	-5.78	116.56	125.53
10	4	623	BCR	C28-C27-C26	-5.76	103.78	114.06
4	2	605	CHL	C4D-ND-C1D	-5.74	100.86	105.22
7	4	622	XAT	C38-C25-C26	-5.73	112.87	122.30
4	3	606	CHL	CHA-C1A-C2A	-5.69	119.94	133.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	3	601	CHL	CHA-C1A-C2A	-5.68	119.96	133.31
8	3	1623	NEX	C15-C14-C13	-5.62	119.40	127.28
8	1	1623	NEX	C27-C28-C29	-5.62	116.81	125.53
7	3	1622	XAT	C15-C14-C13	-5.61	119.41	127.28
4	1	608	CHL	CHA-C1A-C2A	-5.61	120.14	133.31
7	3	1622	XAT	C6-C7-C8	-5.60	114.15	125.99
5	3	611	CLA	C4A-NA-C1A	5.59	109.23	106.68
4	3	605	CHL	CHA-C1A-C2A	-5.59	120.19	133.31
4	1	601	CHL	CHA-C1A-C2A	-5.59	120.19	133.31
4	2	608	CHL	CHA-C1A-C2A	-5.58	120.20	133.31
4	3	608	CHL	CHA-C1A-C2A	-5.56	120.24	133.31
8	2	1623	NEX	C27-C28-C29	-5.51	116.98	125.53
4	4	608	CHL	CHA-C1A-C2A	-5.51	120.38	133.31
5	3	614	CLA	C4A-NA-C1A	5.50	109.19	106.68
4	1	605	CHL	CHA-C1A-C2A	-5.48	120.45	133.31
5	4	603	CLA	C4A-NA-C1A	5.47	109.18	106.68
4	2	606	CHL	CHA-C1A-C2A	-5.47	120.47	133.31
4	2	609	CHL	CHA-C1A-C2A	-5.45	120.50	133.31
4	4	606	CHL	CHA-C1A-C2A	-5.44	120.53	133.31
4	3	606	CHL	C4D-ND-C1D	-5.39	101.12	105.22
7	4	622	XAT	C18-C5-C6	-5.39	113.43	122.30
4	3	607	CHL	C4D-ND-C1D	-5.38	101.14	105.22
6	4	620	LUT	C21-C26-C27	-5.36	106.66	112.83
4	1	606	CHL	CHA-C1A-C2A	-5.34	120.77	133.31
4	4	607	CHL	C4D-ND-C1D	-5.33	101.17	105.22
5	2	611	CLA	C4A-NA-C1A	5.33	109.11	106.68
4	4	609	CHL	CHA-C1A-C2A	-5.28	120.90	133.31
4	1	609	CHL	C4D-ND-C1D	-5.27	101.22	105.22
4	2	606	CHL	C4D-ND-C1D	-5.26	101.22	105.22
4	1	607	CHL	CHA-C1A-C2A	-5.25	120.98	133.31
4	1	609	CHL	CHA-C1A-C2A	-5.24	121.01	133.31
4	3	607	CHL	C1B-CHB-C4A	5.24	124.69	121.32
5	3	604	CLA	C4A-NA-C1A	5.22	109.06	106.68
6	1	1621	LUT	C26-C27-C28	-5.21	116.47	124.58
4	1	606	CHL	C4D-ND-C1D	-5.20	101.27	105.22
7	4	622	XAT	C26-C27-C28	-5.16	115.09	125.99
4	4	607	CHL	CHA-C1A-C2A	-5.14	121.23	133.31
5	3	613	CLA	C4A-NA-C1A	5.13	109.02	106.68
8	3	1623	NEX	C38-C25-C26	-5.12	113.88	122.30
4	1	605	CHL	C4D-ND-C1D	-5.12	101.33	105.22
5	1	613	CLA	C4A-NA-C1A	5.12	109.01	106.68
7	4	622	XAT	O4-C5-C4	5.08	118.25	113.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	2	1620	LUT	C26-C27-C28	-5.07	116.68	124.58
7	3	1622	XAT	C38-C25-C26	-5.07	113.96	122.30
7	2	1622	XAT	C38-C25-C26	-5.06	113.97	122.30
4	3	609	CHL	CHA-C1A-C2A	-5.05	121.45	133.31
5	2	603	CLA	C4A-NA-C1A	5.05	108.98	106.68
10	4	623	BCR	C11-C10-C9	-5.04	120.21	127.28
10	4	623	BCR	C15-C14-C13	-5.03	120.23	127.28
7	4	622	XAT	O24-C25-C24	5.00	118.18	113.49
6	3	1620	LUT	C26-C27-C28	-5.00	116.80	124.58
7	2	1622	XAT	C15-C35-C34	-4.96	113.38	123.52
7	2	1622	XAT	C31-C30-C29	-4.94	120.35	127.28
6	4	620	LUT	C35-C34-C33	-4.92	120.38	127.28
6	1	1620	LUT	C35-C34-C33	-4.89	120.42	127.28
7	3	1622	XAT	O24-C25-C38	4.88	120.50	115.05
7	3	1622	XAT	C11-C10-C9	-4.88	120.44	127.28
7	1	1622	XAT	C15-C14-C13	-4.86	120.46	127.28
6	4	620	LUT	C7-C8-C9	-4.80	119.13	126.23
4	2	605	CHL	CHA-C1A-C2A	-4.78	122.08	133.31
8	2	1623	NEX	C15-C14-C13	-4.77	120.59	127.28
7	1	1622	XAT	C27-C28-C29	-4.76	118.15	125.53
8	2	1623	NEX	C38-C25-C26	-4.74	114.50	122.30
7	2	1622	XAT	C27-C28-C29	-4.74	118.17	125.53
6	4	620	LUT	C15-C14-C13	-4.72	120.65	127.28
4	3	607	CHL	CHA-C1A-C2A	-4.71	122.25	133.31
7	4	622	XAT	O24-C25-C38	4.71	120.31	115.05
7	2	1622	XAT	C7-C8-C9	-4.70	118.23	125.53
8	1	1623	NEX	C35-C34-C33	-4.66	120.74	127.28
4	3	609	CHL	C4D-ND-C1D	-4.66	101.68	105.22
5	1	612	CLA	C4A-NA-C1A	4.64	108.80	106.68
4	2	607	CHL	CHA-C1A-C2A	-4.60	122.50	133.31
9	4	2630	LHG	O4-P-O5	4.55	133.59	112.44
10	4	623	BCR	C16-C17-C18	-4.53	120.92	127.28
5	4	602	CLA	O2D-CGD-O1D	-4.52	115.05	123.85
6	1	1621	LUT	C15-C14-C13	-4.52	120.94	127.28
8	1	1623	NEX	C38-C25-C26	-4.51	114.89	122.30
7	1	1622	XAT	C31-C30-C29	-4.50	120.97	127.28
4	1	609	CHL	C1B-CHB-C4A	4.49	124.21	121.32
7	4	622	XAT	C35-C34-C33	-4.48	120.99	127.28
7	2	1622	XAT	C15-C14-C13	-4.48	121.00	127.28
7	1	1622	XAT	C38-C25-C26	-4.46	114.95	122.30
5	4	604	CLA	C4A-NA-C1A	4.46	108.71	106.68
7	2	1622	XAT	O24-C25-C38	4.44	120.00	115.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	2	2630	LHG	O4-P-O5	4.42	133.02	112.44
9	1	2630	LHG	O4-P-O5	4.38	132.80	112.44
6	3	1621	LUT	C10-C11-C12	-4.36	110.56	123.20
7	1	1622	XAT	C18-C5-C6	-4.35	115.14	122.30
4	4	608	CHL	CBC-CAC-C3C	-4.35	106.64	112.87
7	3	1622	XAT	O24-C25-C24	4.33	117.55	113.49
6	1	1620	LUT	C26-C27-C28	-4.33	117.84	124.58
7	3	1622	XAT	C4-C3-C2	-4.28	102.78	110.79
7	1	1622	XAT	C11-C10-C9	-4.27	121.29	127.28
5	4	612	CLA	C4A-NA-C1A	4.26	108.62	106.68
6	4	620	LUT	C26-C27-C28	-4.26	117.95	124.58
7	4	622	XAT	C15-C14-C13	-4.25	121.31	127.28
9	3	2630	LHG	O4-P-O5	4.23	132.11	112.44
8	1	1623	NEX	C17-C1-C6	-4.22	106.69	110.47
4	2	607	CHL	C4D-ND-C1D	-4.22	102.02	105.22
7	2	1622	XAT	C18-C5-C6	-4.18	115.42	122.30
4	4	607	CHL	C1B-CHB-C4A	4.18	124.01	121.32
4	2	609	CHL	CBC-CAC-C3C	-4.17	106.90	112.87
5	4	602	CLA	C4A-NA-C1A	4.15	108.57	106.68
6	2	1621	LUT	C35-C34-C33	-4.15	121.46	127.28
5	1	614	CLA	C4A-NA-C1A	4.13	108.56	106.68
7	4	622	XAT	C4-C3-C2	-4.12	103.09	110.79
6	2	1620	LUT	C15-C14-C13	-4.06	121.58	127.28
7	3	1622	XAT	C18-C5-C6	-4.06	115.62	122.30
5	4	611	CLA	C4A-NA-C1A	4.05	108.53	106.68
8	3	1623	NEX	O24-C25-C38	4.04	119.57	115.05
6	2	1621	LUT	C26-C27-C28	-4.04	118.29	124.58
8	3	1623	NEX	C11-C10-C9	-4.01	121.65	127.28
6	1	1621	LUT	C35-C34-C33	-4.00	121.67	127.28
5	1	604	CLA	C4A-NA-C1A	3.99	108.50	106.68
6	1	1620	LUT	C10-C11-C12	-3.98	111.66	123.20
8	2	1623	NEX	O24-C25-C38	3.98	119.49	115.05
5	3	602	CLA	C4A-NA-C1A	3.95	108.48	106.68
8	1	1623	NEX	O24-C25-C38	3.95	119.46	115.05
8	3	1623	NEX	C27-C28-C29	-3.93	119.43	125.53
7	3	1622	XAT	C15-C35-C34	-3.92	115.50	123.52
7	4	622	XAT	O4-C5-C18	3.88	119.39	115.05
4	1	605	CHL	OMC-CMC-C2C	-3.88	118.39	125.12
5	2	604	CLA	C4A-NA-C1A	3.87	108.44	106.68
6	4	620	LUT	C27-C28-C29	-3.87	118.00	126.32
4	1	607	CHL	CMA-C3A-C4A	-3.87	106.28	114.61
4	1	607	CHL	OMC-CMC-C2C	-3.82	118.49	125.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	4	601	CHL	C4D-ND-C1D	-3.80	102.33	105.22
4	2	609	CHL	O1D-CGD-CBD	-3.79	118.97	124.72
5	2	612	CLA	C4A-NA-C1A	3.79	108.41	106.68
4	2	605	CHL	C1B-CHB-C4A	3.78	123.75	121.32
6	3	1620	LUT	C35-C34-C33	-3.76	122.01	127.28
6	3	1620	LUT	C21-C26-C27	-3.76	108.51	112.83
4	3	606	CHL	C1C-CHC-C4B	3.75	129.47	116.07
7	3	1622	XAT	C35-C34-C33	-3.73	122.05	127.28
7	1	1622	XAT	C6-C7-C8	-3.73	118.12	125.99
7	2	1622	XAT	C11-C10-C9	-3.72	122.05	127.28
6	1	1620	LUT	C35-C15-C14	-3.72	115.91	123.52
5	2	614	CLA	O2D-CGD-O1D	-3.72	116.61	123.85
6	3	1620	LUT	C16-C1-C6	-3.70	104.44	110.24
7	4	622	XAT	C24-C23-C22	-3.67	103.92	110.79
7	1	1622	XAT	C4-C3-C2	-3.67	103.93	110.79
6	1	1620	LUT	C15-C14-C13	-3.66	122.14	127.28
7	3	1622	XAT	O4-C5-C18	3.65	119.13	115.05
4	4	609	CHL	OMC-CMC-C2C	-3.65	118.79	125.12
5	2	602	CLA	C4A-NA-C1A	3.63	108.33	106.68
5	4	602	CLA	O2D-CGD-CBD	3.62	117.56	111.23
4	3	609	CHL	O1D-CGD-CBD	-3.62	119.23	124.72
6	3	1621	LUT	C26-C27-C28	-3.61	118.96	124.58
8	1	1623	NEX	C15-C35-C34	-3.61	116.14	123.52
7	1	1622	XAT	O4-C5-C18	3.60	119.07	115.05
6	4	620	LUT	C31-C30-C29	-3.59	122.24	127.28
6	1	1621	LUT	C19-C9-C8	3.59	123.57	118.09
7	2	1622	XAT	O4-C5-C18	3.59	119.06	115.05
6	1	1620	LUT	C30-C31-C32	-3.58	112.84	123.20
4	1	608	CHL	C1C-CHC-C4B	3.56	128.82	116.07
4	3	601	CHL	O1D-CGD-CBD	-3.53	119.37	124.72
8	2	1623	NEX	C11-C10-C9	-3.50	122.37	127.28
5	3	612	CLA	O2D-CGD-O1D	-3.50	117.04	123.85
7	3	1622	XAT	C35-C15-C14	-3.49	116.39	123.52
4	2	605	CHL	O1D-CGD-CBD	-3.48	119.44	124.72
7	1	1622	XAT	O24-C25-C38	3.48	118.94	115.05
6	1	1621	LUT	C18-C5-C6	-3.45	120.72	124.48
4	3	605	CHL	C3D-C4D-ND	3.44	118.46	112.94
4	1	609	CHL	CMB-C2B-C3B	3.44	131.55	124.68
6	1	1621	LUT	C10-C11-C12	-3.44	113.25	123.20
6	3	1621	LUT	C7-C8-C9	-3.43	121.16	126.23
6	2	1620	LUT	C35-C34-C33	-3.43	122.47	127.28
7	3	1622	XAT	C30-C31-C32	-3.42	113.29	123.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	1	609	CHL	O1D-CGD-CBD	-3.42	119.54	124.72
5	2	612	CLA	CAA-C2A-C3A	-3.41	103.77	113.00
6	3	1620	LUT	C11-C10-C9	-3.41	122.49	127.28
10	4	623	BCR	C11-C12-C13	-3.41	117.02	126.36
4	3	606	CHL	O1D-CGD-CBD	-3.40	119.57	124.72
8	1	1623	NEX	C11-C10-C9	-3.38	122.54	127.28
7	4	622	XAT	C10-C11-C12	-3.38	113.42	123.20
6	3	1620	LUT	C30-C31-C32	-3.37	113.43	123.20
4	3	606	CHL	C3D-C4D-ND	3.37	118.34	112.94
5	2	613	CLA	C4A-NA-C1A	3.37	108.22	106.68
8	1	1623	NEX	C24-C23-C22	-3.36	104.50	110.79
8	2	1623	NEX	C24-C23-C22	-3.36	104.51	110.79
4	1	606	CHL	O1D-CGD-CBD	-3.36	119.63	124.72
5	3	603	CLA	CHB-C4A-NA	3.35	129.23	124.40
4	3	601	CHL	C3D-C4D-ND	3.35	118.31	112.94
4	4	608	CHL	C1C-CHC-C4B	3.34	128.01	116.07
5	4	612	CLA	CAA-C2A-C3A	-3.34	103.98	113.00
4	4	601	CHL	CHA-C1A-C2A	-3.33	125.48	133.31
4	4	609	CHL	O1D-CGD-CBD	-3.33	119.67	124.72
6	3	1621	LUT	C35-C34-C33	-3.32	122.62	127.28
4	3	601	CHL	CMA-C3A-C4A	-3.31	107.47	114.61
6	3	1620	LUT	C35-C15-C14	-3.31	116.74	123.52
5	3	603	CLA	CBC-CAC-C3C	-3.29	103.49	112.42
6	1	1621	LUT	C15-C35-C34	-3.29	116.79	123.52
4	1	609	CHL	CMA-C3A-C4A	-3.28	107.53	114.61
6	4	620	LUT	C11-C10-C9	-3.28	122.67	127.28
5	3	610	CLA	C3B-C4B-NB	-3.28	107.60	110.53
5	3	614	CLA	O2D-CGD-O1D	-3.28	117.47	123.85
6	3	1620	LUT	C15-C14-C13	-3.28	122.69	127.28
4	2	601	CHL	C1C-CHC-C4B	3.27	127.78	116.07
5	2	611	CLA	O2D-CGD-O1D	-3.27	117.49	123.85
4	3	608	CHL	C1C-CHC-C4B	3.25	127.70	116.07
4	1	607	CHL	C3D-C4D-ND	3.24	118.14	112.94
6	2	1621	LUT	C10-C11-C12	-3.22	113.86	123.20
6	1	1620	LUT	C31-C30-C29	-3.22	122.76	127.28
4	2	609	CHL	C3D-C4D-ND	3.21	118.09	112.94
6	2	1621	LUT	C30-C31-C32	-3.21	113.90	123.20
4	1	601	CHL	O1D-CGD-CBD	-3.21	119.86	124.72
4	3	607	CHL	C3D-C4D-ND	3.20	118.07	112.94
4	2	601	CHL	C3D-C4D-ND	3.20	118.07	112.94
4	1	608	CHL	O1D-CGD-CBD	-3.17	119.91	124.72
8	1	1623	NEX	C15-C14-C13	-3.17	122.84	127.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	2	605	CHL	C3D-C4D-ND	3.16	118.01	112.94
6	4	620	LUT	C30-C31-C32	-3.16	114.04	123.20
5	1	612	CLA	CAA-C2A-C3A	-3.16	104.46	113.00
5	2	603	CLA	CHB-C4A-NA	3.16	128.96	124.40
4	3	609	CHL	C1B-CHB-C4A	3.16	123.35	121.32
4	1	605	CHL	C3D-C4D-ND	3.15	118.00	112.94
4	4	608	CHL	C3D-C4D-ND	3.15	118.00	112.94
4	1	609	CHL	C3D-C4D-ND	3.15	117.99	112.94
5	2	603	CLA	CAA-C2A-C3A	-3.14	104.50	113.00
5	3	612	CLA	CAA-C2A-C3A	-3.14	104.50	113.00
4	3	609	CHL	C1C-CHC-C4B	3.14	127.30	116.07
7	1	1622	XAT	C35-C34-C33	-3.13	122.89	127.28
8	2	1623	NEX	C35-C34-C33	-3.12	122.91	127.28
8	3	1623	NEX	C11-C12-C13	-3.12	117.82	126.36
8	3	1623	NEX	C19-C9-C10	-3.12	117.77	122.82
4	4	608	CHL	O1D-CGD-CBD	-3.12	120.00	124.72
4	3	609	CHL	CBC-CAC-C3C	-3.11	108.42	112.87
6	3	1620	LUT	C10-C11-C12	-3.10	114.20	123.20
4	4	606	CHL	C3D-C4D-ND	3.10	117.92	112.94
4	2	608	CHL	C3D-C4D-ND	3.09	117.89	112.94
7	3	1622	XAT	C40-C33-C32	3.09	122.81	118.09
4	1	607	CHL	C1C-CHC-C4B	3.09	127.11	116.07
4	4	607	CHL	C3D-C4D-ND	3.08	117.88	112.94
4	4	609	CHL	C1C-CHC-C4B	3.08	127.08	116.07
7	1	1622	XAT	C15-C35-C34	-3.08	117.22	123.52
7	3	1622	XAT	C24-C23-C22	-3.08	105.04	110.79
4	1	601	CHL	C3D-C4D-ND	3.07	117.86	112.94
7	1	1622	XAT	C7-C8-C9	-3.07	120.77	125.53
4	3	607	CHL	OMC-CMC-C2C	-3.06	119.80	125.12
4	3	608	CHL	C3D-C4D-ND	3.06	117.85	112.94
5	1	604	CLA	O2D-CGD-O1D	-3.06	117.89	123.85
4	4	608	CHL	OMC-CMC-C2C	-3.05	119.82	125.12
6	2	1620	LUT	C21-C26-C27	-3.04	109.33	112.83
6	1	1620	LUT	C21-C26-C27	-3.04	109.33	112.83
6	4	620	LUT	C15-C35-C34	-3.04	117.31	123.52
7	3	1622	XAT	C7-C8-C9	-3.03	120.82	125.53
6	2	1620	LUT	C30-C31-C32	-3.03	114.42	123.20
4	2	608	CHL	C1C-CHC-C4B	3.03	126.91	116.07
6	4	620	LUT	C18-C5-C6	-3.03	121.18	124.48
4	4	606	CHL	O1D-CGD-CBD	-3.02	120.14	124.72
4	3	607	CHL	CMA-C3A-C4A	-3.02	108.10	114.61
4	4	609	CHL	C3D-C4D-ND	3.00	117.76	112.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	1	603	CLA	CHB-C4A-NA	3.00	128.73	124.40
5	4	604	CLA	CHB-C4A-NA	3.00	128.73	124.40
6	2	1620	LUT	C31-C30-C29	-3.00	123.07	127.28
4	2	609	CHL	C1C-CHC-C4B	3.00	126.80	116.07
5	3	602	CLA	CMB-C2B-C1B	-2.99	120.87	125.42
4	1	606	CHL	CMA-C3A-C4A	-2.99	108.18	114.61
5	2	604	CLA	O2D-CGD-O1D	-2.98	118.04	123.85
5	3	602	CLA	O2D-CGD-O1D	-2.98	118.05	123.85
6	2	1621	LUT	C15-C14-C13	-2.97	123.11	127.28
6	3	1621	LUT	C15-C14-C13	-2.96	123.12	127.28
7	2	1622	XAT	C40-C33-C32	2.96	122.61	118.09
4	1	609	CHL	C1C-CHC-C4B	2.96	126.66	116.07
4	1	608	CHL	C3D-C4D-ND	2.96	117.68	112.94
4	4	606	CHL	C1C-CHC-C4B	2.95	126.63	116.07
7	4	622	XAT	C31-C30-C29	-2.95	123.14	127.28
4	2	606	CHL	C3D-C4D-ND	2.95	117.67	112.94
4	2	601	CHL	CMD-C2D-C3D	-2.94	118.81	124.68
6	4	620	LUT	C18-C5-C4	2.94	119.83	114.42
6	2	1620	LUT	C10-C11-C12	-2.94	114.69	123.20
8	1	1623	NEX	C30-C31-C32	-2.93	114.71	123.20
10	4	623	BCR	C20-C21-C22	-2.93	123.17	127.28
4	2	605	CHL	C1C-CHC-C4B	2.93	126.54	116.07
4	4	607	CHL	C1C-CHC-C4B	2.92	126.53	116.07
4	1	609	CHL	O2A-CGA-CBA	2.92	120.74	111.83
5	3	613	CLA	CHB-C4A-NA	2.92	128.61	124.40
4	3	605	CHL	C1C-CHC-C4B	2.91	126.48	116.07
8	1	1623	NEX	C26-C27-C28	-2.91	119.84	125.99
5	1	614	CLA	CHB-C4A-NA	2.91	128.59	124.40
5	4	602	CLA	CHB-C4A-NA	2.90	128.58	124.40
7	3	1622	XAT	C38-C25-C24	2.89	117.49	114.24
4	1	605	CHL	CMA-C3A-C4A	-2.89	108.38	114.61
4	2	605	CHL	OMC-CMC-C2C	-2.89	120.10	125.12
8	1	1623	NEX	C39-C29-C30	-2.88	118.16	122.82
5	2	614	CLA	CHB-C4A-NA	2.87	128.55	124.40
4	2	609	CHL	O2A-CGA-CBA	2.87	120.58	111.83
5	2	612	CLA	CHB-C4A-NA	2.87	128.53	124.40
7	4	622	XAT	C18-C5-C4	2.86	117.45	114.24
7	3	1622	XAT	C10-C11-C12	-2.86	114.92	123.20
6	1	1620	LUT	C8-C9-C10	-2.86	114.52	119.01
5	2	613	CLA	O2D-CGD-O1D	-2.85	118.30	123.85
4	4	601	CHL	CMA-C3A-C4A	-2.85	108.47	114.61
4	2	606	CHL	C1C-CHC-C4B	2.85	126.27	116.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	2	1623	NEX	C39-C29-C30	-2.84	118.22	122.82
4	4	606	CHL	CBC-CAC-C3C	-2.84	108.81	112.87
8	1	1623	NEX	C31-C30-C29	-2.83	123.30	127.28
5	2	604	CLA	CHB-C4A-NA	2.83	128.49	124.40
8	3	1623	NEX	C39-C29-C30	-2.83	118.23	122.82
4	3	606	CHL	CMD-C2D-C3D	-2.83	119.03	124.68
4	3	605	CHL	O1D-CGD-CBD	-2.83	120.44	124.72
6	2	1620	LUT	C15-C35-C34	-2.82	117.74	123.52
4	3	609	CHL	C3D-C4D-ND	2.82	117.46	112.94
7	2	1622	XAT	C35-C34-C33	-2.82	123.32	127.28
5	4	603	CLA	CHB-C4A-NA	2.82	128.47	124.40
4	2	607	CHL	C3D-C4D-ND	2.82	117.45	112.94
8	1	1623	NEX	C11-C12-C13	-2.81	118.66	126.36
8	3	1623	NEX	C28-C29-C30	2.81	123.43	119.01
4	2	609	CHL	CMB-C2B-C3B	2.81	130.29	124.68
5	2	610	CLA	C3B-C4B-NB	-2.80	108.03	110.53
6	2	1621	LUT	C39-C29-C28	2.80	122.37	118.09
8	3	1623	NEX	C15-C35-C34	-2.80	117.79	123.52
4	4	606	CHL	CMA-C3A-C4A	-2.79	108.59	114.61
8	3	1623	NEX	C24-C23-C22	-2.79	105.58	110.79
5	1	603	CLA	CBC-CAC-C3C	-2.79	104.87	112.42
5	1	612	CLA	O2D-CGD-O1D	-2.78	118.43	123.85
5	4	603	CLA	CAA-C2A-C3A	-2.78	105.49	113.00
4	1	606	CHL	C3D-C4D-ND	2.78	117.39	112.94
4	4	607	CHL	O1D-CGD-CBD	-2.78	120.51	124.72
8	3	1623	NEX	C35-C34-C33	-2.78	123.38	127.28
4	2	606	CHL	O1D-CGD-CBD	-2.77	120.52	124.72
4	2	606	CHL	CMA-C3A-C4A	-2.77	108.63	114.61
6	1	1620	LUT	C19-C9-C8	2.77	122.32	118.09
4	2	607	CHL	C1C-CHC-C4B	2.77	125.98	116.07
4	1	605	CHL	O2D-CGD-O1D	-2.77	118.46	123.85
4	3	605	CHL	CBC-CAC-C3C	-2.77	108.91	112.87
5	2	602	CLA	C1-C2-C3	-2.76	121.67	126.20
6	3	1621	LUT	C30-C31-C32	-2.76	115.20	123.20
4	1	605	CHL	C1C-CHC-C4B	2.76	125.94	116.07
4	2	606	CHL	CBC-CAC-C3C	-2.76	108.93	112.87
6	3	1620	LUT	C22-C23-C24	2.75	115.27	111.18
4	1	601	CHL	C1C-CHC-C4B	2.75	125.91	116.07
5	2	610	CLA	O2D-CGD-O1D	-2.75	118.50	123.85
4	1	606	CHL	C1C-CHC-C4B	2.75	125.90	116.07
4	2	607	CHL	CAA-C2A-C3A	-2.74	105.58	113.00
6	2	1621	LUT	C22-C23-C24	-2.74	107.11	111.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	1	1621	LUT	C31-C30-C29	-2.74	123.44	127.28
5	1	612	CLA	CHB-C4A-NA	2.74	128.35	124.40
4	1	601	CHL	CMD-C2D-C3D	-2.74	119.21	124.68
4	4	601	CHL	CBC-CAC-C3C	-2.74	108.95	112.87
6	4	620	LUT	C35-C15-C14	-2.73	117.92	123.52
4	3	609	CHL	O2A-CGA-CBA	2.73	120.17	111.83
5	2	602	CLA	CMB-C2B-C1B	-2.73	121.27	125.42
4	2	601	CHL	O1D-CGD-CBD	-2.73	120.59	124.72
4	1	607	CHL	O2A-CGA-CBA	2.72	120.14	111.83
4	2	607	CHL	O1D-CGD-CBD	-2.72	120.59	124.72
6	1	1621	LUT	C30-C31-C32	-2.71	115.35	123.20
10	4	623	BCR	C1-C6-C5	-2.71	118.93	122.64
5	1	613	CLA	CHB-C4A-NA	2.71	128.31	124.40
4	1	609	CHL	C4-C3-C5	2.70	119.92	115.23
4	1	607	CHL	CAA-C2A-C3A	-2.70	105.69	113.00
4	2	606	CHL	CMB-C2B-C3B	2.70	130.08	124.68
5	3	603	CLA	CAA-C2A-C3A	-2.70	105.70	113.00
7	4	622	XAT	C15-C35-C34	-2.70	118.00	123.52
4	3	608	CHL	O1D-CGD-CBD	-2.70	120.64	124.72
4	2	608	CHL	O1D-CGD-CBD	-2.69	120.64	124.72
4	1	606	CHL	CBC-CAC-C3C	-2.69	109.02	112.87
5	4	612	CLA	CMB-C2B-C1B	-2.69	121.32	125.42
5	1	612	CLA	C2A-C1A-CHA	2.69	128.53	123.87
4	4	601	CHL	C1C-CHC-C4B	2.69	125.69	116.07
4	3	606	CHL	CMA-C3A-C4A	-2.69	108.82	114.61
5	2	602	CLA	O2D-CGD-O1D	-2.69	118.62	123.85
5	1	603	CLA	CAA-C2A-C3A	-2.68	105.75	113.00
5	1	611	CLA	CHB-C4A-NA	2.68	128.27	124.40
10	4	623	BCR	C7-C8-C9	-2.68	122.27	126.23
4	4	606	CHL	CMB-C2B-C3B	2.68	130.04	124.68
4	3	607	CHL	C1C-CHC-C4B	2.68	125.66	116.07
5	2	614	CLA	C4A-NA-C1A	2.68	107.90	106.68
5	1	602	CLA	O2D-CGD-O1D	-2.68	118.64	123.85
5	4	610	CLA	O2D-CGD-O1D	-2.67	118.64	123.85
5	2	603	CLA	O2D-CGD-O1D	-2.66	118.66	123.85
4	3	607	CHL	O2A-CGA-CBA	2.66	119.96	111.83
5	4	604	CLA	O2D-CGD-O1D	-2.66	118.67	123.85
6	1	1621	LUT	C8-C7-C6	-2.66	119.90	127.00
4	2	609	CHL	C1-C2-C3	-2.65	121.85	126.20
6	2	1621	LUT	C21-C26-C27	-2.65	109.78	112.83
4	1	607	CHL	O1D-CGD-CBD	-2.65	120.70	124.72
5	3	612	CLA	CHB-C4A-NA	2.65	128.23	124.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	1	608	CHL	CBC-CAC-C3C	-2.65	109.08	112.87
4	2	607	CHL	O2A-CGA-CBA	2.65	119.90	111.83
9	2	2630	LHG	O8-C23-C24	2.64	119.89	111.83
8	1	1623	NEX	C16-C1-C6	2.63	112.83	110.47
5	2	612	CLA	O2D-CGD-O1D	-2.63	118.72	123.85
5	3	611	CLA	O2D-CGD-O1D	-2.63	118.73	123.85
4	3	606	CHL	CMB-C2B-C3B	2.62	129.92	124.68
4	4	601	CHL	OMC-CMC-C2C	-2.61	120.58	125.12
7	3	1622	XAT	C19-C9-C8	2.61	122.08	118.09
7	2	1622	XAT	C10-C11-C12	-2.61	115.65	123.20
4	4	609	CHL	CMA-C3A-C4A	-2.61	108.99	114.61
4	2	607	CHL	CED-O2D-CGD	2.60	121.82	115.92
5	2	613	CLA	CHB-C4A-NA	2.60	128.15	124.40
4	3	609	CHL	CMA-C3A-C4A	-2.60	109.01	114.61
4	2	607	CHL	C1-C2-C3	-2.60	121.94	126.20
4	3	608	CHL	CMB-C2B-C3B	2.60	129.87	124.68
4	2	607	CHL	CBA-CAA-C2A	-2.60	106.13	113.78
5	4	610	CLA	C4A-NA-C1A	2.60	107.86	106.68
6	2	1620	LUT	C8-C7-C6	-2.59	120.07	127.00
6	4	620	LUT	C8-C7-C6	-2.59	120.08	127.00
6	3	1620	LUT	C31-C30-C29	-2.59	123.65	127.28
8	2	1623	NEX	C16-C1-C6	2.59	112.79	110.47
4	2	609	CHL	C4-C3-C5	2.59	119.72	115.23
5	2	611	CLA	O2D-CGD-CBD	2.58	115.75	111.23
5	1	610	CLA	C3B-C4B-NB	-2.58	108.22	110.53
5	4	602	CLA	CMB-C2B-C1B	-2.58	121.48	125.42
5	3	604	CLA	O2D-CGD-O1D	-2.58	118.82	123.85
4	3	605	CHL	OMC-CMC-C2C	-2.58	120.64	125.12
4	4	607	CHL	OMC-CMC-C2C	-2.58	120.64	125.12
4	2	609	CHL	CMA-C3A-C4A	-2.58	109.06	114.61
7	2	1622	XAT	C6-C7-C8	-2.57	120.55	125.99
4	3	607	CHL	O1D-CGD-CBD	-2.57	120.82	124.72
9	1	2630	LHG	O8-C23-C24	2.57	119.67	111.83
5	1	604	CLA	CHB-C4A-NA	2.57	128.10	124.40
4	1	601	CHL	CBC-CAC-C3C	-2.56	109.20	112.87
5	3	614	CLA	C3B-C4B-NB	-2.56	108.24	110.53
4	3	607	CHL	C4-C3-C5	2.56	119.67	115.23
5	4	611	CLA	O2D-CGD-O1D	-2.56	118.87	123.85
4	1	601	CHL	CMA-C3A-C4A	-2.55	109.11	114.61
5	4	603	CLA	O2D-CGD-O1D	-2.55	118.88	123.85
6	3	1620	LUT	C37-C21-C22	-2.55	104.65	109.41
5	3	604	CLA	CHB-C4A-NA	2.55	128.08	124.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	3	611	CLA	CHB-C4A-NA	2.55	128.08	124.40
9	3	2630	LHG	C11-C10-C9	-2.55	101.49	114.37
6	2	1620	LUT	C19-C9-C8	2.54	121.97	118.09
10	4	623	BCR	C27-C26-C25	-2.54	119.27	122.70
4	3	609	CHL	CMB-C2B-C3B	2.54	129.75	124.68
6	1	1620	LUT	C39-C29-C28	2.53	121.95	118.09
5	3	614	CLA	CHB-C4A-NA	2.53	128.04	124.40
4	3	609	CHL	O2D-CGD-O1D	-2.51	118.96	123.85
5	1	602	CLA	C3B-C4B-NB	-2.51	108.29	110.53
5	4	612	CLA	O2D-CGD-O1D	-2.51	118.97	123.85
10	4	623	BCR	C29-C30-C25	2.50	114.08	110.44
5	3	612	CLA	C4A-NA-C1A	2.50	107.82	106.68
5	2	604	CLA	O2A-CGA-O1A	-2.50	116.91	123.33
7	1	1622	XAT	C24-C23-C22	-2.50	106.12	110.79
5	4	612	CLA	CHB-C4A-NA	2.50	128.00	124.40
6	4	620	LUT	C1-C6-C5	-2.49	119.23	122.64
5	4	610	CLA	C3B-C4B-NB	-2.49	108.31	110.53
5	2	604	CLA	CMB-C2B-C1B	-2.48	121.64	125.42
5	1	611	CLA	O2D-CGD-O1D	-2.48	119.02	123.85
6	2	1621	LUT	C8-C7-C6	-2.48	120.38	127.00
5	1	613	CLA	O2D-CGD-O1D	-2.48	119.02	123.85
4	4	607	CHL	CMB-C2B-C3B	2.48	129.63	124.68
4	3	606	CHL	O2D-CGD-O1D	-2.48	119.03	123.85
6	2	1620	LUT	C35-C15-C14	-2.47	118.46	123.52
4	4	601	CHL	O2D-CGD-CBD	2.47	119.97	113.90
4	3	601	CHL	C1C-CHC-C4B	2.47	124.91	116.07
5	3	603	CLA	O2D-CGD-O1D	-2.47	119.04	123.85
5	3	602	CLA	CMB-C2B-C3B	2.47	132.36	126.55
6	2	1621	LUT	C35-C15-C14	-2.47	118.47	123.52
5	3	612	CLA	CMD-C2D-C1D	-2.46	120.39	124.73
5	4	611	CLA	C3B-C4B-NB	-2.46	108.33	110.53
5	2	602	CLA	CAA-CBA-CGA	-2.46	106.22	113.21
5	2	610	CLA	CHB-C4A-NA	2.46	127.95	124.40
4	2	606	CHL	CAA-C2A-C3A	-2.46	106.35	113.00
6	3	1620	LUT	C20-C13-C12	2.46	121.84	118.09
9	1	2630	LHG	C18-C17-C16	-2.46	101.96	114.37
4	3	609	CHL	OMC-CMC-C2C	-2.45	120.86	125.12
7	1	1622	XAT	C31-C32-C33	-2.45	119.64	126.36
5	4	602	CLA	C3B-C4B-NB	-2.45	108.34	110.53
5	3	610	CLA	C1-C2-C3	-2.45	122.18	126.20
9	3	2630	LHG	O8-C6-C5	-2.45	101.33	108.40
5	1	604	CLA	CHB-C1B-NB	2.45	127.72	124.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	1	1623	NEX	C2-C1-C6	2.45	111.59	109.21
5	4	612	CLA	CMB-C2B-C3B	2.45	132.31	126.55
4	1	607	CHL	CMB-C2B-C3B	2.44	129.57	124.68
6	3	1621	LUT	C15-C35-C34	-2.44	118.53	123.52
6	2	1621	LUT	C31-C30-C29	-2.44	123.86	127.28
5	1	610	CLA	O2D-CGD-O1D	-2.44	119.10	123.85
4	1	605	CHL	CMB-C2B-C3B	2.43	129.54	124.68
4	2	607	CHL	CMB-C2B-C3B	2.43	129.54	124.68
5	4	602	CLA	CMB-C2B-C3B	2.43	132.26	126.55
6	1	1621	LUT	C7-C8-C9	-2.43	122.65	126.23
5	2	602	CLA	C3B-C4B-NB	-2.42	108.37	110.53
5	1	604	CLA	C3B-C4B-NB	-2.42	108.37	110.53
4	4	607	CHL	CMA-C3A-C4A	-2.41	109.41	114.61
6	3	1620	LUT	C39-C29-C28	2.41	121.77	118.09
5	2	611	CLA	C3B-C4B-NB	-2.41	108.38	110.53
4	3	605	CHL	CAA-C2A-C3A	-2.41	106.49	113.00
5	3	603	CLA	C2A-C1A-CHA	2.41	128.05	123.87
5	1	602	CLA	C1-C2-C3	-2.41	122.25	126.20
5	1	604	CLA	C1-C2-C3	-2.41	122.87	126.76
5	1	610	CLA	CHB-C4A-NA	2.41	127.87	124.40
4	4	606	CHL	OMC-CMC-C2C	-2.41	120.94	125.12
5	3	610	CLA	O2D-CGD-O1D	-2.40	119.17	123.85
6	1	1621	LUT	O3-C3-C2	-2.40	104.83	109.75
5	2	602	CLA	CMB-C2B-C3B	2.40	132.20	126.55
9	1	2630	LHG	C20-C19-C18	-2.40	102.24	114.37
7	4	622	XAT	C19-C9-C8	2.40	121.75	118.09
4	4	601	CHL	CMB-C2B-C3B	2.40	129.47	124.68
6	2	1621	LUT	C27-C28-C29	-2.39	121.17	126.32
8	2	1623	NEX	C28-C29-C30	2.39	122.77	119.01
4	1	606	CHL	CMB-C2B-C3B	2.39	129.46	124.68
5	2	610	CLA	C4A-NA-C1A	2.38	107.77	106.68
5	1	614	CLA	CMB-C2B-C1B	-2.38	121.79	125.42
5	1	604	CLA	O2A-CGA-O1A	-2.38	117.68	123.63
7	2	1622	XAT	C4-C3-C2	-2.38	106.34	110.79
6	1	1620	LUT	C40-C33-C32	2.38	121.72	118.09
6	2	1621	LUT	C7-C8-C9	-2.38	122.72	126.23
8	2	1623	NEX	C8-C7-C6	-2.38	174.86	177.66
7	1	1622	XAT	C18-C5-C4	2.37	116.91	114.24
4	3	601	CHL	CMD-C2D-C3D	-2.37	119.94	124.68
6	4	620	LUT	C10-C11-C12	-2.37	116.32	123.20
9	2	2630	LHG	C11-C10-C9	-2.37	102.39	114.37
5	1	613	CLA	CHC-C4B-NB	2.37	127.60	124.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	2	612	CLA	C2A-C1A-CHA	2.37	127.98	123.87
6	2	1620	LUT	C39-C29-C28	2.37	121.70	118.09
7	4	622	XAT	C38-C25-C24	2.37	116.90	114.24
8	3	1623	NEX	C12-C13-C14	2.36	122.73	119.01
6	1	1621	LUT	C8-C9-C10	-2.36	115.29	119.01
4	3	606	CHL	CAC-C3C-C4C	2.36	130.09	124.03
5	2	602	CLA	CHB-C4A-NA	2.36	127.81	124.40
5	3	613	CLA	O2D-CGD-O1D	-2.36	119.26	123.85
6	3	1620	LUT	C18-C5-C6	-2.35	121.92	124.48
4	2	609	CHL	C6-C5-C3	-2.35	107.74	113.47
5	2	614	CLA	C3B-C4B-NB	-2.35	108.43	110.53
6	3	1621	LUT	C18-C5-C6	-2.35	121.92	124.48
5	4	610	CLA	CHB-C4A-NA	2.34	127.78	124.40
6	3	1621	LUT	C1-C2-C3	2.34	118.73	113.59
4	4	601	CHL	C3D-C4D-ND	2.34	116.69	112.94
5	1	610	CLA	O2A-CGA-O1A	-2.34	117.78	123.63
5	4	604	CLA	O2A-CGA-O1A	-2.34	117.32	123.33
5	1	612	CLA	CHA-C1A-NA	-2.34	121.10	126.39
5	1	612	CLA	CHC-C4B-NB	2.33	127.55	124.05
9	1	2630	LHG	C11-C10-C9	-2.33	102.59	114.37
4	1	607	CHL	CMD-C2D-C3D	-2.33	120.03	124.68
4	2	601	CHL	O2A-CGA-CBA	2.33	121.35	114.00
4	3	606	CHL	CBC-CAC-C3C	-2.32	109.55	112.87
7	3	1622	XAT	C32-C33-C34	-2.32	115.35	119.01
7	3	1622	XAT	C26-C27-C28	-2.32	121.08	125.99
4	4	607	CHL	CBC-CAC-C3C	-2.31	109.56	112.87
6	1	1620	LUT	C8-C7-C6	-2.31	120.83	127.00
5	2	613	CLA	CHB-C1B-NB	2.30	127.50	124.05
5	1	610	CLA	CMB-C2B-C3B	2.30	131.95	126.55
5	2	611	CLA	CHB-C4A-NA	2.30	127.71	124.40
5	4	612	CLA	C3B-C4B-NB	-2.30	108.48	110.53
7	2	1622	XAT	C24-C23-C22	-2.29	106.51	110.79
4	3	605	CHL	CMB-C2B-C3B	2.29	129.25	124.68
4	4	607	CHL	CAC-C3C-C4C	2.29	129.91	124.03
5	4	610	CLA	CMB-C2B-C1B	-2.29	121.94	125.42
6	2	1621	LUT	C36-C21-C26	2.29	113.01	109.55
4	2	605	CHL	CMB-C2B-C3B	2.28	129.25	124.68
4	3	601	CHL	CMB-C2B-C3B	2.28	129.25	124.68
5	3	611	CLA	CAA-C2A-C3A	-2.28	106.84	113.00
5	3	611	CLA	CAA-CBA-CGA	2.28	119.68	113.21
5	1	614	CLA	O2D-CGD-O1D	-2.28	119.42	123.85
5	3	602	CLA	C1-C2-C3	-2.27	122.48	126.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	3	1620	LUT	C27-C28-C29	-2.27	121.44	126.32
5	1	603	CLA	C2A-C1A-CHA	2.27	127.80	123.87
4	2	601	CHL	CED-O2D-CGD	2.27	121.06	115.92
4	2	608	CHL	CMB-C2B-C3B	2.26	129.19	124.68
8	2	1623	NEX	C30-C31-C32	-2.25	116.67	123.20
6	3	1620	LUT	C17-C1-C6	2.25	113.77	110.24
5	1	612	CLA	O2D-CGD-CBD	2.25	115.16	111.23
10	4	623	BCR	C21-C20-C19	-2.25	116.69	123.20
4	1	601	CHL	C1C-C2C-CMC	2.25	130.85	126.80
5	1	610	CLA	CHB-C1B-NB	2.25	127.42	124.05
5	1	613	CLA	CHB-C1B-NB	2.25	127.42	124.05
4	4	609	CHL	CMD-C2D-C3D	-2.25	120.19	124.68
4	1	609	CHL	CBC-CAC-C3C	-2.24	109.66	112.87
5	4	611	CLA	CHB-C4A-NA	2.24	127.64	124.40
8	3	1623	NEX	C30-C31-C32	-2.24	116.71	123.20
7	4	622	XAT	C8-C9-C10	-2.24	115.48	119.01
4	3	608	CHL	CMA-C3A-C4A	-2.24	109.79	114.61
4	3	607	CHL	CAC-C3C-C4C	2.23	129.77	124.03
4	1	608	CHL	OBD-CAD-C3D	-2.23	124.41	127.89
5	3	610	CLA	O2A-CGA-O1A	-2.23	118.05	123.63
4	3	601	CHL	C4-C3-C5	2.23	119.10	115.23
4	4	608	CHL	O2A-CGA-CBA	2.23	121.04	114.00
4	4	606	CHL	O2A-CGA-CBA	2.23	121.04	114.00
5	2	604	CLA	O1A-CGA-CBA	2.23	130.15	123.09
5	3	613	CLA	C1-C2-C3	-2.22	122.55	126.20
4	4	609	CHL	CMB-C2B-C3B	2.22	129.13	124.68
6	1	1620	LUT	C20-C13-C12	2.22	121.47	118.09
5	1	613	CLA	CMB-C2B-C1B	-2.22	122.05	125.42
4	4	608	CHL	CMA-C3A-C2A	-2.21	105.59	114.13
4	2	607	CHL	C4-C3-C5	2.21	119.06	115.23
4	4	601	CHL	C3D-C4D-CHA	2.21	111.89	108.54
5	1	603	CLA	O2D-CGD-O1D	-2.20	119.56	123.85
5	4	611	CLA	O2A-CGA-O1A	-2.20	117.67	123.33
4	3	608	CHL	O2A-CGA-CBA	2.20	120.96	114.00
5	3	602	CLA	CHB-C4A-NA	2.20	127.58	124.40
6	3	1621	LUT	C38-C25-C24	-2.20	118.15	123.36
6	2	1621	LUT	C16-C1-C6	-2.19	106.80	110.24
4	3	601	CHL	O2A-CGA-CBA	2.19	118.53	111.83
4	4	608	CHL	CMA-C3A-C4A	-2.19	109.89	114.61
4	3	609	CHL	C4-C3-C5	2.19	119.03	115.23
9	3	2630	LHG	C18-C17-C16	-2.19	103.28	114.37
4	1	608	CHL	CMB-C2B-C3B	2.19	129.06	124.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	1	607	CHL	C1-C2-C3	-2.19	122.61	126.20
5	4	603	CLA	CMB-C2B-C1B	-2.19	122.08	125.42
5	1	614	CLA	C3B-C4B-NB	-2.19	108.58	110.53
5	3	612	CLA	CMD-C2D-C3D	2.19	132.70	127.69
5	3	610	CLA	CHB-C1B-NB	2.18	127.33	124.05
10	4	623	BCR	C3-C4-C5	-2.18	110.17	114.06
4	2	606	CHL	CED-O2D-CGD	2.17	120.85	115.92
5	2	604	CLA	CHB-C1B-NB	2.17	127.31	124.05
4	3	601	CHL	C1C-C2C-CMC	2.17	130.72	126.80
5	3	603	CLA	CHA-C1A-NA	-2.17	121.47	126.39
5	3	614	CLA	CMB-C2B-C1B	-2.17	122.11	125.42
4	4	609	CHL	O2A-CGA-CBA	2.17	120.85	114.00
4	2	609	CHL	CAC-C3C-C4C	2.17	129.60	124.03
5	1	602	CLA	CHA-C1A-NA	-2.17	121.48	126.39
4	3	607	CHL	CMB-C2B-C3B	2.17	129.01	124.68
7	2	1622	XAT	C16-C1-C2	-2.16	105.16	108.97
5	1	602	CLA	CHB-C4A-NA	2.16	127.52	124.40
4	2	609	CHL	OMC-CMC-C2C	-2.16	121.36	125.12
4	4	608	CHL	O2A-CGA-O1A	-2.16	117.78	123.33
7	1	1622	XAT	C10-C11-C12	-2.16	116.94	123.20
4	4	609	CHL	CAC-C3C-C4C	2.16	129.58	124.03
6	2	1620	LUT	C8-C9-C10	-2.16	115.61	119.01
6	3	1620	LUT	C38-C25-C24	-2.16	118.26	123.36
6	3	1620	LUT	C15-C35-C34	-2.16	119.11	123.52
6	3	1621	LUT	C39-C29-C28	2.15	121.38	118.09
10	4	623	BCR	C4-C5-C6	-2.15	119.80	122.70
9	2	2630	LHG	C27-C26-C25	-2.15	103.50	114.37
4	3	601	CHL	C4C-CHD-C1D	2.15	123.76	116.07
4	2	601	CHL	CMB-C2B-C3B	2.15	128.98	124.68
4	3	608	CHL	C4C-CHD-C1D	2.15	123.76	116.07
5	2	604	CLA	CMB-C2B-C3B	2.15	131.60	126.55
4	1	601	CHL	C4C-CHD-C1D	2.15	123.75	116.07
5	1	614	CLA	CMB-C2B-C3B	2.15	131.60	126.55
6	1	1621	LUT	C31-C32-C33	-2.14	120.48	126.36
5	4	610	CLA	CHB-C1B-NB	2.14	127.27	124.05
7	4	622	XAT	C30-C31-C32	-2.14	116.99	123.20
6	2	1621	LUT	C11-C10-C9	-2.14	124.28	127.28
5	4	610	CLA	CMB-C2B-C3B	2.14	131.58	126.55
4	2	601	CHL	C4C-CHD-C1D	2.14	123.71	116.07
9	1	2630	LHG	C27-C26-C25	-2.14	103.57	114.37
4	1	607	CHL	CAC-C3C-C4C	2.13	129.52	124.03
4	4	606	CHL	C4C-CHD-C1D	2.13	123.70	116.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	1	603	CLA	CHA-C1A-NA	-2.13	121.56	126.39
5	2	603	CLA	O2A-CGA-O1A	-2.13	118.30	123.63
4	2	601	CHL	CMA-C3A-C4A	-2.13	110.02	114.61
4	1	606	CHL	O2A-CGA-CBA	2.13	120.73	114.00
4	1	601	CHL	CAC-C3C-C4C	2.12	129.49	124.03
4	4	607	CHL	CAA-CBA-CGA	-2.12	106.83	112.49
5	2	613	CLA	O1D-CGD-CBD	2.12	128.70	124.52
4	4	601	CHL	CAC-C3C-C4C	2.12	129.49	124.03
4	1	609	CHL	C1-C2-C3	-2.12	122.72	126.20
4	3	601	CHL	C1-O2A-CGA	2.12	121.78	116.65
6	4	620	LUT	C20-C13-C12	2.12	121.33	118.09
4	1	607	CHL	CBA-CAA-C2A	-2.12	107.54	113.78
4	4	607	CHL	O2D-CGD-O1D	-2.12	119.73	123.85
4	2	607	CHL	CAC-C3C-C4C	2.12	129.47	124.03
4	2	605	CHL	CMA-C3A-C4A	-2.11	110.06	114.61
5	3	610	CLA	CHB-C4A-NA	2.11	127.45	124.40
5	1	602	CLA	CMB-C2B-C3B	2.11	131.51	126.55
4	4	608	CHL	OBD-CAD-C3D	-2.11	124.60	127.89
6	2	1621	LUT	C3-C4-C5	-2.11	106.93	112.18
4	1	601	CHL	CMB-C2B-C3B	2.11	128.89	124.68
7	2	1622	XAT	O24-C25-C26	-2.10	57.26	58.93
4	1	607	CHL	C1C-C2C-CMC	-2.10	123.02	126.80
7	3	1622	XAT	C11-C12-C13	-2.10	120.61	126.36
5	1	602	CLA	C4A-NA-C1A	2.10	107.64	106.68
4	3	605	CHL	C4C-CHD-C1D	2.09	123.56	116.07
5	1	614	CLA	O2A-CGA-O1A	-2.09	117.95	123.33
5	4	604	CLA	C3B-C4B-NB	-2.09	108.66	110.53
5	2	612	CLA	CHB-C1B-C2B	-2.09	121.36	127.43
4	3	601	CHL	O2D-CGD-O1D	-2.09	119.78	123.85
5	2	604	CLA	CAC-C3C-C4C	2.09	127.50	124.79
7	1	1622	XAT	O24-C25-C26	-2.09	57.28	58.93
5	1	610	CLA	CHB-C1B-C2B	-2.09	121.37	127.43
5	3	602	CLA	O1D-CGD-CBD	2.08	128.63	124.52
4	2	605	CHL	CAA-C2A-C3A	-2.08	107.37	113.00
5	2	612	CLA	CHB-C1B-NB	2.08	127.17	124.05
5	2	612	CLA	CHA-C1A-NA	-2.08	121.68	126.39
5	3	603	CLA	CHC-C4B-NB	2.08	127.17	124.05
4	3	607	CHL	O2D-CGD-O1D	-2.08	119.80	123.85
5	2	614	CLA	O2D-CGD-CBD	2.08	114.87	111.23
5	2	603	CLA	C2A-C1A-CHA	2.08	127.47	123.87
5	1	610	CLA	C4A-NA-C1A	2.08	107.63	106.68
4	1	607	CHL	CAA-CBA-CGA	-2.07	107.32	113.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	4	2630	LHG	C5-O7-C7	-2.07	114.19	117.85
8	3	1623	NEX	C20-C13-C14	-2.07	119.46	122.82
10	4	623	BCR	C15-C16-C17	-2.07	119.28	123.52
4	3	601	CHL	C1-C2-C3	-2.07	122.81	126.20
5	1	610	CLA	CMB-C2B-C1B	-2.07	122.27	125.42
4	2	606	CHL	O2A-CGA-CBA	2.07	120.54	114.00
5	2	613	CLA	CHB-C1B-C2B	-2.07	121.42	127.43
7	2	1622	XAT	C20-C13-C12	2.07	121.24	118.09
5	3	614	CLA	O2D-CGD-CBD	2.06	114.84	111.23
5	3	611	CLA	CHC-C4B-NB	2.06	127.14	124.05
4	3	608	CHL	OBD-CAD-C3D	-2.06	124.67	127.89
5	2	604	CLA	CHC-C4B-NB	2.06	127.14	124.05
5	3	611	CLA	O2A-CGA-O1A	-2.06	118.47	123.63
4	4	609	CHL	CED-O2D-CGD	2.06	120.59	115.92
7	4	622	XAT	C39-C29-C28	2.06	121.23	118.09
4	1	607	CHL	C4-C3-C5	2.05	118.79	115.23
4	3	609	CHL	CMD-C2D-C3D	-2.05	120.58	124.68
4	2	601	CHL	C1C-C2C-CMC	2.05	130.50	126.80
5	3	610	CLA	C7-C6-C5	-2.05	107.80	113.26
4	2	605	CHL	CBC-CAC-C3C	-2.05	109.94	112.87
4	2	608	CHL	C4C-CHD-C1D	2.04	123.39	116.07
4	2	601	CHL	CBC-CAC-C3C	-2.04	109.95	112.87
4	3	605	CHL	CGD-CBD-CAD	-2.04	104.23	110.85
6	3	1621	LUT	C31-C30-C29	-2.04	124.41	127.28
4	3	607	CHL	C1-C2-C3	-2.04	122.85	126.20
8	1	1623	NEX	C19-C9-C10	-2.04	119.51	122.82
5	3	604	CLA	O1A-CGA-CBA	2.04	129.56	123.09
4	2	601	CHL	CAC-C3C-C4C	2.04	129.27	124.03
4	2	609	CHL	O2D-CGD-O1D	-2.03	119.89	123.85
8	2	1623	NEX	C15-C35-C34	-2.03	119.36	123.52
7	2	1622	XAT	C31-C32-C33	-2.03	120.79	126.36
5	1	612	CLA	CHC-C1C-NC	2.03	127.37	124.31
5	2	614	CLA	O1D-CGD-CBD	2.03	128.52	124.52
5	4	611	CLA	CHC-C4B-NB	2.03	127.09	124.05
5	1	613	CLA	CHB-C1B-C2B	-2.03	121.54	127.43
5	2	611	CLA	CHB-C1B-NB	2.03	127.09	124.05
4	1	609	CHL	O2D-CGD-O1D	-2.03	119.90	123.85
5	2	614	CLA	CMB-C2B-C1B	-2.03	122.33	125.42
4	1	608	CHL	O2A-CGA-CBA	2.03	120.40	114.00
6	2	1620	LUT	C11-C10-C9	-2.02	124.44	127.28
5	1	611	CLA	O2A-CGA-O1A	-2.02	118.13	123.33
4	3	608	CHL	CAC-C3C-C4C	2.02	129.22	124.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	1	613	CLA	O2D-CGD-CBD	2.02	114.75	111.23
6	2	1621	LUT	C18-C5-C6	-2.01	122.29	124.48
6	1	1621	LUT	C36-C21-C26	2.01	112.59	109.55
4	2	608	CHL	O2A-CGA-CBA	2.01	120.35	114.00
7	3	1622	XAT	C39-C29-C30	-2.01	119.56	122.82
5	4	602	CLA	CHB-C1B-C2B	-2.01	121.59	127.43
6	3	1620	LUT	C36-C21-C26	2.01	112.59	109.55
4	4	608	CHL	CED-O2D-CGD	2.01	120.47	115.92
5	3	611	CLA	CHB-C1B-NB	2.01	127.06	124.05
4	1	608	CHL	C4C-CHD-C1D	2.00	123.23	116.07
5	4	604	CLA	O1A-CGA-CBA	2.00	129.44	123.09

All (100) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	1	601	CHL	NA
4	1	601	CHL	NC
4	1	601	CHL	ND
4	1	605	CHL	NA
4	1	605	CHL	NC
4	1	605	CHL	ND
4	1	606	CHL	NA
4	1	606	CHL	NC
4	1	606	CHL	ND
4	1	607	CHL	NA
4	1	607	CHL	C8
4	1	607	CHL	NC
4	1	607	CHL	ND
4	1	608	CHL	NA
4	1	608	CHL	NC
4	1	608	CHL	ND
4	1	609	CHL	NA
4	1	609	CHL	C8
4	1	609	CHL	NC
4	1	609	CHL	ND
4	2	601	CHL	NA
4	2	601	CHL	NC
4	2	601	CHL	ND
4	2	605	CHL	NA
4	2	605	CHL	NC
4	2	605	CHL	ND
4	2	606	CHL	NA

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Mol	Chain	Res	Type	Atom
4	2	606	CHL	NC
4	2	606	CHL	ND
4	2	607	CHL	NA
4	2	607	CHL	C8
4	2	607	CHL	NC
4	2	607	CHL	ND
4	2	608	CHL	NA
4	2	608	CHL	NC
4	2	608	CHL	ND
4	2	609	CHL	NA
4	2	609	CHL	C8
4	2	609	CHL	NC
4	2	609	CHL	ND
4	3	601	CHL	NA
4	3	601	CHL	C8
4	3	601	CHL	NC
4	3	601	CHL	ND
4	3	605	CHL	NA
4	3	605	CHL	NC
4	3	605	CHL	ND
4	3	606	CHL	NA
4	3	606	CHL	NC
4	3	606	CHL	ND
4	3	607	CHL	NA
4	3	607	CHL	NC
4	3	607	CHL	ND
4	3	608	CHL	NA
4	3	608	CHL	NC
4	3	608	CHL	ND
4	3	609	CHL	NA
4	3	609	CHL	C8
4	3	609	CHL	NC
4	3	609	CHL	ND
4	4	601	CHL	NA
4	4	601	CHL	NC
4	4	601	CHL	ND
4	4	606	CHL	NA
4	4	606	CHL	NC
4	4	606	CHL	ND
4	4	607	CHL	NA
4	4	607	CHL	NC
4	4	607	CHL	ND

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Mol	Chain	Res	Type	Atom
4	4	608	CHL	NA
4	4	608	CHL	NC
4	4	608	CHL	ND
4	4	609	CHL	NA
4	4	609	CHL	NC
4	4	609	CHL	ND
5	1	602	CLA	ND
5	1	603	CLA	ND
5	1	604	CLA	ND
5	1	610	CLA	ND
5	1	611	CLA	ND
5	1	612	CLA	ND
5	1	614	CLA	ND
5	2	602	CLA	ND
5	2	603	CLA	ND
5	2	610	CLA	ND
5	2	611	CLA	ND
5	2	612	CLA	ND
5	3	602	CLA	ND
5	3	603	CLA	ND
5	3	604	CLA	ND
5	3	610	CLA	ND
5	3	611	CLA	ND
5	3	612	CLA	ND
5	3	613	CLA	ND
5	3	614	CLA	ND
5	4	602	CLA	ND
5	4	603	CLA	ND
5	4	610	CLA	ND
5	4	611	CLA	ND
5	4	612	CLA	ND

All (555) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	1	605	CHL	C1A-C2A-CAA-CBA
4	1	605	CHL	C1C-C2C-CMC-OMC
4	1	605	CHL	C3C-C2C-CMC-OMC
4	1	606	CHL	C3C-C2C-CMC-OMC
4	1	606	CHL	CBD-CGD-O2D-CED
4	1	607	CHL	C1C-C2C-CMC-OMC
4	1	607	CHL	C3C-C2C-CMC-OMC

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Mol	Chain	Res	Type	Atoms
4	1	607	CHL	C4C-C3C-CAC-CBC
4	1	608	CHL	C1A-C2A-CAA-CBA
4	1	609	CHL	C1C-C2C-CMC-OMC
4	1	609	CHL	C3C-C2C-CMC-OMC
4	1	609	CHL	C4C-C3C-CAC-CBC
4	1	609	CHL	CHA-CBD-CGD-O1D
4	2	601	CHL	C3A-C2A-CAA-CBA
4	2	605	CHL	C1A-C2A-CAA-CBA
4	2	605	CHL	C1C-C2C-CMC-OMC
4	2	605	CHL	C3C-C2C-CMC-OMC
4	2	606	CHL	C3C-C2C-CMC-OMC
4	2	607	CHL	C1C-C2C-CMC-OMC
4	2	607	CHL	C3C-C2C-CMC-OMC
4	2	607	CHL	C4C-C3C-CAC-CBC
4	2	607	CHL	CHA-CBD-CGD-O1D
4	2	607	CHL	CHA-CBD-CGD-O2D
4	2	608	CHL	C1C-C2C-CMC-OMC
4	2	608	CHL	C3C-C2C-CMC-OMC
4	2	608	CHL	C4C-C3C-CAC-CBC
4	2	609	CHL	C1C-C2C-CMC-OMC
4	2	609	CHL	C3C-C2C-CMC-OMC
4	3	605	CHL	C1A-C2A-CAA-CBA
4	3	605	CHL	C1C-C2C-CMC-OMC
4	3	605	CHL	C3C-C2C-CMC-OMC
4	3	606	CHL	C3A-C2A-CAA-CBA
4	3	607	CHL	C1C-C2C-CMC-OMC
4	3	607	CHL	C3C-C2C-CMC-OMC
4	3	609	CHL	C1C-C2C-CMC-OMC
4	3	609	CHL	C3C-C2C-CMC-OMC
4	3	609	CHL	C4C-C3C-CAC-CBC
4	4	601	CHL	C1C-C2C-CMC-OMC
4	4	601	CHL	C3C-C2C-CMC-OMC
4	4	606	CHL	C3A-C2A-CAA-CBA
4	4	606	CHL	C1C-C2C-CMC-OMC
4	4	607	CHL	C1C-C2C-CMC-OMC
4	4	607	CHL	C3C-C2C-CMC-OMC
4	4	608	CHL	C1A-C2A-CAA-CBA
4	4	608	CHL	C1C-C2C-CMC-OMC
4	4	608	CHL	C3C-C2C-CMC-OMC
4	4	608	CHL	CHA-CBD-CGD-O1D
4	4	608	CHL	CHA-CBD-CGD-O2D
4	4	608	CHL	CBD-CGD-O2D-CED

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Mol	Chain	Res	Type	Atoms
4	4	609	CHL	C1C-C2C-CMC-OMC
4	4	609	CHL	C3C-C2C-CMC-OMC
4	4	609	CHL	C4C-C3C-CAC-CBC
5	1	603	CLA	C4B-C3B-CAB-CBB
5	1	603	CLA	CBD-CGD-O2D-CED
5	1	604	CLA	CAD-CBD-CGD-O1D
5	1	604	CLA	CAD-CBD-CGD-O2D
5	1	611	CLA	C2B-C3B-CAB-CBB
5	1	611	CLA	C4B-C3B-CAB-CBB
5	1	612	CLA	CHA-CBD-CGD-O1D
5	1	612	CLA	CHA-CBD-CGD-O2D
5	1	613	CLA	CBD-CGD-O2D-CED
5	1	614	CLA	CBD-CGD-O2D-CED
5	2	603	CLA	C2B-C3B-CAB-CBB
5	2	603	CLA	C4B-C3B-CAB-CBB
5	2	603	CLA	CBD-CGD-O2D-CED
5	2	604	CLA	C1A-C2A-CAA-CBA
5	2	604	CLA	C3A-C2A-CAA-CBA
5	2	604	CLA	CAD-CBD-CGD-O1D
5	2	604	CLA	CAD-CBD-CGD-O2D
5	2	604	CLA	CBD-CGD-O2D-CED
5	2	610	CLA	C2B-C3B-CAB-CBB
5	2	610	CLA	C4B-C3B-CAB-CBB
5	2	611	CLA	C2B-C3B-CAB-CBB
5	2	611	CLA	C4B-C3B-CAB-CBB
5	2	612	CLA	CHA-CBD-CGD-O1D
5	2	612	CLA	CHA-CBD-CGD-O2D
5	2	613	CLA	CBD-CGD-O2D-CED
5	2	614	CLA	C1A-C2A-CAA-CBA
5	2	614	CLA	C3A-C2A-CAA-CBA
5	3	602	CLA	CHA-CBD-CGD-O2D
5	3	603	CLA	O1A-CGA-O2A-C1
5	3	603	CLA	C4B-C3B-CAB-CBB
5	3	610	CLA	C4B-C3B-CAB-CBB
5	3	611	CLA	C1A-C2A-CAA-CBA
5	3	611	CLA	C4B-C3B-CAB-CBB
5	3	612	CLA	C2B-C3B-CAB-CBB
5	3	612	CLA	C4B-C3B-CAB-CBB
5	3	612	CLA	CAD-CBD-CGD-O1D
5	3	612	CLA	CAD-CBD-CGD-O2D
5	3	613	CLA	C2B-C3B-CAB-CBB
5	3	613	CLA	C4B-C3B-CAB-CBB

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Mol	Chain	Res	Type	Atoms
5	3	614	CLA	CAD-CBD-CGD-O1D
5	3	614	CLA	CAD-CBD-CGD-O2D
5	3	614	CLA	CBD-CGD-O2D-CED
5	4	602	CLA	CBD-CGD-O2D-CED
5	4	604	CLA	C1A-C2A-CAA-CBA
5	4	610	CLA	C1A-C2A-CAA-CBA
5	4	611	CLA	C2B-C3B-CAB-CBB
5	4	611	CLA	C4B-C3B-CAB-CBB
5	4	611	CLA	CBD-CGD-O2D-CED
6	4	620	LUT	C11-C12-C13-C14
6	4	620	LUT	C11-C12-C13-C20
7	1	1622	XAT	C31-C32-C33-C34
7	1	1622	XAT	C31-C32-C33-C40
7	2	1622	XAT	O4-C6-C7-C8
7	2	1622	XAT	C31-C32-C33-C34
8	1	1623	NEX	C11-C12-C13-C14
8	1	1623	NEX	C11-C12-C13-C20
9	2	2630	LHG	C4-O6-P-O5
9	3	2630	LHG	O1-C1-C2-C3
9	3	2630	LHG	C4-O6-P-O3
9	3	2630	LHG	C4-O6-P-O5
9	4	2630	LHG	O1-C1-C2-O2
9	4	2630	LHG	C3-O3-P-O5
9	4	2630	LHG	C3-O3-P-O6
10	4	623	BCR	C11-C12-C13-C14
10	4	623	BCR	C11-C12-C13-C35
10	4	623	BCR	C21-C22-C23-C24
10	4	623	BCR	C37-C22-C23-C24
5	2	612	CLA	O1D-CGD-O2D-CED
5	4	604	CLA	O1D-CGD-O2D-CED
4	1	608	CHL	O1D-CGD-O2D-CED
5	2	613	CLA	O1D-CGD-O2D-CED
5	4	611	CLA	O1D-CGD-O2D-CED
9	4	2630	LHG	C24-C23-O8-C6
4	1	608	CHL	CBD-CGD-O2D-CED
4	2	606	CHL	CBD-CGD-O2D-CED
4	2	608	CHL	CBD-CGD-O2D-CED
4	3	607	CHL	CBD-CGD-O2D-CED
4	3	608	CHL	CBD-CGD-O2D-CED
4	4	606	CHL	CBD-CGD-O2D-CED
5	2	602	CLA	CBD-CGD-O2D-CED
5	2	612	CLA	CBD-CGD-O2D-CED

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Mol	Chain	Res	Type	Atoms
5	2	614	CLA	CBD-CGD-O2D-CED
5	3	603	CLA	CBD-CGD-O2D-CED
5	3	612	CLA	CBD-CGD-O2D-CED
5	3	613	CLA	CBD-CGD-O2D-CED
5	4	604	CLA	CBD-CGD-O2D-CED
4	3	608	CHL	O1D-CGD-O2D-CED
4	4	606	CHL	O1D-CGD-O2D-CED
5	3	613	CLA	O1D-CGD-O2D-CED
5	3	603	CLA	CBA-CGA-O2A-C1
4	1	606	CHL	O1D-CGD-O2D-CED
5	1	613	CLA	O1D-CGD-O2D-CED
5	2	603	CLA	O1D-CGD-O2D-CED
5	4	602	CLA	O1D-CGD-O2D-CED
4	4	608	CHL	O1D-CGD-O2D-CED
5	1	614	CLA	O1D-CGD-O2D-CED
5	2	604	CLA	O1D-CGD-O2D-CED
5	3	614	CLA	O1D-CGD-O2D-CED
5	3	604	CLA	CBD-CGD-O2D-CED
4	1	607	CHL	C3-C5-C6-C7
4	1	609	CHL	C3-C5-C6-C7
4	2	607	CHL	C3-C5-C6-C7
4	3	601	CHL	C3-C5-C6-C7
5	1	603	CLA	C3-C5-C6-C7
4	2	601	CHL	CBD-CGD-O2D-CED
4	2	605	CHL	CBD-CGD-O2D-CED
4	2	609	CHL	CBD-CGD-O2D-CED
4	3	605	CHL	CBD-CGD-O2D-CED
5	1	604	CLA	CBD-CGD-O2D-CED
5	4	612	CLA	CBD-CGD-O2D-CED
4	2	606	CHL	O1D-CGD-O2D-CED
5	1	603	CLA	O1D-CGD-O2D-CED
5	2	602	CLA	O1D-CGD-O2D-CED
5	3	612	CLA	O1D-CGD-O2D-CED
5	1	613	CLA	C3-C5-C6-C7
9	4	2630	LHG	O10-C23-O8-C6
10	4	623	BCR	C19-C20-C21-C22
5	2	603	CLA	O1A-CGA-O2A-C1
5	2	602	CLA	C3-C5-C6-C7
5	2	603	CLA	C3-C5-C6-C7
9	2	2630	LHG	O2-C2-C3-O3
4	2	608	CHL	O1D-CGD-O2D-CED
5	2	614	CLA	O1D-CGD-O2D-CED

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Mol	Chain	Res	Type	Atoms
5	3	603	CLA	O1D-CGD-O2D-CED
4	3	607	CHL	O1D-CGD-O2D-CED
5	2	603	CLA	CBA-CGA-O2A-C1
5	1	611	CLA	C2A-CAA-CBA-CGA
4	1	605	CHL	CBD-CGD-O2D-CED
4	3	601	CHL	CBD-CGD-O2D-CED
4	4	607	CHL	CBD-CGD-O2D-CED
9	2	2630	LHG	C1-C2-C3-O3
5	3	604	CLA	O1D-CGD-O2D-CED
5	4	612	CLA	O1D-CGD-O2D-CED
4	2	605	CHL	O1D-CGD-O2D-CED
4	3	605	CHL	O1D-CGD-O2D-CED
4	2	601	CHL	O1D-CGD-O2D-CED
7	2	1622	XAT	C31-C32-C33-C40
10	4	623	BCR	C7-C8-C9-C10
5	3	613	CLA	C3-C5-C6-C7
4	3	601	CHL	C5-C6-C7-C8
5	1	602	CLA	C10-C11-C12-C13
4	2	609	CHL	O1D-CGD-O2D-CED
7	3	1622	XAT	C29-C30-C31-C32
5	1	604	CLA	O1D-CGD-O2D-CED
4	3	609	CHL	CBD-CGD-O2D-CED
4	2	607	CHL	C10-C11-C12-C13
4	2	605	CHL	C2A-CAA-CBA-CGA
4	3	606	CHL	C2A-CAA-CBA-CGA
5	3	603	CLA	C2A-CAA-CBA-CGA
5	3	604	CLA	C2A-CAA-CBA-CGA
5	3	611	CLA	C2A-CAA-CBA-CGA
5	4	604	CLA	C2A-CAA-CBA-CGA
4	3	601	CHL	C10-C11-C12-C13
9	1	2630	LHG	C23-C24-C25-C26
5	2	610	CLA	CBD-CGD-O2D-CED
4	2	607	CHL	C8-C10-C11-C12
4	3	601	CHL	C8-C10-C11-C12
4	3	601	CHL	C13-C15-C16-C17
5	1	602	CLA	CBD-CGD-O2D-CED
5	2	602	CLA	C10-C11-C12-C13
5	3	602	CLA	CBD-CGD-O2D-CED
4	1	607	CHL	C5-C6-C7-C8
4	3	609	CHL	C10-C11-C12-C13
5	3	611	CLA	C5-C6-C7-C8
5	3	611	CLA	C3-C5-C6-C7

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Mol	Chain	Res	Type	Atoms
4	1	607	CHL	C2A-CAA-CBA-CGA
4	2	607	CHL	C2A-CAA-CBA-CGA
5	2	604	CLA	C2A-CAA-CBA-CGA
5	2	613	CLA	C2A-CAA-CBA-CGA
5	3	611	CLA	CBA-CGA-O2A-C1
4	2	609	CHL	C8-C10-C11-C12
5	3	610	CLA	C10-C11-C12-C13
4	3	609	CHL	C8-C10-C11-C12
4	1	607	CHL	C13-C15-C16-C17
10	4	623	BCR	C7-C8-C9-C34
4	3	605	CHL	C2A-CAA-CBA-CGA
4	4	601	CHL	C2A-CAA-CBA-CGA
9	2	2630	LHG	O1-C1-C2-C3
9	4	2630	LHG	O1-C1-C2-C3
4	1	607	CHL	C2C-C3C-CAC-CBC
4	2	609	CHL	C11-C12-C13-C15
4	3	601	CHL	O1D-CGD-O2D-CED
4	1	605	CHL	O1D-CGD-O2D-CED
4	1	609	CHL	CBA-CGA-O2A-C1
9	1	2630	LHG	C12-C13-C14-C15
4	4	607	CHL	O1D-CGD-O2D-CED
5	1	604	CLA	C4B-C3B-CAB-CBB
5	2	613	CLA	C4B-C3B-CAB-CBB
5	2	614	CLA	C4B-C3B-CAB-CBB
5	3	604	CLA	C4B-C3B-CAB-CBB
5	4	603	CLA	C4B-C3B-CAB-CBB
9	3	2630	LHG	C23-C24-C25-C26
4	2	609	CHL	C11-C12-C13-C14
4	3	608	CHL	C2A-CAA-CBA-CGA
9	2	2630	LHG	C23-C24-C25-C26
5	3	611	CLA	O1A-CGA-O2A-C1
4	1	605	CHL	C3A-C2A-CAA-CBA
4	1	607	CHL	C3A-C2A-CAA-CBA
4	2	605	CHL	C3A-C2A-CAA-CBA
4	2	607	CHL	C3A-C2A-CAA-CBA
4	3	605	CHL	C3A-C2A-CAA-CBA
4	4	608	CHL	C3A-C2A-CAA-CBA
5	1	611	CLA	C3A-C2A-CAA-CBA
9	3	2630	LHG	C28-C29-C30-C31
4	2	609	CHL	CBA-CGA-O2A-C1
9	3	2630	LHG	C14-C15-C16-C17
5	1	603	CLA	C2B-C3B-CAB-CBB

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Mol	Chain	Res	Type	Atoms
5	3	611	CLA	C2B-C3B-CAB-CBB
5	4	603	CLA	C2B-C3B-CAB-CBB
6	1	1620	LUT	C1-C6-C7-C8
6	1	1620	LUT	C5-C6-C7-C8
6	2	1620	LUT	C1-C6-C7-C8
6	2	1620	LUT	C5-C6-C7-C8
4	3	607	CHL	C2A-CAA-CBA-CGA
9	2	2630	LHG	C12-C13-C14-C15
4	1	607	CHL	C8-C10-C11-C12
9	3	2630	LHG	C10-C11-C12-C13
9	1	2630	LHG	C9-C10-C11-C12
6	1	1621	LUT	C29-C30-C31-C32
5	3	602	CLA	C11-C12-C13-C15
5	3	602	CLA	C11-C12-C13-C14
4	2	607	CHL	C4-C3-C5-C6
4	3	609	CHL	C2C-C3C-CAC-CBC
4	1	609	CHL	O1A-CGA-O2A-C1
4	1	609	CHL	C2C-C3C-CAC-CBC
9	2	2630	LHG	C11-C12-C13-C14
9	3	2630	LHG	O7-C5-C6-O8
9	4	2630	LHG	O7-C5-C6-O8
9	2	2630	LHG	C24-C25-C26-C27
4	1	609	CHL	C8-C10-C11-C12
4	2	607	CHL	C2-C3-C5-C6
5	4	603	CLA	CBD-CGD-O2D-CED
5	1	604	CLA	CBA-CGA-O2A-C1
9	3	2630	LHG	C33-C34-C35-C36
9	2	2630	LHG	O1-C1-C2-O2
9	3	2630	LHG	O1-C1-C2-O2
4	2	609	CHL	O1A-CGA-O2A-C1
5	1	604	CLA	C1A-C2A-CAA-CBA
5	1	610	CLA	C1A-C2A-CAA-CBA
5	1	611	CLA	C1A-C2A-CAA-CBA
5	2	602	CLA	C1A-C2A-CAA-CBA
5	2	610	CLA	C1A-C2A-CAA-CBA
5	3	604	CLA	C1A-C2A-CAA-CBA
5	3	610	CLA	C1A-C2A-CAA-CBA
5	4	611	CLA	C1A-C2A-CAA-CBA
5	3	602	CLA	C6-C7-C8-C10
4	1	607	CHL	C2-C3-C5-C6
4	3	607	CHL	C2-C3-C5-C6
5	3	602	CLA	C2A-CAA-CBA-CGA

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Mol	Chain	Res	Type	Atoms
4	1	609	CHL	C11-C10-C8-C9
9	2	2630	LHG	C4-C5-C6-O8
9	3	2630	LHG	C4-C5-C6-O8
5	1	602	CLA	O1D-CGD-O2D-CED
4	1	609	CHL	CHA-CBD-CGD-O2D
4	3	608	CHL	CHA-CBD-CGD-O1D
4	1	607	CHL	C4-C3-C5-C6
4	3	607	CHL	C4-C3-C5-C6
5	2	610	CLA	O1D-CGD-O2D-CED
5	1	603	CLA	CBA-CGA-O2A-C1
4	2	607	CHL	C5-C6-C7-C8
4	3	606	CHL	C3C-C2C-CMC-OMC
4	4	606	CHL	C3C-C2C-CMC-OMC
8	2	1623	NEX	C9-C10-C11-C12
5	1	604	CLA	O1A-CGA-O2A-C1
9	1	2630	LHG	C27-C28-C29-C30
9	1	2630	LHG	C26-C27-C28-C29
4	2	607	CHL	C2C-C3C-CAC-CBC
9	3	2630	LHG	C16-C17-C18-C19
5	3	610	CLA	CBA-CGA-O2A-C1
4	1	606	CHL	C1C-C2C-CMC-OMC
4	2	606	CHL	C1C-C2C-CMC-OMC
4	3	606	CHL	C1C-C2C-CMC-OMC
9	3	2630	LHG	C35-C36-C37-C38
5	3	602	CLA	O1D-CGD-O2D-CED
5	1	613	CLA	C6-C7-C8-C10
5	1	603	CLA	C4-C3-C5-C6
4	3	609	CHL	CBA-CGA-O2A-C1
4	1	607	CHL	C6-C7-C8-C9
5	3	602	CLA	C6-C7-C8-C9
5	1	610	CLA	C4B-C3B-CAB-CBB
5	1	612	CLA	C4B-C3B-CAB-CBB
5	1	614	CLA	C4B-C3B-CAB-CBB
5	4	604	CLA	C4B-C3B-CAB-CBB
5	4	610	CLA	C4B-C3B-CAB-CBB
4	1	607	CHL	C6-C7-C8-C10
4	2	607	CHL	C6-C7-C8-C10
4	3	609	CHL	O1D-CGD-O2D-CED
9	3	2630	LHG	O8-C23-C24-C25
5	3	602	CLA	C8-C10-C11-C12
9	1	2630	LHG	C11-C12-C13-C14
4	2	601	CHL	C1A-C2A-CAA-CBA

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Mol	Chain	Res	Type	Atoms
4	2	607	CHL	C1A-C2A-CAA-CBA
4	2	608	CHL	C1A-C2A-CAA-CBA
4	3	606	CHL	C1A-C2A-CAA-CBA
4	3	608	CHL	C1A-C2A-CAA-CBA
4	4	606	CHL	C1A-C2A-CAA-CBA
5	1	603	CLA	C2-C3-C5-C6
5	3	610	CLA	C2B-C3B-CAB-CBB
6	1	1621	LUT	C1-C6-C7-C8
10	4	623	BCR	C1-C6-C7-C8
9	2	2630	LHG	O7-C5-C6-O8
5	1	603	CLA	O1A-CGA-O2A-C1
5	1	610	CLA	C5-C6-C7-C8
5	1	613	CLA	C6-C7-C8-C9
4	2	607	CHL	C6-C7-C8-C9
4	3	601	CHL	C14-C13-C15-C16
9	3	2630	LHG	C24-C25-C26-C27
4	3	601	CHL	C15-C16-C17-C18
5	3	610	CLA	O1A-CGA-O2A-C1
4	3	601	CHL	C11-C10-C8-C7
5	1	602	CLA	C6-C7-C8-C10
6	4	620	LUT	C27-C28-C29-C30
5	2	602	CLA	C2A-CAA-CBA-CGA
4	3	609	CHL	O1A-CGA-O2A-C1
9	1	2630	LHG	C15-C16-C17-C18
9	4	2630	LHG	C4-C5-C6-O8
5	1	602	CLA	C6-C7-C8-C9
5	4	603	CLA	O1D-CGD-O2D-CED
9	3	2630	LHG	O10-C23-O8-C6
5	1	614	CLA	C1A-C2A-CAA-CBA
5	2	612	CLA	C4B-C3B-CAB-CBB
5	3	602	CLA	C1A-C2A-CAA-CBA
5	2	603	CLA	C4-C3-C5-C6
9	2	2630	LHG	O10-C23-O8-C6
5	2	602	CLA	C11-C12-C13-C15
4	3	601	CHL	C11-C10-C8-C9
4	1	607	CHL	C10-C11-C12-C13
4	4	609	CHL	C2C-C3C-CAC-CBC
8	1	1623	NEX	C33-C34-C35-C15
5	1	603	CLA	CAD-CBD-CGD-O2D
5	2	603	CLA	CAD-CBD-CGD-O2D
5	3	603	CLA	CAD-CBD-CGD-O2D
5	3	604	CLA	CAD-CBD-CGD-O2D

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Mol	Chain	Res	Type	Atoms
5	1	602	CLA	C2A-CAA-CBA-CGA
5	4	602	CLA	C2A-CAA-CBA-CGA
4	2	608	CHL	C2C-C3C-CAC-CBC
4	1	607	CHL	CHA-CBD-CGD-O2D
4	2	606	CHL	CHA-CBD-CGD-O2D
4	3	607	CHL	CHA-CBD-CGD-O1D
4	3	607	CHL	CHA-CBD-CGD-O2D
4	3	608	CHL	CHA-CBD-CGD-O2D
5	1	603	CLA	CAD-CBD-CGD-O1D
5	2	603	CLA	CAD-CBD-CGD-O1D
5	3	602	CLA	CHA-CBD-CGD-O1D
5	3	603	CLA	CAD-CBD-CGD-O1D
5	3	604	CLA	CAD-CBD-CGD-O1D
5	4	612	CLA	CHA-CBD-CGD-O1D
5	4	612	CLA	CHA-CBD-CGD-O2D
8	3	1623	NEX	C9-C10-C11-C12
9	3	2630	LHG	C4-O6-P-O4
5	1	604	CLA	C2B-C3B-CAB-CBB
5	1	610	CLA	C2B-C3B-CAB-CBB
5	1	612	CLA	C2B-C3B-CAB-CBB
5	2	613	CLA	C2B-C3B-CAB-CBB
5	2	614	CLA	C2B-C3B-CAB-CBB
5	3	603	CLA	C2B-C3B-CAB-CBB
5	3	604	CLA	C2B-C3B-CAB-CBB
5	4	604	CLA	C2B-C3B-CAB-CBB
9	3	2630	LHG	C27-C28-C29-C30
6	4	620	LUT	C27-C28-C29-C39
10	4	623	BCR	C36-C18-C19-C20
9	1	2630	LHG	C14-C15-C16-C17
6	4	620	LUT	C29-C30-C31-C32
5	3	602	CLA	C11-C10-C8-C9
5	3	603	CLA	C5-C6-C7-C8
6	3	1621	LUT	C29-C30-C31-C32
4	1	606	CHL	C2A-CAA-CBA-CGA
8	1	1623	NEX	C31-C32-C33-C40
10	4	623	BCR	C9-C10-C11-C12
4	3	609	CHL	C4-C3-C5-C6
5	2	602	CLA	C4-C3-C5-C6
5	2	603	CLA	C2-C3-C5-C6
5	3	602	CLA	C10-C11-C12-C13
8	1	1623	NEX	C29-C30-C31-C32
5	4	610	CLA	C3A-C2A-CAA-CBA

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Mol	Chain	Res	Type	Atoms
8	1	1623	NEX	C39-C29-C30-C31
8	2	1623	NEX	C39-C29-C30-C31
8	3	1623	NEX	C39-C29-C30-C31
4	2	609	CHL	C5-C6-C7-C8
9	3	2630	LHG	C15-C16-C17-C18
9	3	2630	LHG	C29-C30-C31-C32
5	2	614	CLA	CAA-CBA-CGA-O2A
4	2	609	CHL	C6-C7-C8-C9
5	1	612	CLA	CAA-CBA-CGA-O1A
4	1	607	CHL	C1A-C2A-CAA-CBA
4	1	609	CHL	C1A-C2A-CAA-CBA
4	3	609	CHL	C1A-C2A-CAA-CBA
4	4	607	CHL	C1A-C2A-CAA-CBA
5	1	612	CLA	CAA-CBA-CGA-O2A
4	4	607	CHL	CAA-CBA-CGA-O1A
5	2	614	CLA	CAA-CBA-CGA-O1A
5	4	603	CLA	CAA-CBA-CGA-O1A
8	1	1623	NEX	C28-C29-C30-C31
8	2	1623	NEX	C28-C29-C30-C31
8	3	1623	NEX	C28-C29-C30-C31
5	3	604	CLA	CAA-CBA-CGA-O2A
5	1	613	CLA	C2B-C3B-CAB-CBB
5	2	612	CLA	C2B-C3B-CAB-CBB
6	1	1621	LUT	C5-C6-C7-C8
6	3	1621	LUT	C1-C6-C7-C8
10	4	623	BCR	C5-C6-C7-C8
4	3	609	CHL	C2-C3-C5-C6
4	4	609	CHL	CAA-CBA-CGA-O1A
5	1	613	CLA	C2A-CAA-CBA-CGA
4	4	609	CHL	CAA-CBA-CGA-O2A
5	4	612	CLA	CAA-CBA-CGA-O2A
4	3	608	CHL	CAA-CBA-CGA-O2A
7	4	622	XAT	C27-C28-C29-C39
5	4	603	CLA	CAA-CBA-CGA-O2A
4	4	607	CHL	CAA-CBA-CGA-O2A
5	4	604	CLA	CAA-CBA-CGA-O2A
5	4	612	CLA	CAA-CBA-CGA-O1A
4	1	605	CHL	CAA-CBA-CGA-O2A
5	3	604	CLA	CAA-CBA-CGA-O1A
9	3	2630	LHG	C5-C4-O6-P
4	3	605	CHL	CAA-CBA-CGA-O2A
5	4	611	CLA	CAA-CBA-CGA-O2A

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Mol	Chain	Res	Type	Atoms
5	4	602	CLA	C4B-C3B-CAB-CBB
5	3	614	CLA	O2A-C1-C2-C3
4	1	605	CHL	CAA-CBA-CGA-O1A
4	1	601	CHL	CHA-CBD-CGD-O1D
4	1	601	CHL	CHA-CBD-CGD-O2D
4	1	607	CHL	CHA-CBD-CGD-O1D
4	2	609	CHL	C4-C3-C5-C6
4	1	601	CHL	CAA-CBA-CGA-O2A
5	4	604	CLA	CAA-CBA-CGA-O1A
4	3	607	CHL	C3A-C2A-CAA-CBA
4	1	606	CHL	CAA-CBA-CGA-O2A
4	4	608	CHL	CAA-CBA-CGA-O2A
4	3	608	CHL	CAA-CBA-CGA-O1A
4	3	605	CHL	CAA-CBA-CGA-O1A
5	1	611	CLA	CAA-CBA-CGA-O2A
5	4	610	CLA	CAA-CBA-CGA-O2A
4	1	601	CHL	C3C-C2C-CMC-OMC
5	4	611	CLA	CAA-CBA-CGA-O1A
9	1	2630	LHG	C18-C19-C20-C21
4	1	606	CHL	CAA-CBA-CGA-O1A
7	4	622	XAT	O4-C6-C7-C8
8	1	1623	NEX	O24-C26-C27-C28
8	2	1623	NEX	O24-C26-C27-C28
8	3	1623	NEX	O24-C26-C27-C28
5	4	610	CLA	CAA-CBA-CGA-O1A
4	4	607	CHL	C2A-CAA-CBA-CGA
4	4	608	CHL	CAA-CBA-CGA-O1A
6	1	1621	LUT	C9-C10-C11-C12
5	1	611	CLA	CAA-CBA-CGA-O1A
5	2	602	CLA	C11-C12-C13-C14
5	2	603	CLA	CAA-CBA-CGA-O2A
5	3	611	CLA	CAA-CBA-CGA-O2A
4	2	601	CHL	CAA-CBA-CGA-O1A
4	3	601	CHL	C11-C12-C13-C15
5	3	613	CLA	C6-C7-C8-C10
4	1	601	CHL	CAA-CBA-CGA-O1A
5	2	611	CLA	CAA-CBA-CGA-O2A
5	1	602	CLA	C2B-C3B-CAB-CBB
5	1	614	CLA	C2B-C3B-CAB-CBB
5	4	610	CLA	C2B-C3B-CAB-CBB
6	3	1620	LUT	C5-C6-C7-C8
6	3	1621	LUT	C5-C6-C7-C8

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Mol	Chain	Res	Type	Atoms
5	1	604	CLA	C2-C1-O2A-CGA
9	2	2630	LHG	C13-C14-C15-C16
5	2	602	CLA	C2-C3-C5-C6
4	1	601	CHL	C2A-CAA-CBA-CGA
5	3	602	CLA	CAA-CBA-CGA-O2A
5	2	612	CLA	C1A-C2A-CAA-CBA
5	4	612	CLA	C1A-C2A-CAA-CBA
4	1	607	CHL	CAA-CBA-CGA-O2A
4	3	607	CHL	CAA-CBA-CGA-O2A
5	1	602	CLA	CAA-CBA-CGA-O2A
6	1	1621	LUT	C7-C8-C9-C10
7	2	1622	XAT	C13-C14-C15-C35
4	2	607	CHL	CAA-CBA-CGA-O2A
5	2	602	CLA	CAA-CBA-CGA-O2A
5	3	613	CLA	C2A-CAA-CBA-CGA
4	2	609	CHL	C2-C1-O2A-CGA
4	3	609	CHL	C2-C1-O2A-CGA
4	2	607	CHL	C11-C10-C8-C7
5	2	602	CLA	C6-C7-C8-C10
5	2	603	CLA	C5-C6-C7-C8
4	1	609	CHL	C5-C6-C7-C8
5	2	611	CLA	CAA-CBA-CGA-O1A
5	2	612	CLA	C3A-C2A-CAA-CBA
5	3	611	CLA	C3A-C2A-CAA-CBA
5	4	604	CLA	C3A-C2A-CAA-CBA
5	4	612	CLA	C3A-C2A-CAA-CBA
9	3	2630	LHG	C9-C10-C11-C12
5	3	613	CLA	C6-C7-C8-C9
5	2	603	CLA	CAA-CBA-CGA-O1A
4	2	601	CHL	CAA-CBA-CGA-O2A
10	4	623	BCR	C17-C18-C19-C20
4	3	607	CHL	CAA-CBA-CGA-O1A
5	3	611	CLA	CAA-CBA-CGA-O1A
5	2	602	CLA	CAA-CBA-CGA-O1A
4	2	607	CHL	CAA-CBA-CGA-O1A
5	3	602	CLA	CAA-CBA-CGA-O1A
9	3	2630	LHG	C17-C18-C19-C20
5	1	602	CLA	CAD-CBD-CGD-O2D
5	4	611	CLA	CAD-CBD-CGD-O2D
9	2	2630	LHG	C10-C11-C12-C13
4	3	607	CHL	C1A-C2A-CAA-CBA
4	2	605	CHL	CAA-CBA-CGA-O2A

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Mol	Chain	Res	Type	Atoms
4	1	607	CHL	CAA-CBA-CGA-O1A
5	1	602	CLA	CAA-CBA-CGA-O1A

There are no ring outliers.

58 monomers are involved in 119 short contacts:

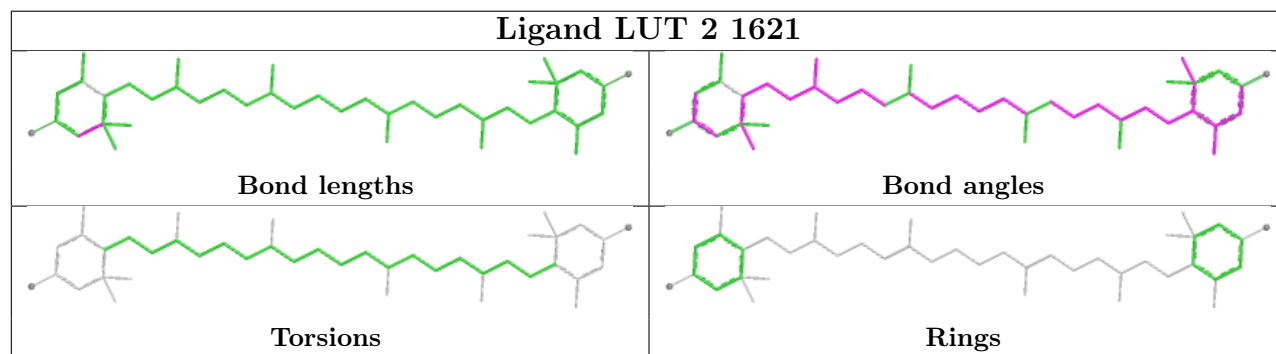
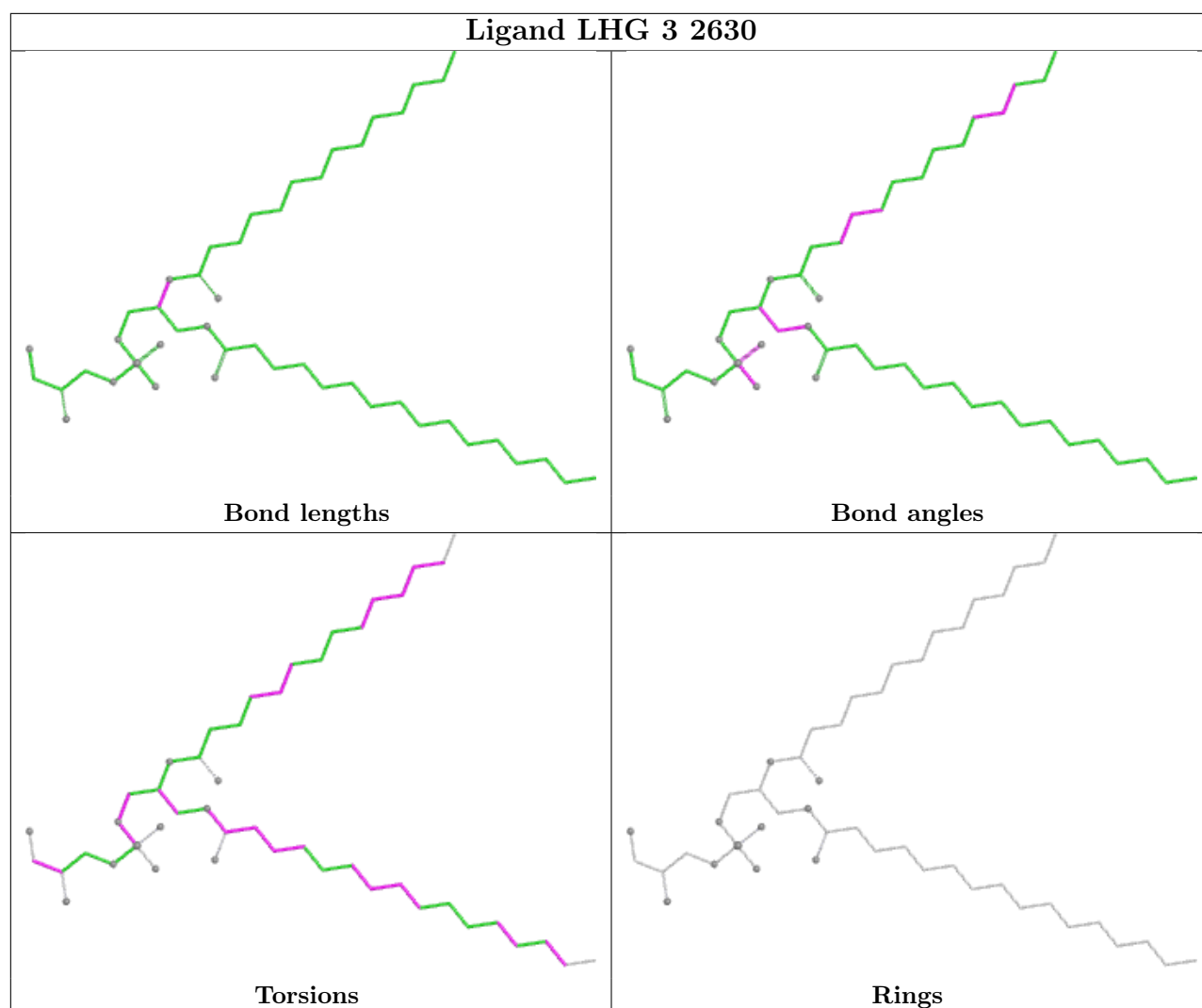
Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	3	2630	LHG	3	0
6	2	1621	LUT	4	0
5	2	613	CLA	2	0
9	4	2630	LHG	1	0
4	3	606	CHL	4	0
4	1	606	CHL	2	0
4	4	601	CHL	1	0
5	1	613	CLA	1	0
6	1	1620	LUT	5	0
5	2	612	CLA	2	0
7	3	1622	XAT	5	0
5	2	602	CLA	3	0
4	2	608	CHL	2	0
4	2	607	CHL	2	0
5	3	602	CLA	4	0
6	2	1620	LUT	4	0
5	1	604	CLA	1	0
5	2	610	CLA	3	0
7	4	622	XAT	4	0
4	3	605	CHL	1	0
6	1	1621	LUT	4	0
6	4	620	LUT	4	0
5	4	603	CLA	1	0
8	3	1623	NEX	3	0
7	2	1622	XAT	6	0
8	2	1623	NEX	1	0
4	1	609	CHL	5	0
4	3	601	CHL	2	0
5	4	612	CLA	2	0
8	1	1623	NEX	2	0
4	3	608	CHL	3	0
5	3	613	CLA	3	0
4	3	607	CHL	6	0
4	2	609	CHL	4	0
4	4	607	CHL	1	0

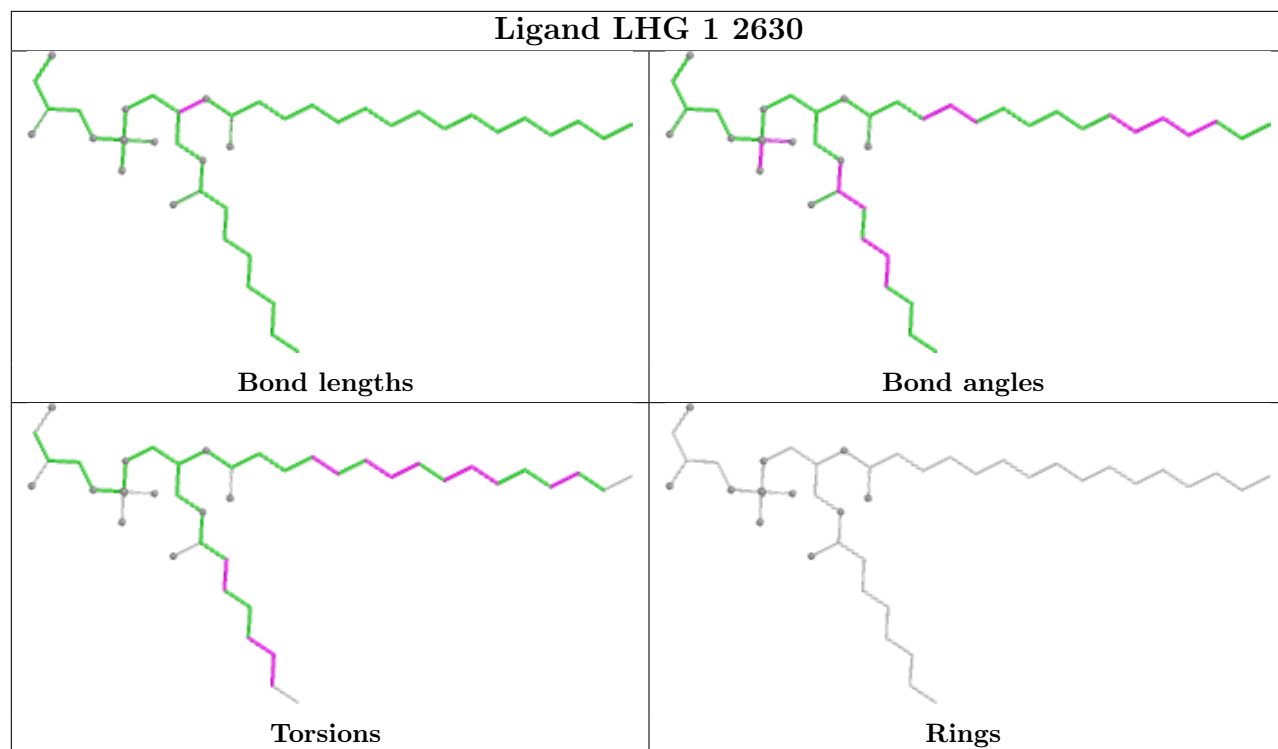
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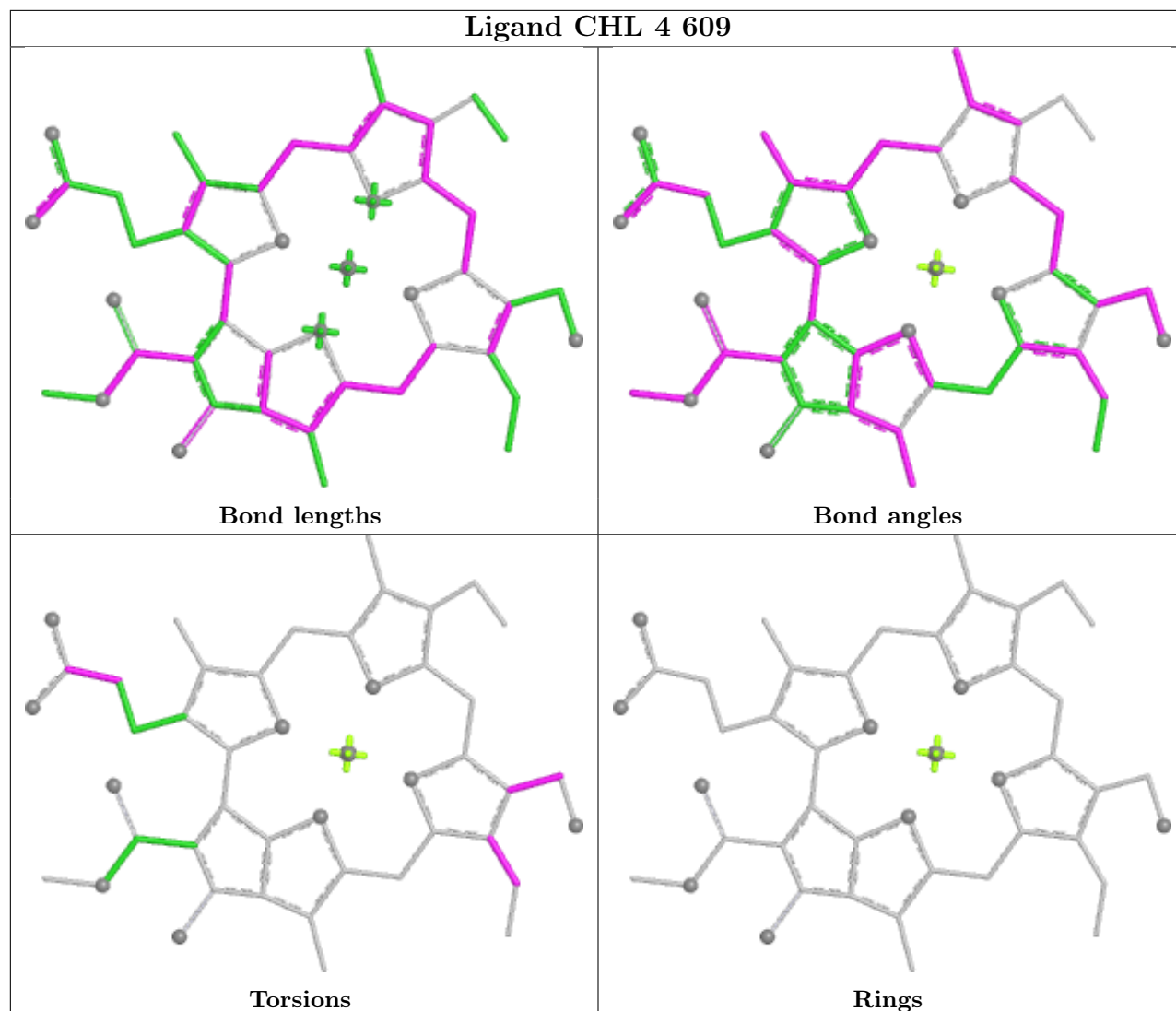
Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	2	2630	LHG	2	0
5	3	611	CLA	2	0
10	4	623	BCR	3	0
4	1	608	CHL	1	0
5	1	612	CLA	2	0
4	1	607	CHL	4	0
5	3	603	CLA	1	0
5	1	610	CLA	2	0
4	4	608	CHL	1	0
4	4	606	CHL	1	0
4	2	606	CHL	3	0
5	1	603	CLA	2	0
5	4	610	CLA	1	0
4	3	609	CHL	4	0
5	3	610	CLA	2	0
6	3	1620	LUT	7	0
5	2	603	CLA	2	0
5	1	602	CLA	4	0
5	2	604	CLA	1	0
7	1	1622	XAT	4	0
6	3	1621	LUT	3	0
4	2	605	CHL	1	0
5	4	611	CLA	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.

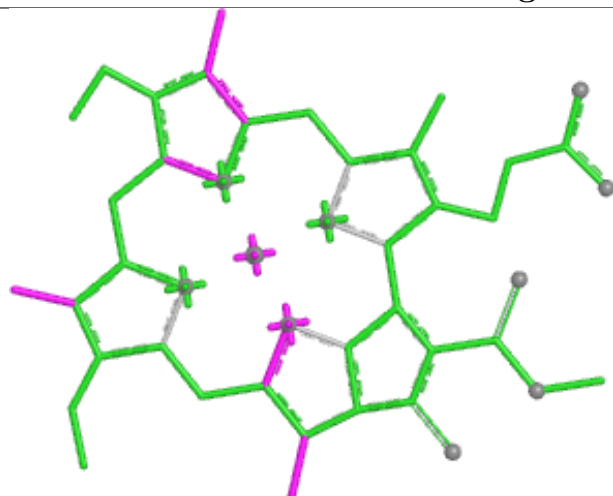




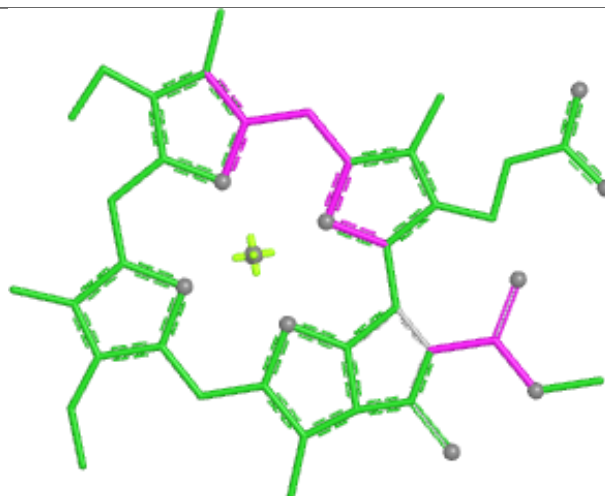
Ligand CHL 4 609



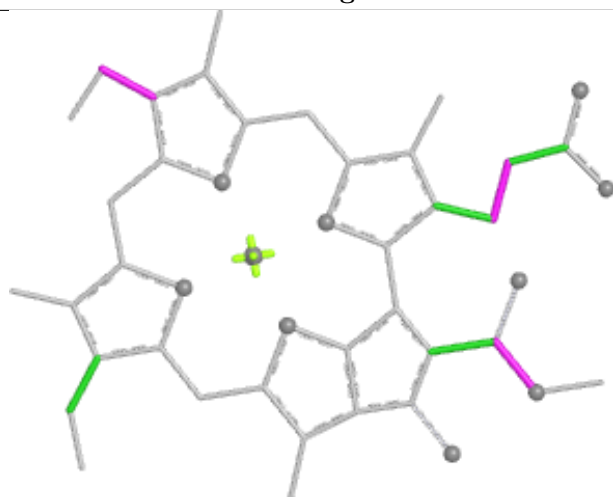
Ligand CLA 2 613



Bond lengths



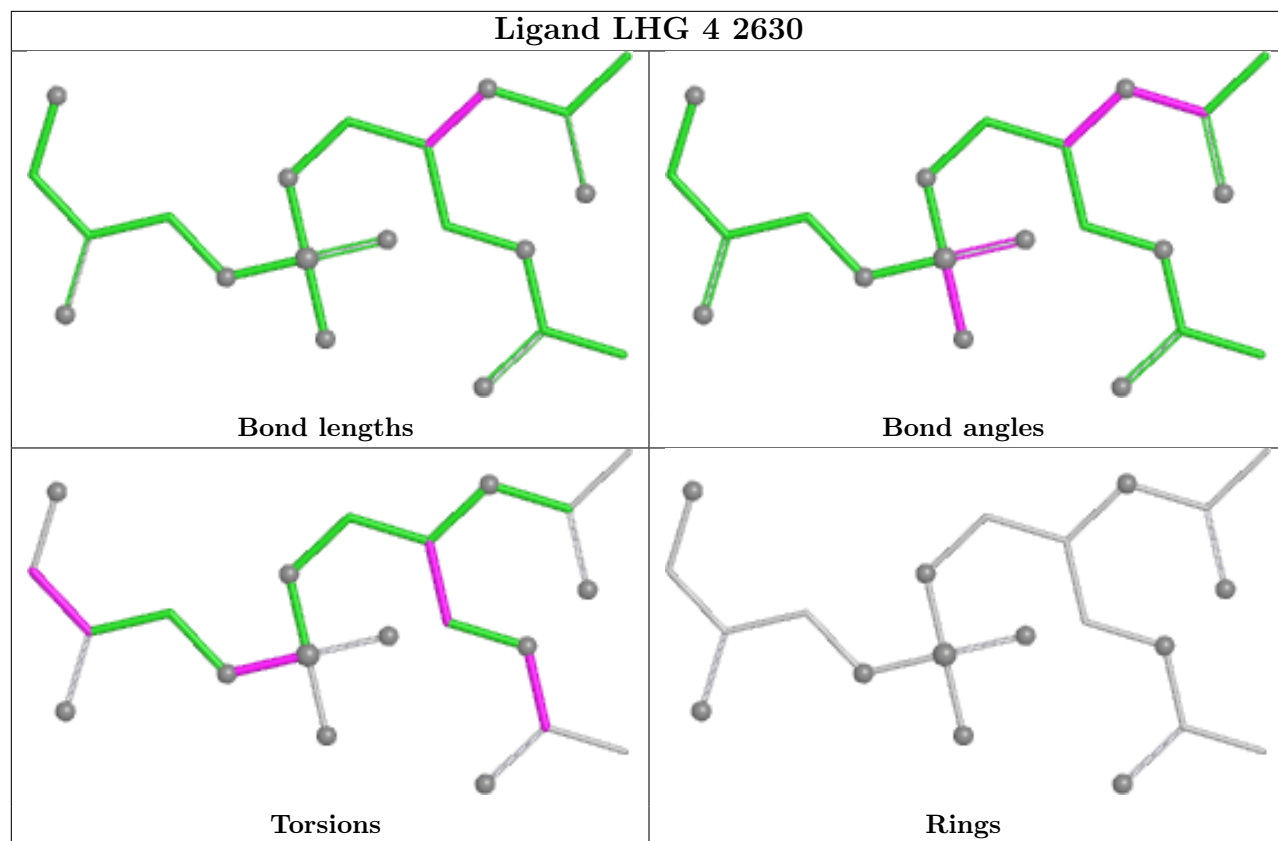
Bond angles



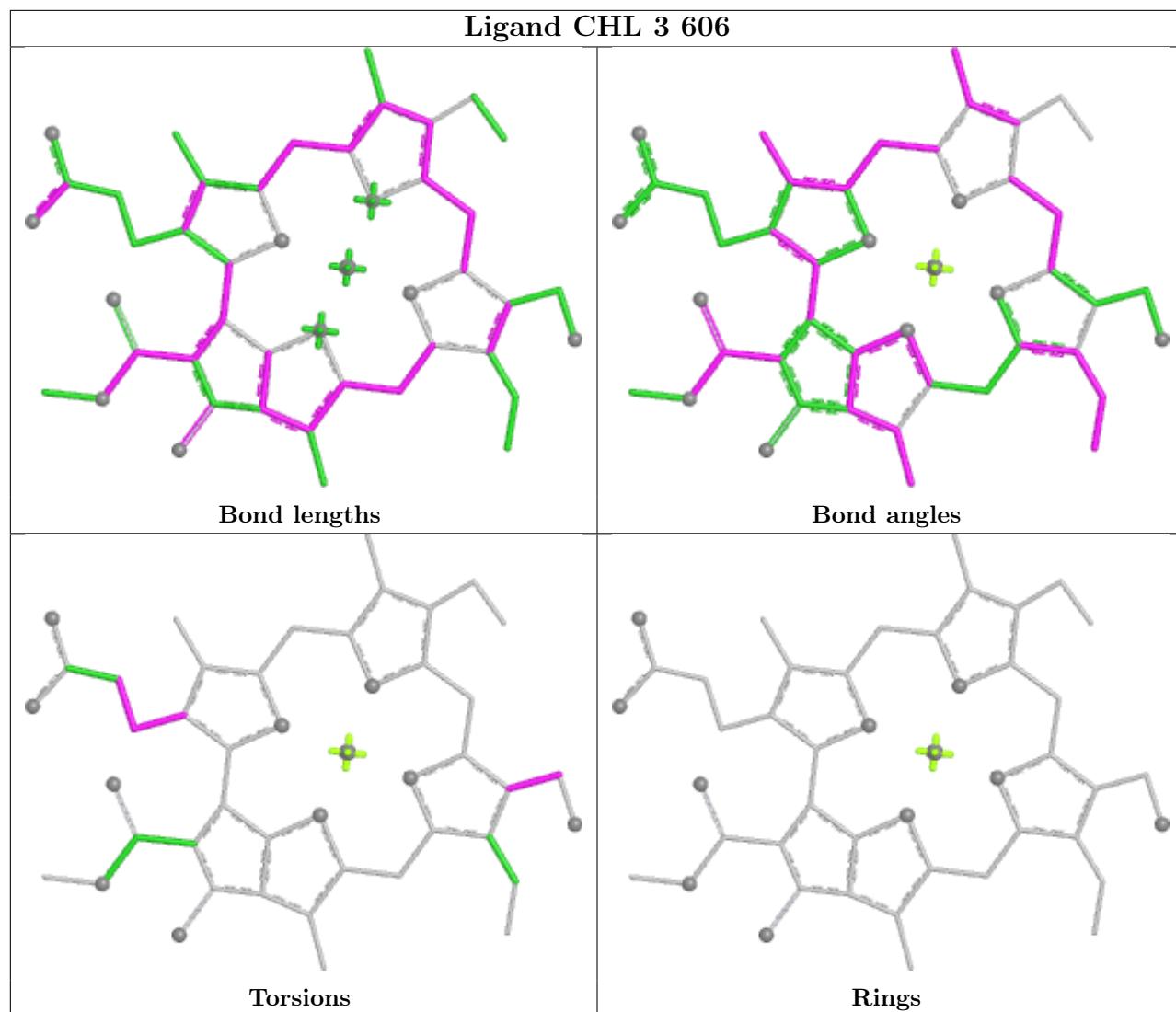
Torsions



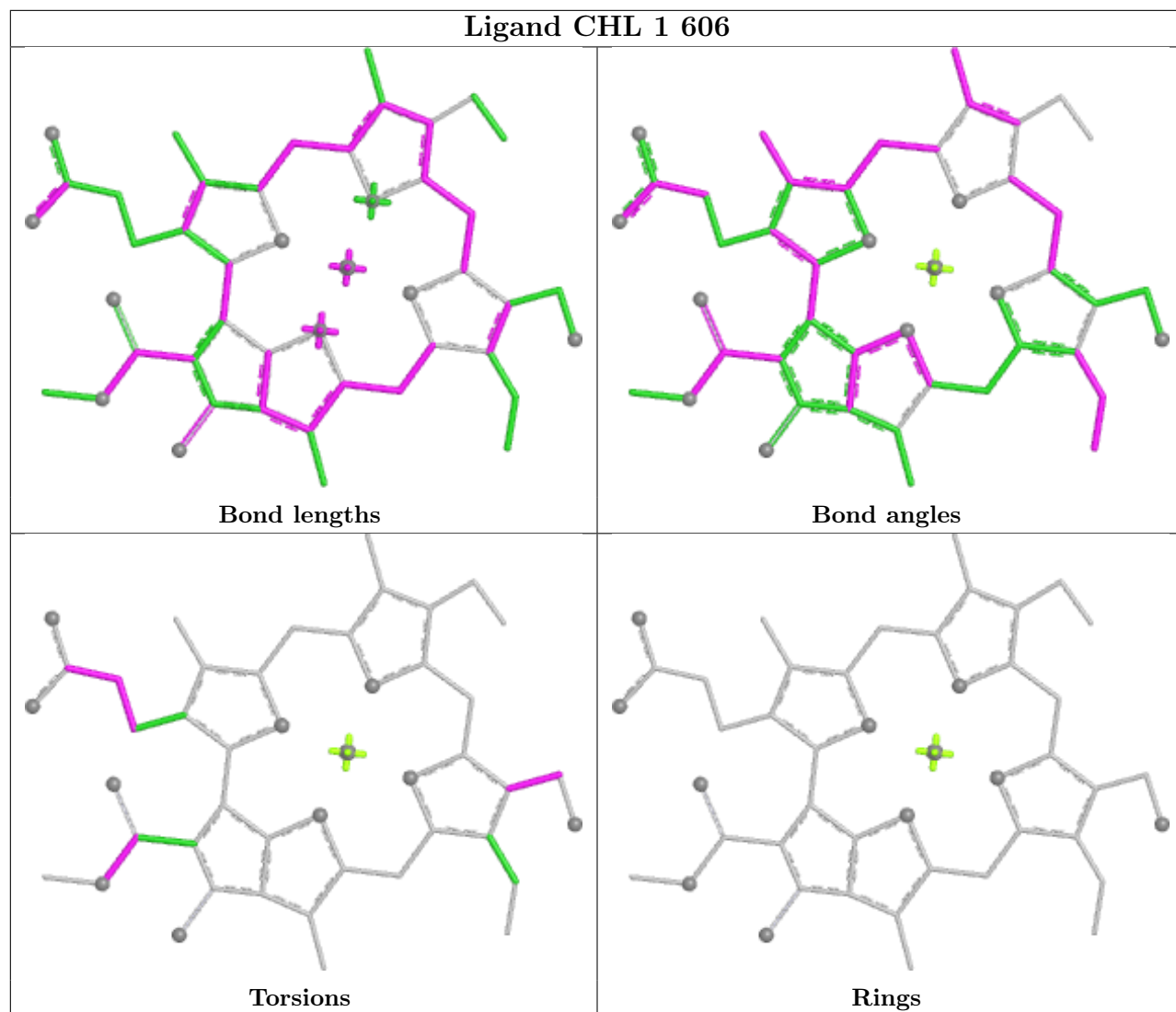
Rings



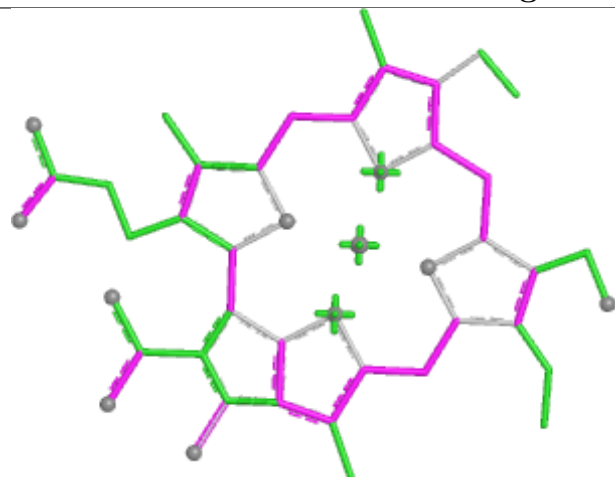
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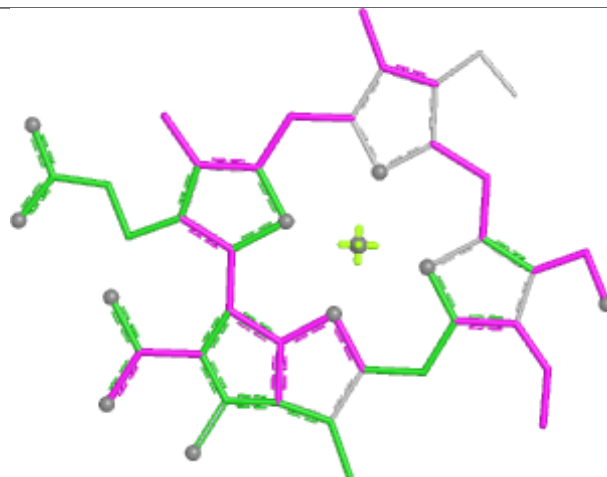
Ligand CHL 1 606



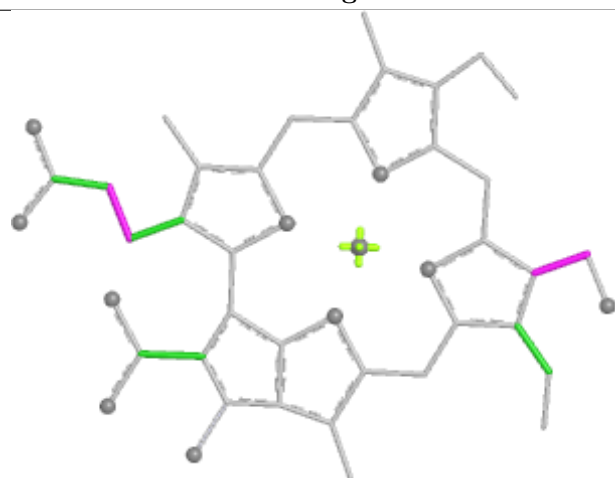
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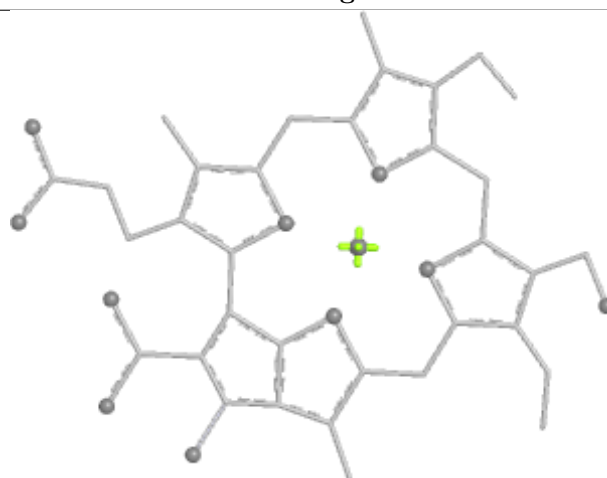
Bond lengths



Bond angles

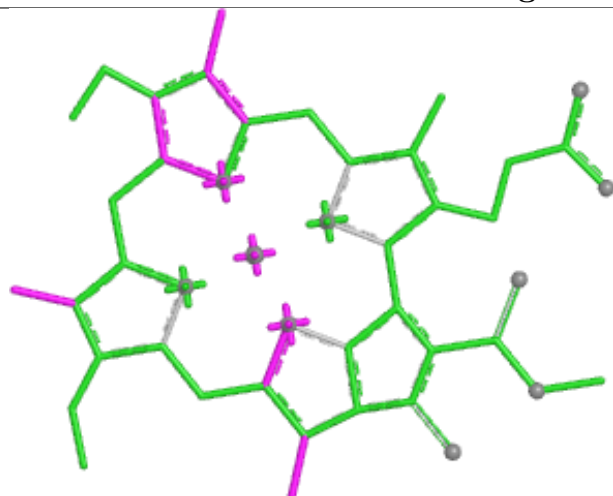


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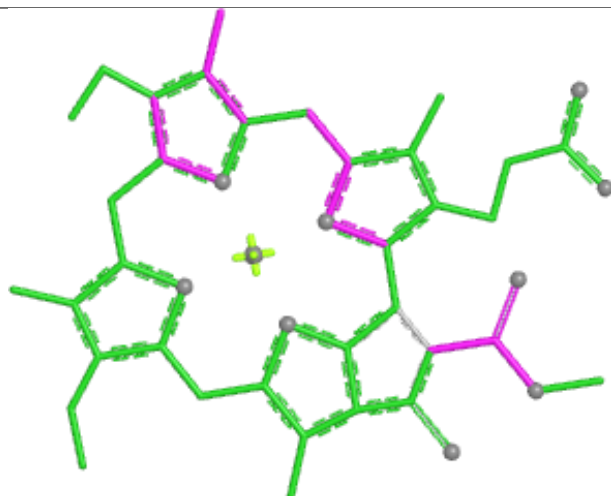


Rings

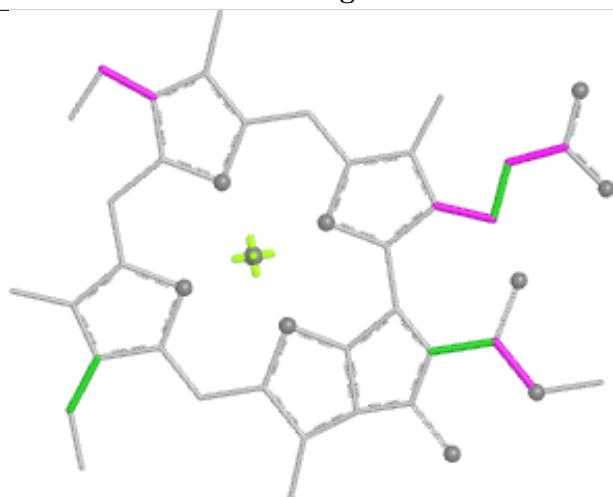
Ligand CLA 2 614



Bond lengths



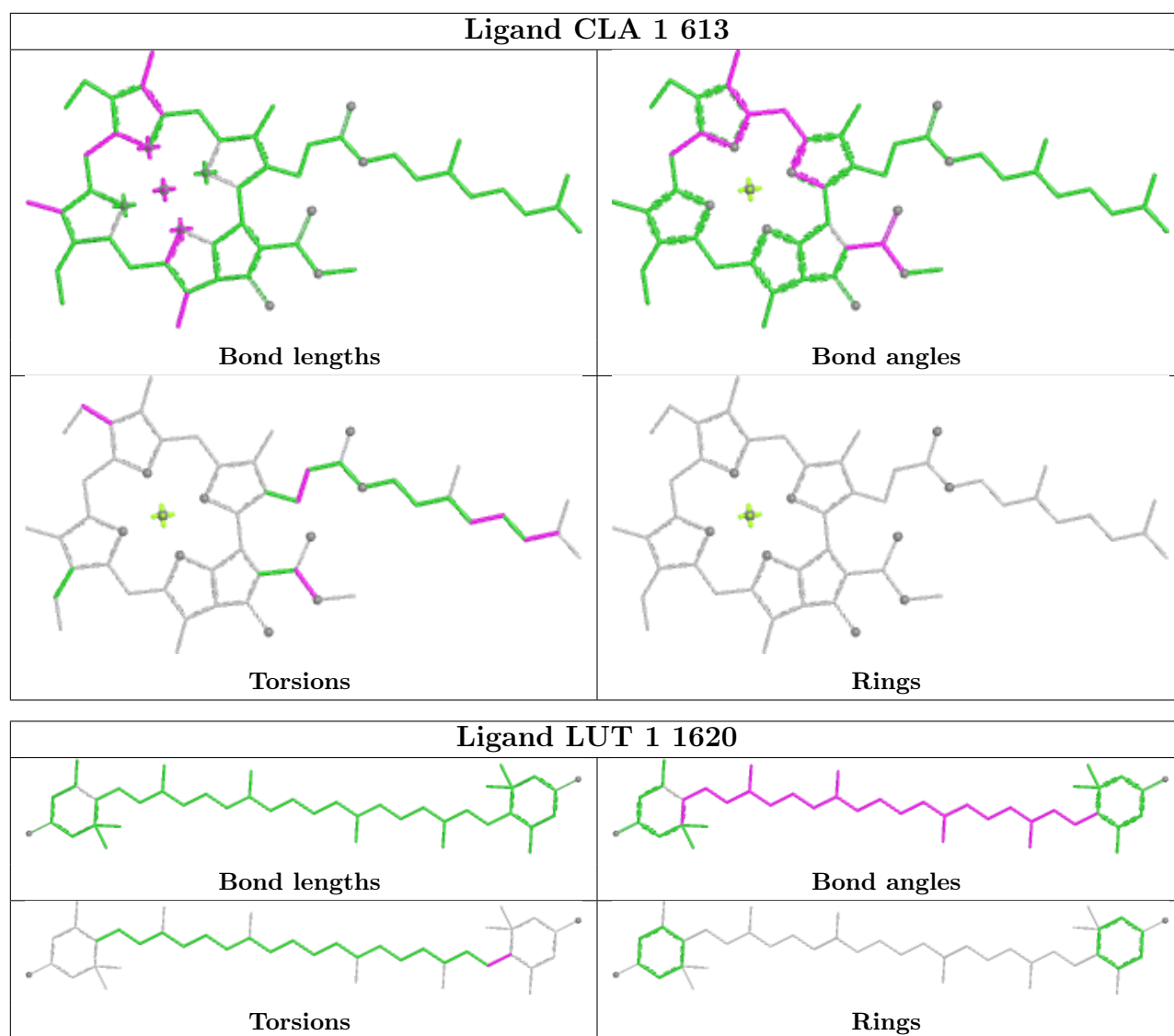
Bond angles



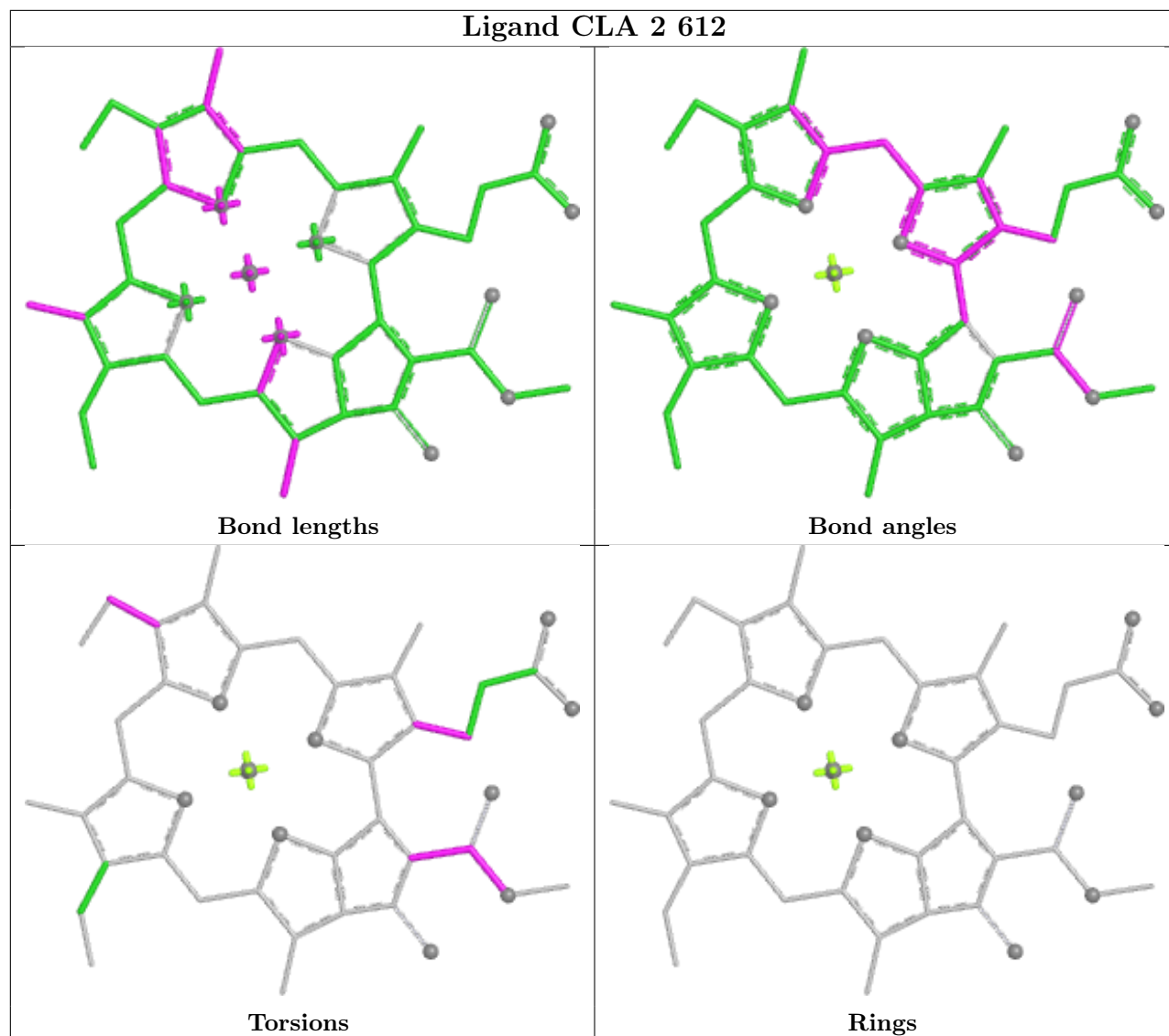
Torsions



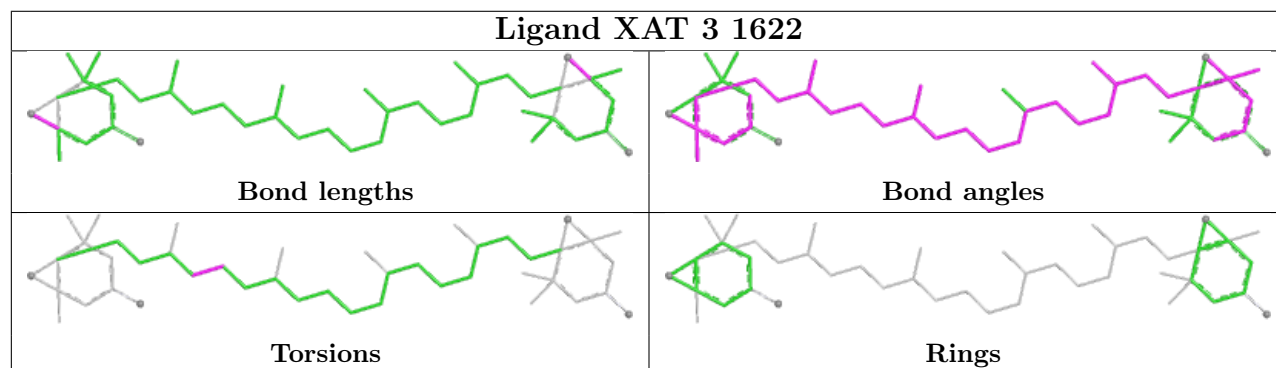
Rings



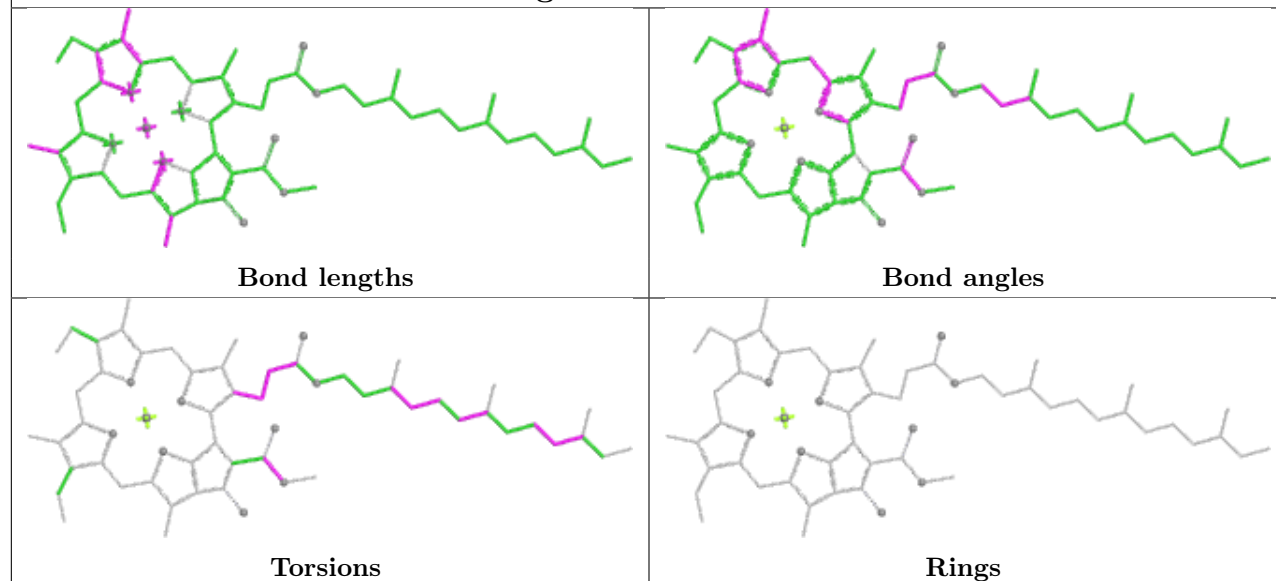
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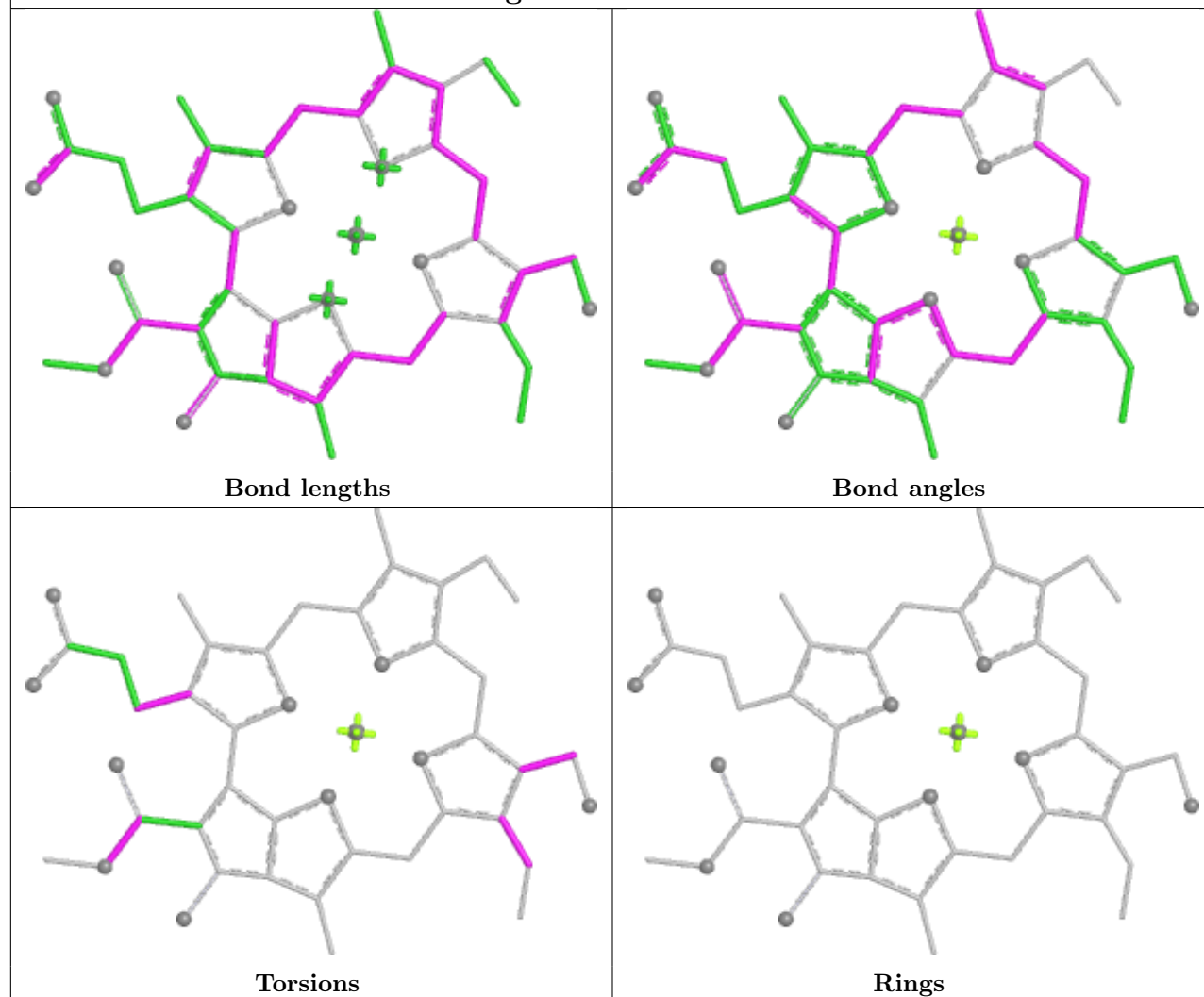
Ligand XAT 3 1622



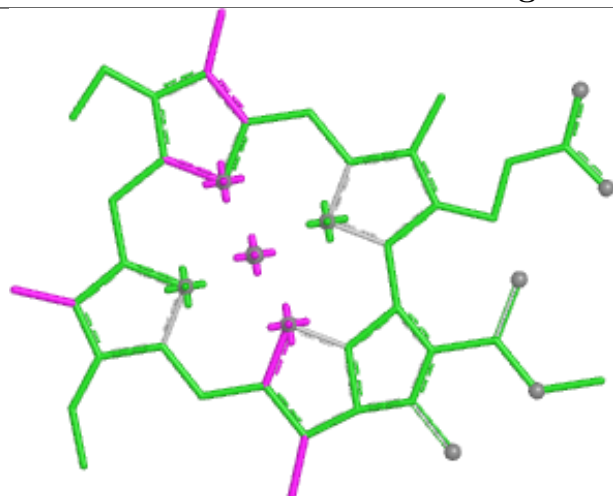
Ligand CLA 2 602



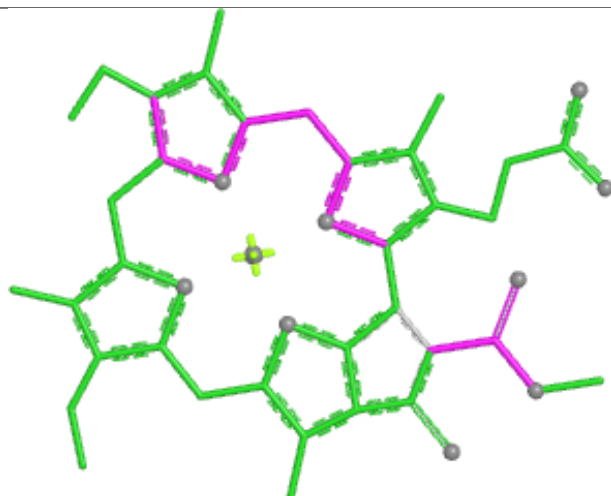
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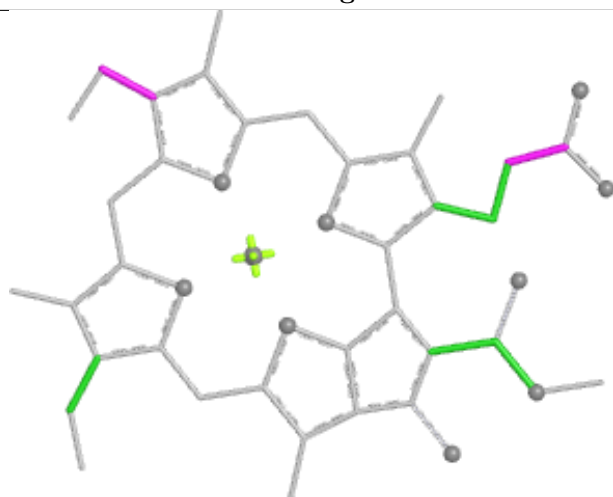
Ligand CLA 2 611



Bond lengths



Bond angles

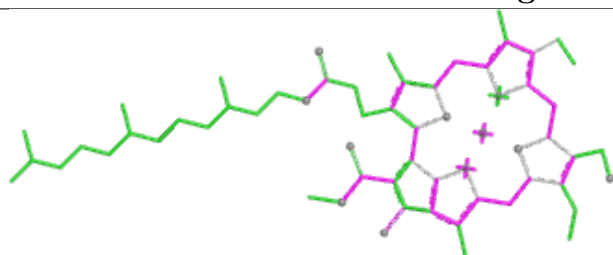


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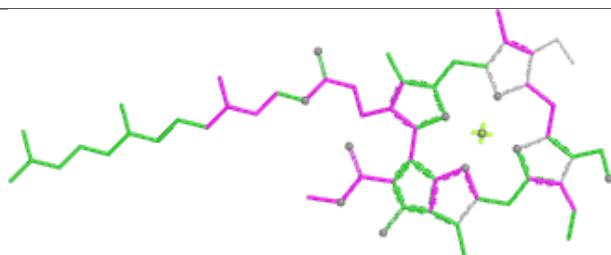


Rings

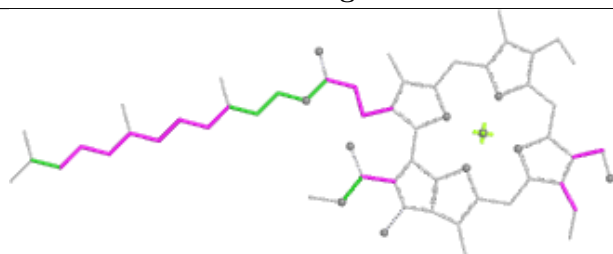
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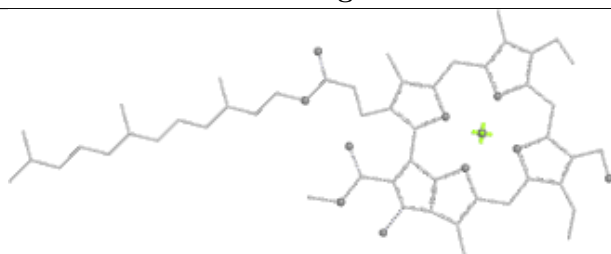
Bond lengths



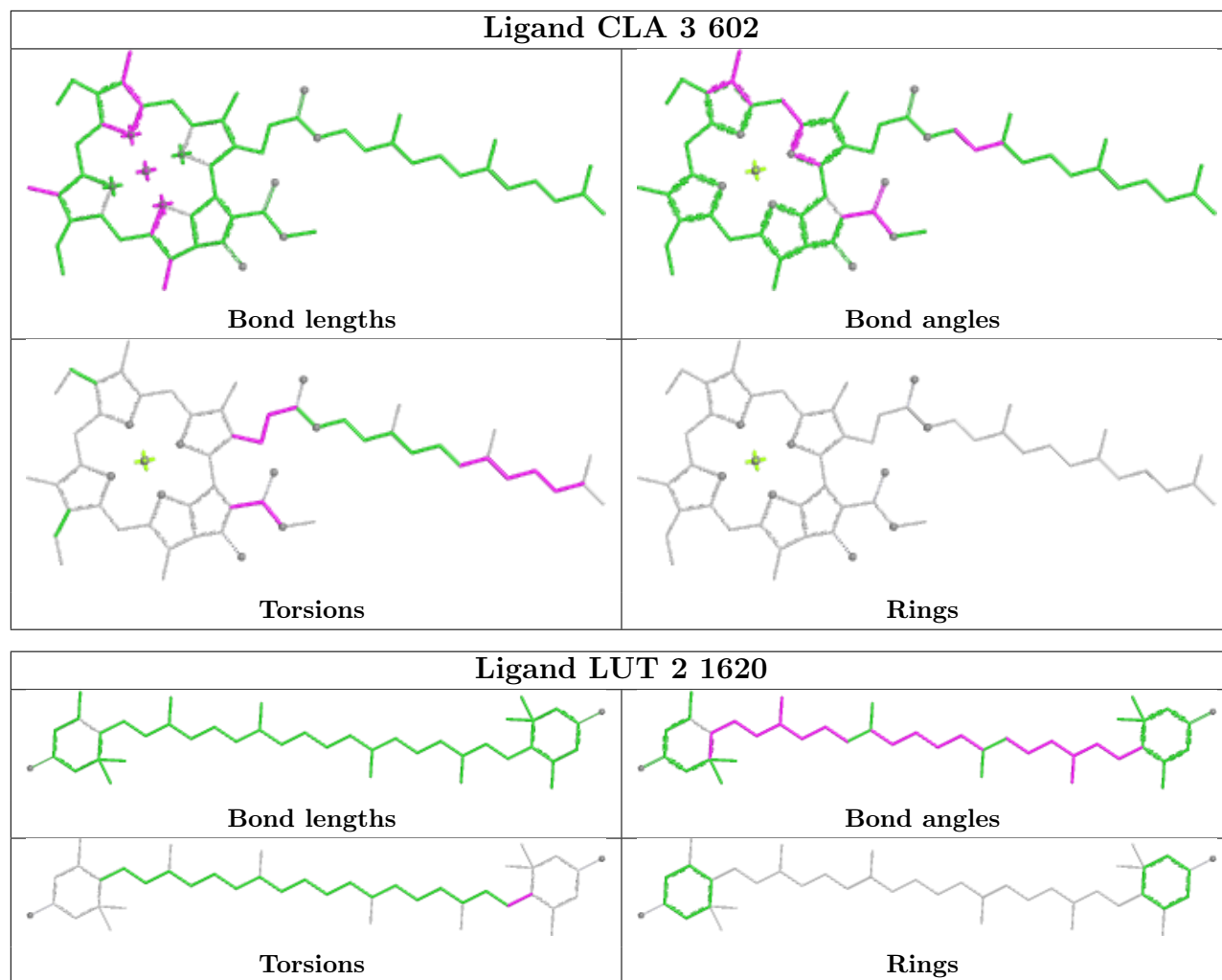
Bond angles

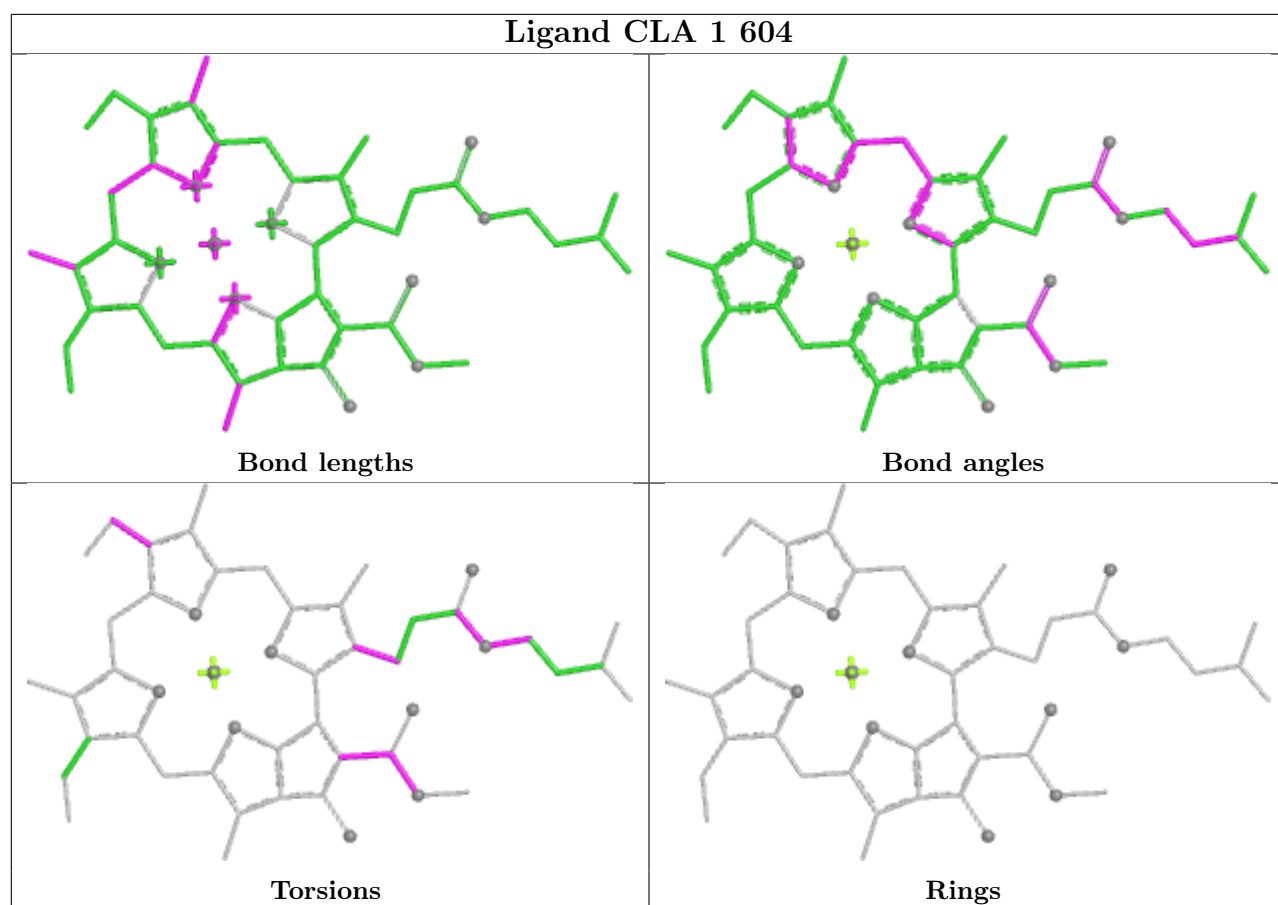


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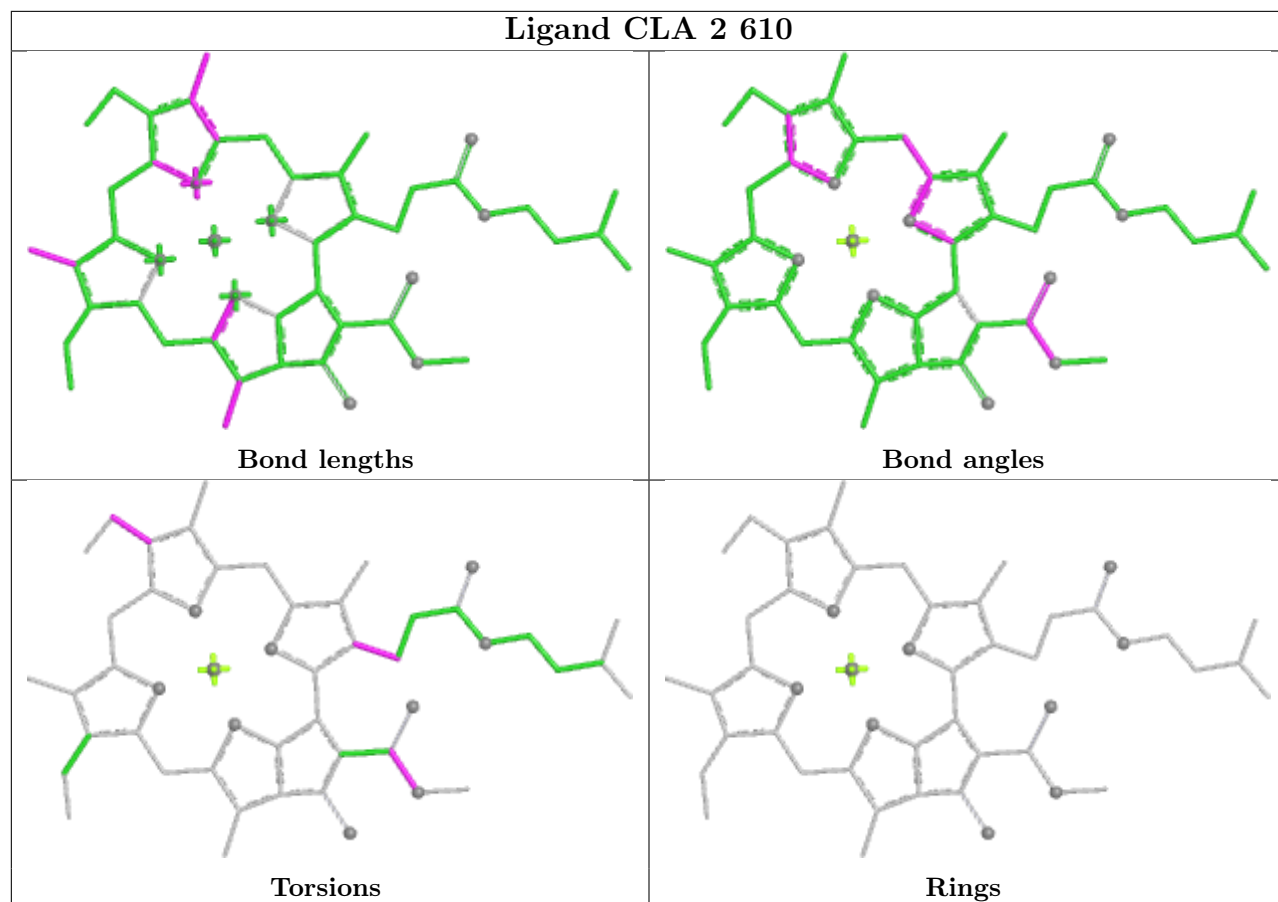


Rings

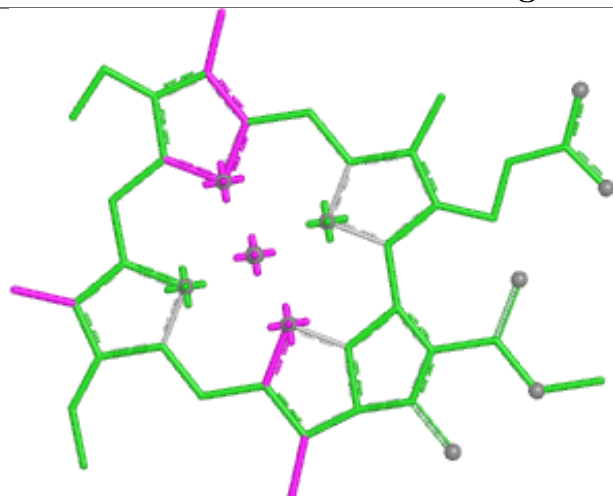




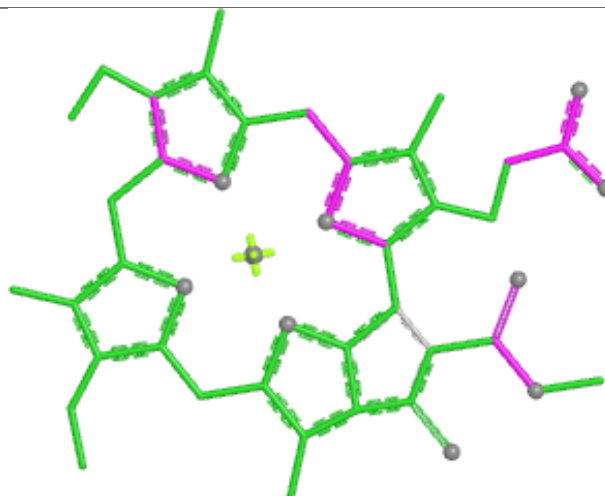
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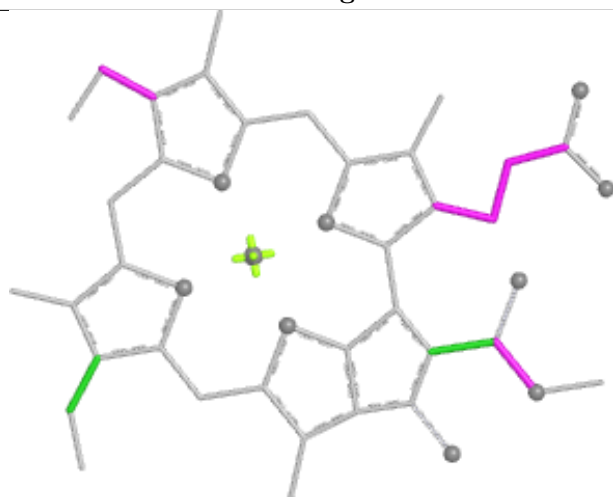
Ligand CLA 4 604



Bond lengths



Bond angles

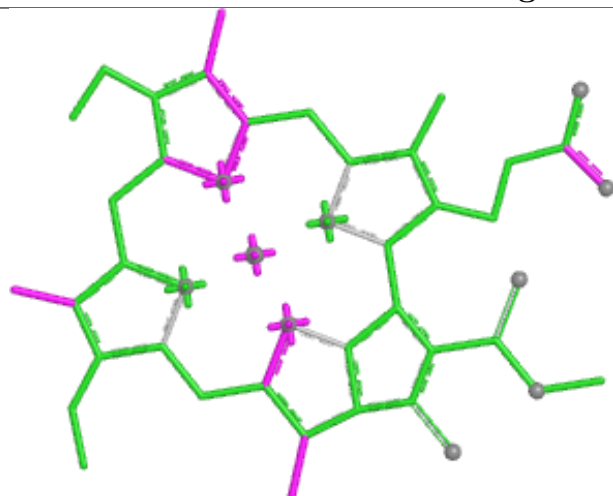


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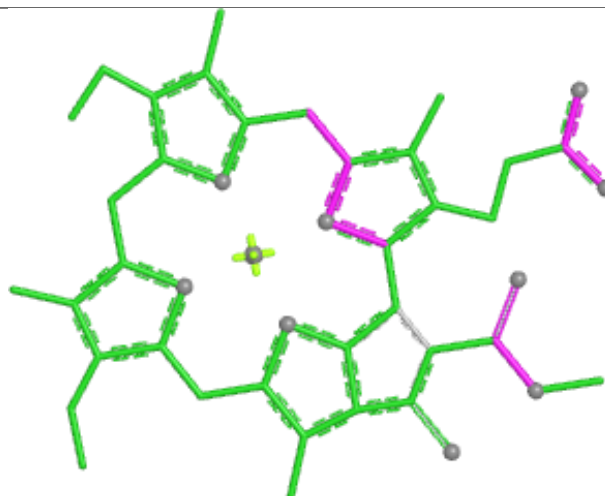


Rings

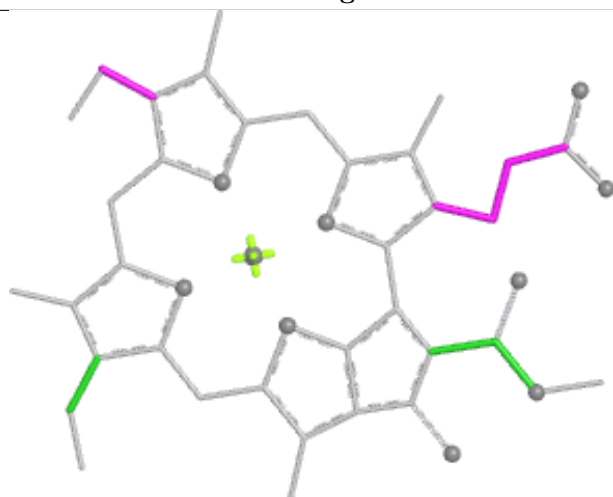
Ligand CLA 1 611



Bond lengths



Bond angles

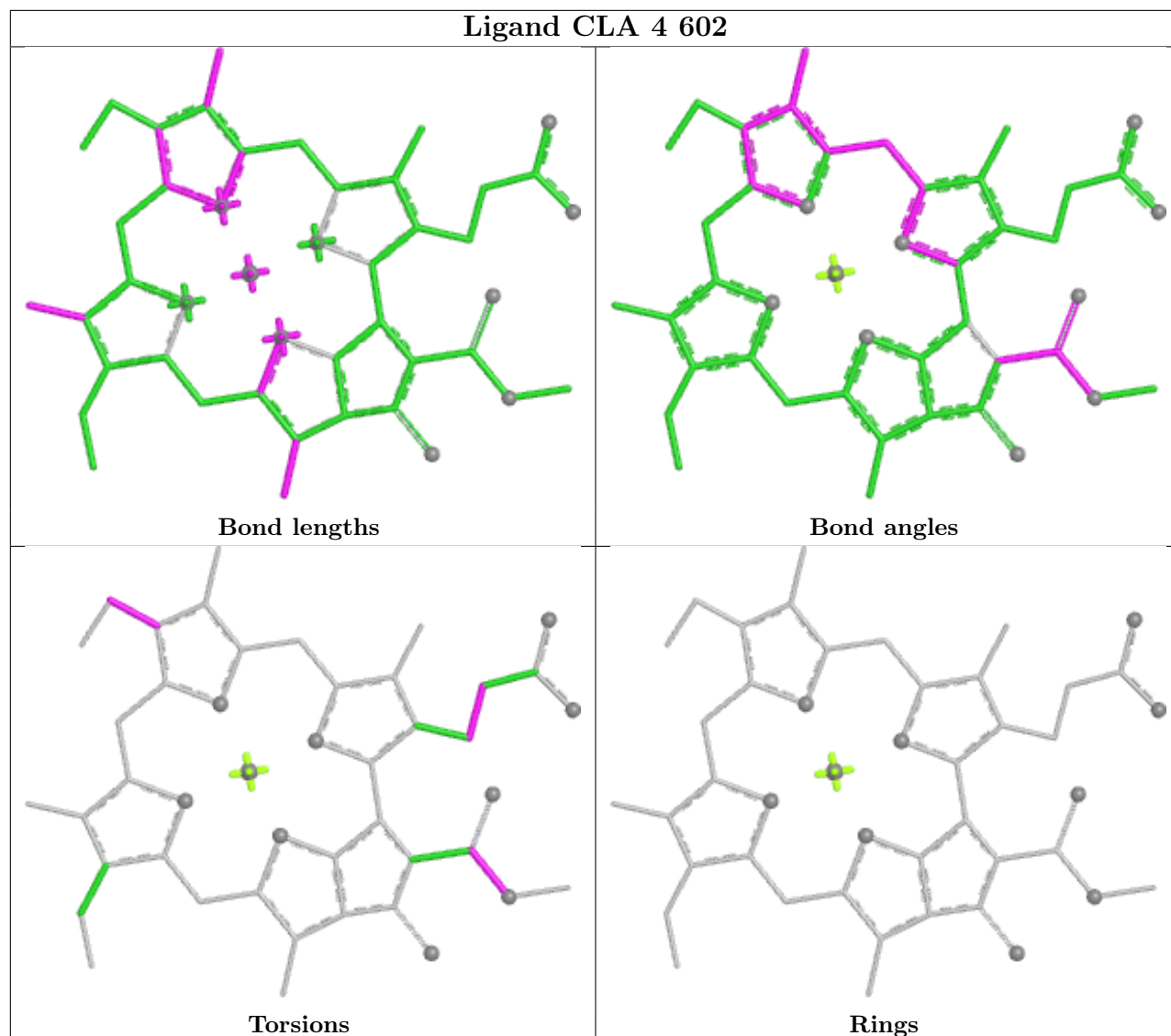


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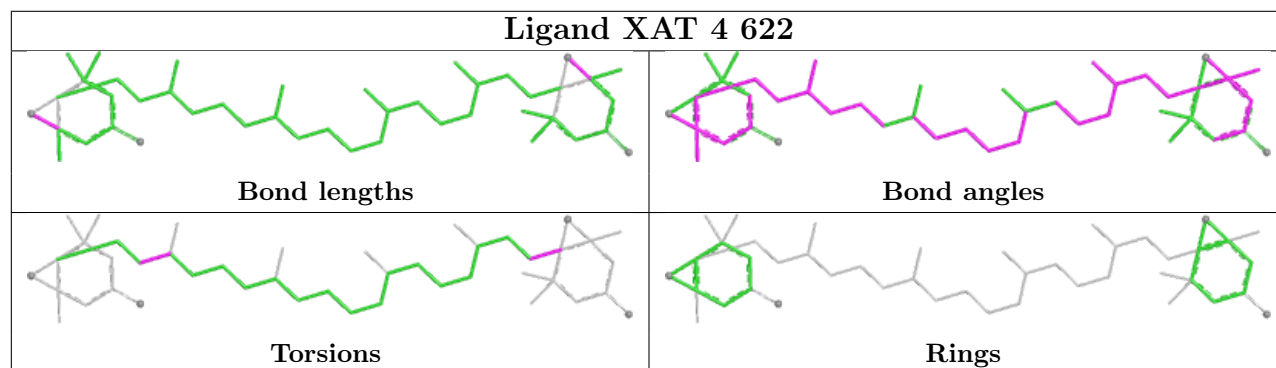


Rings

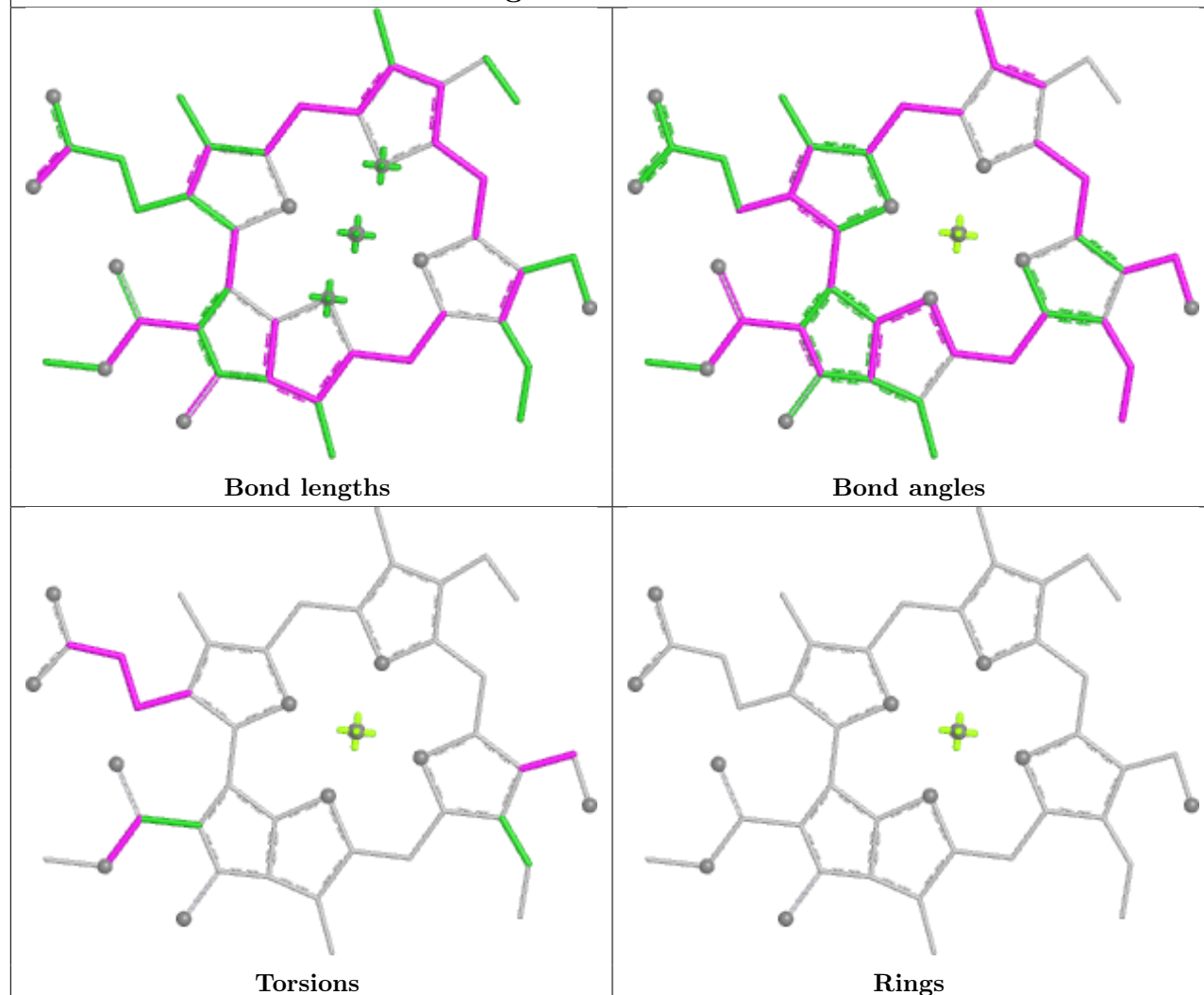
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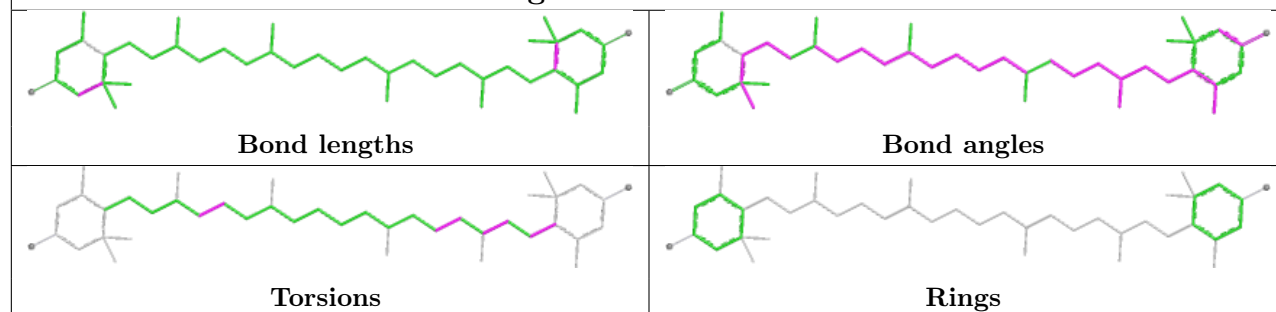
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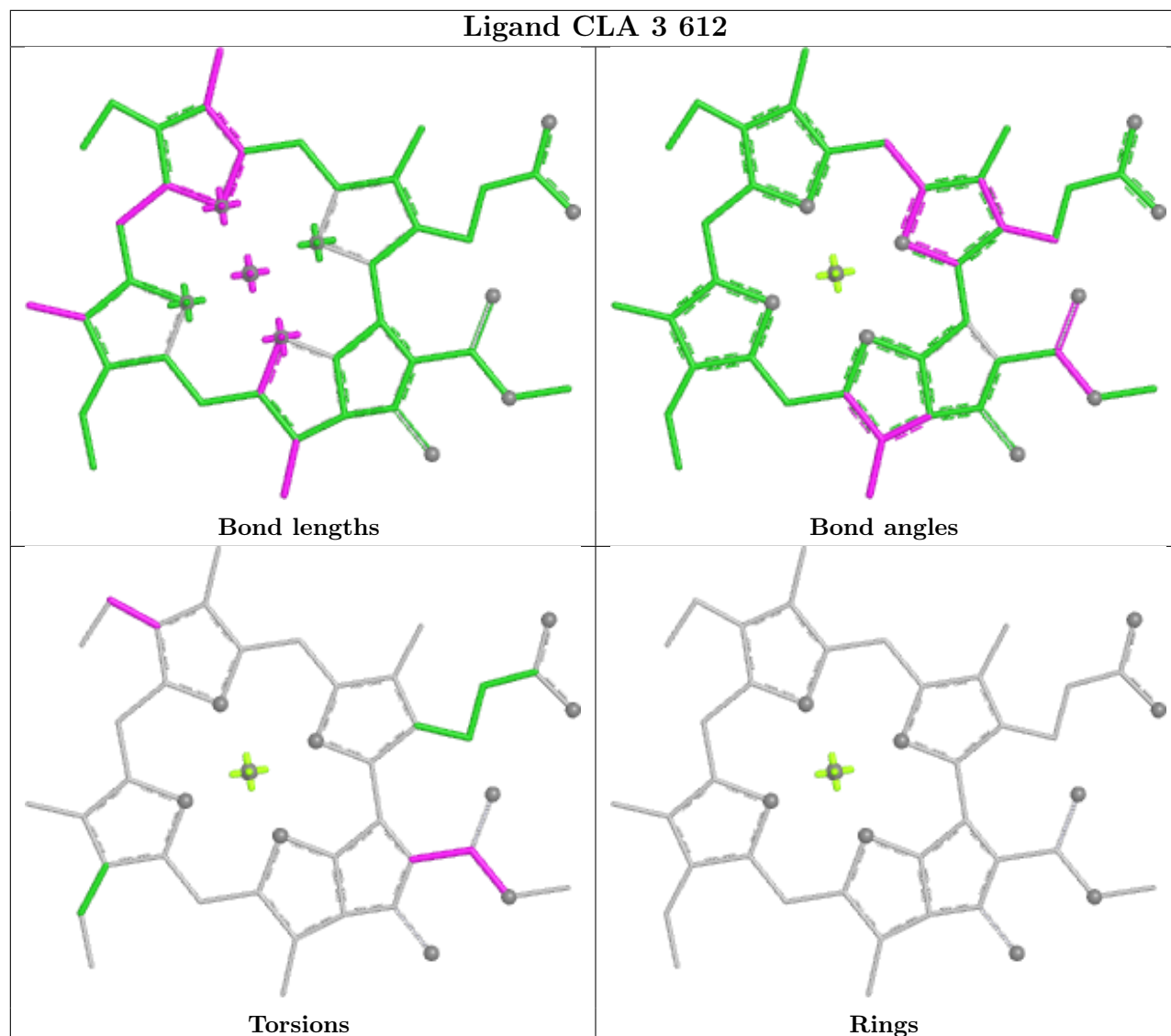
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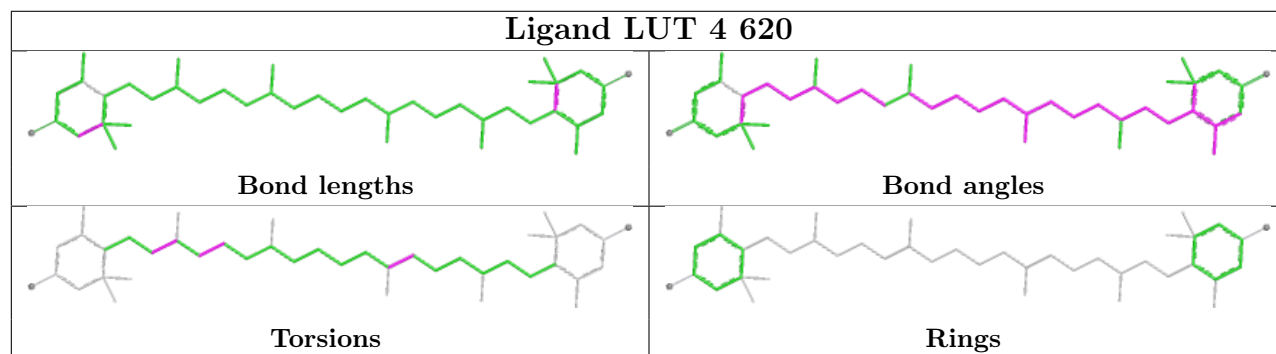
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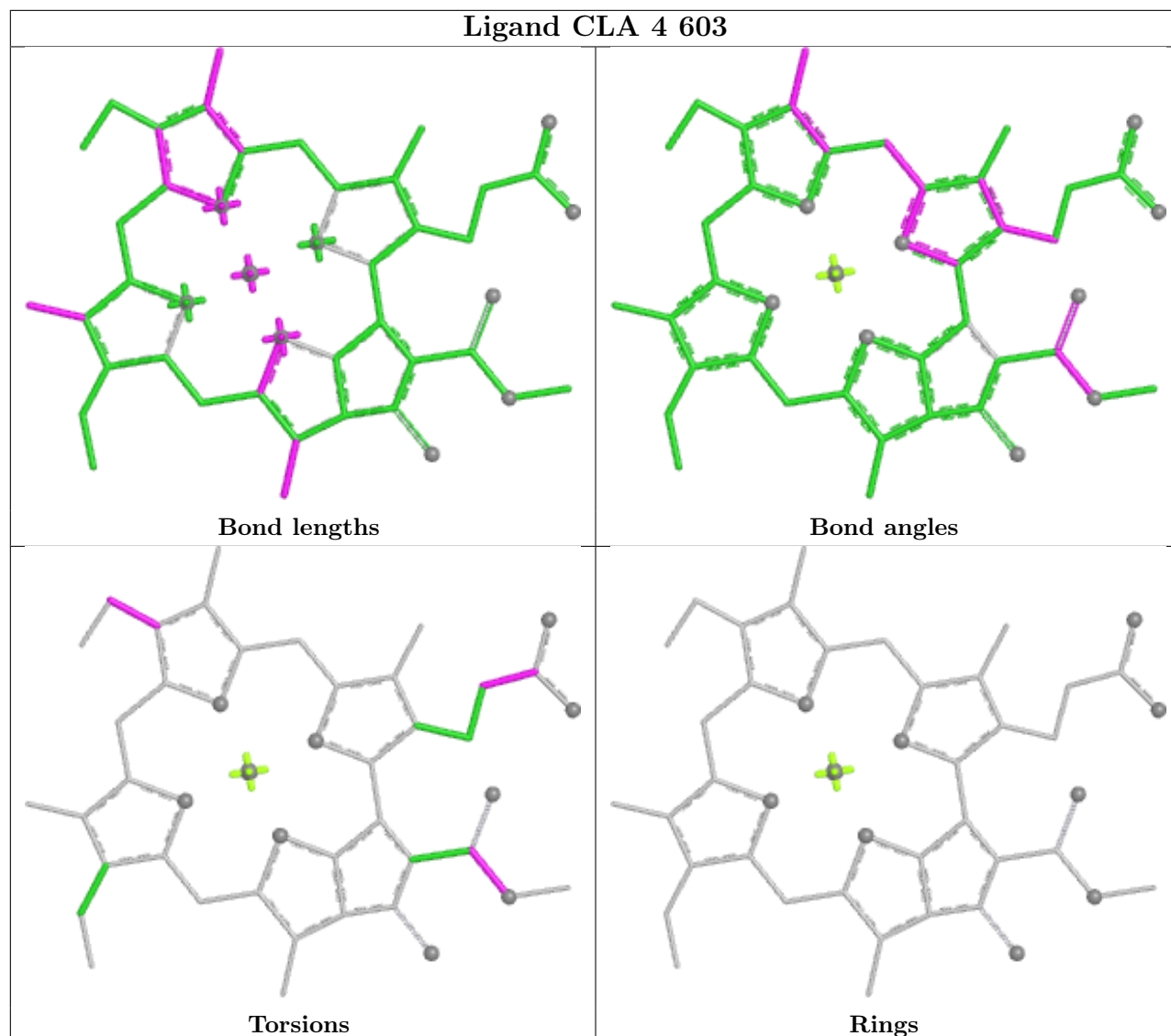
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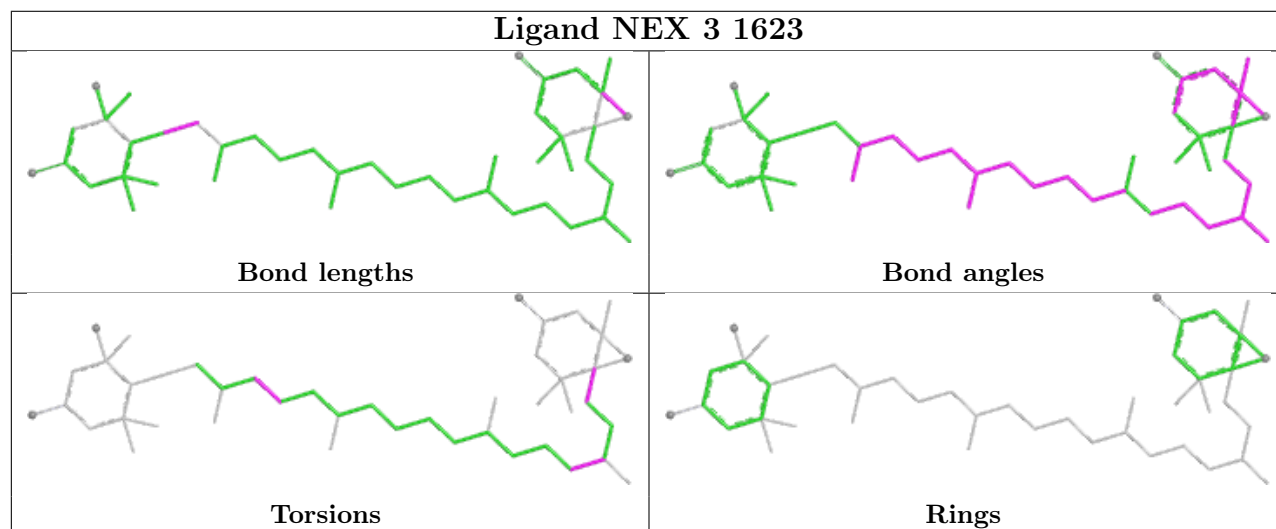
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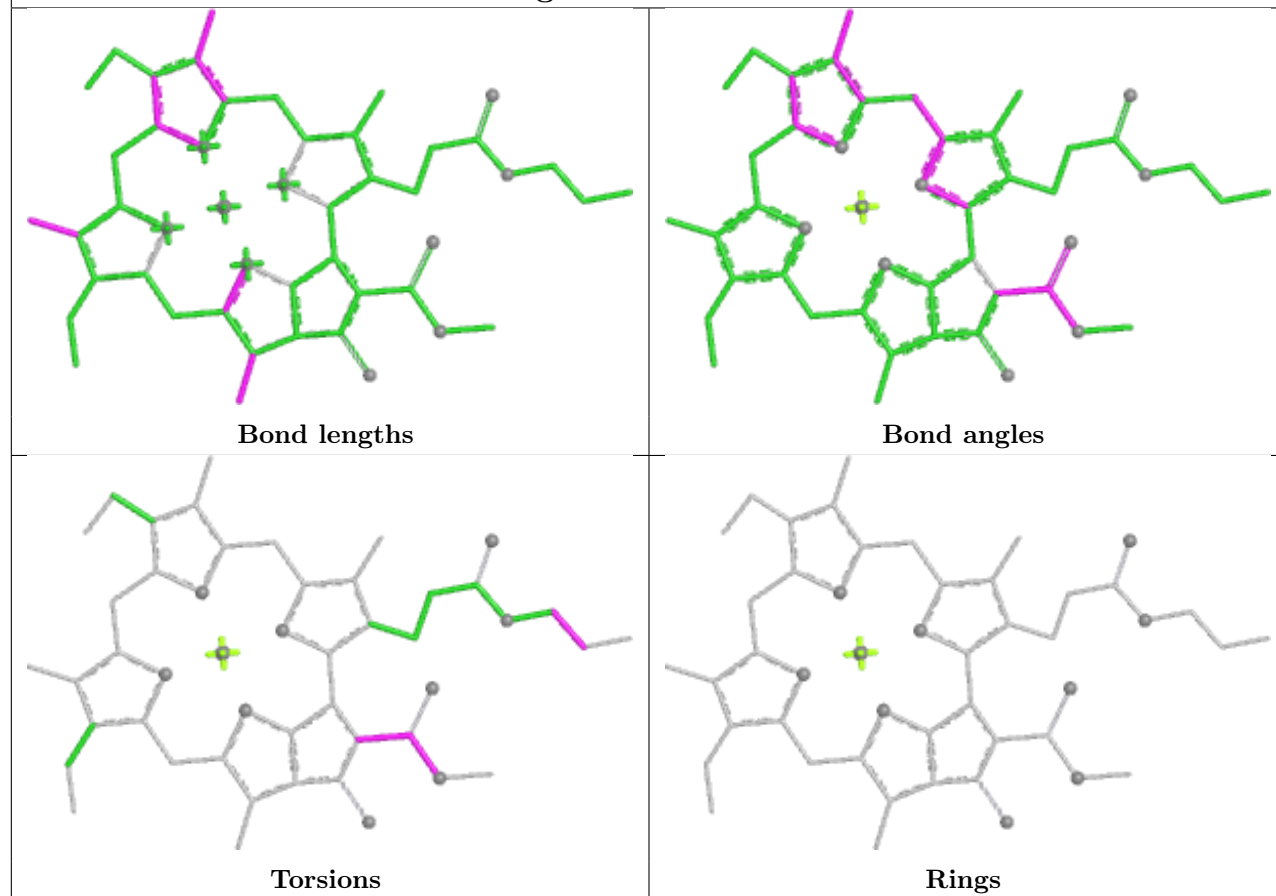
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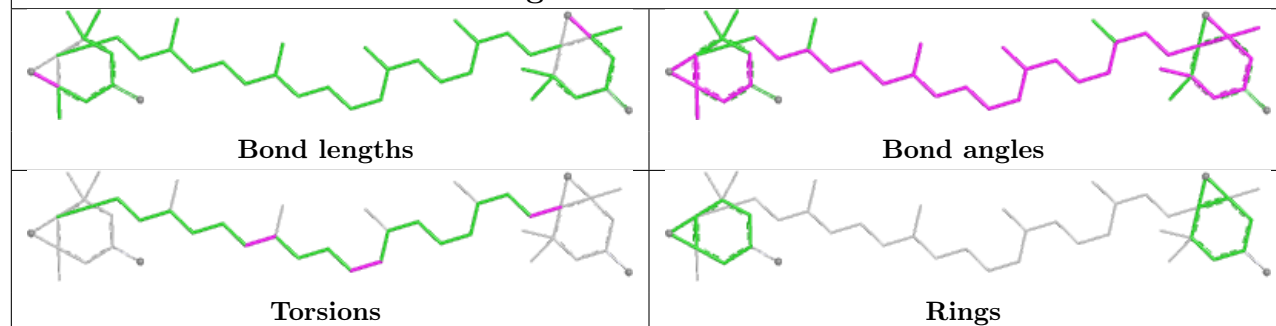
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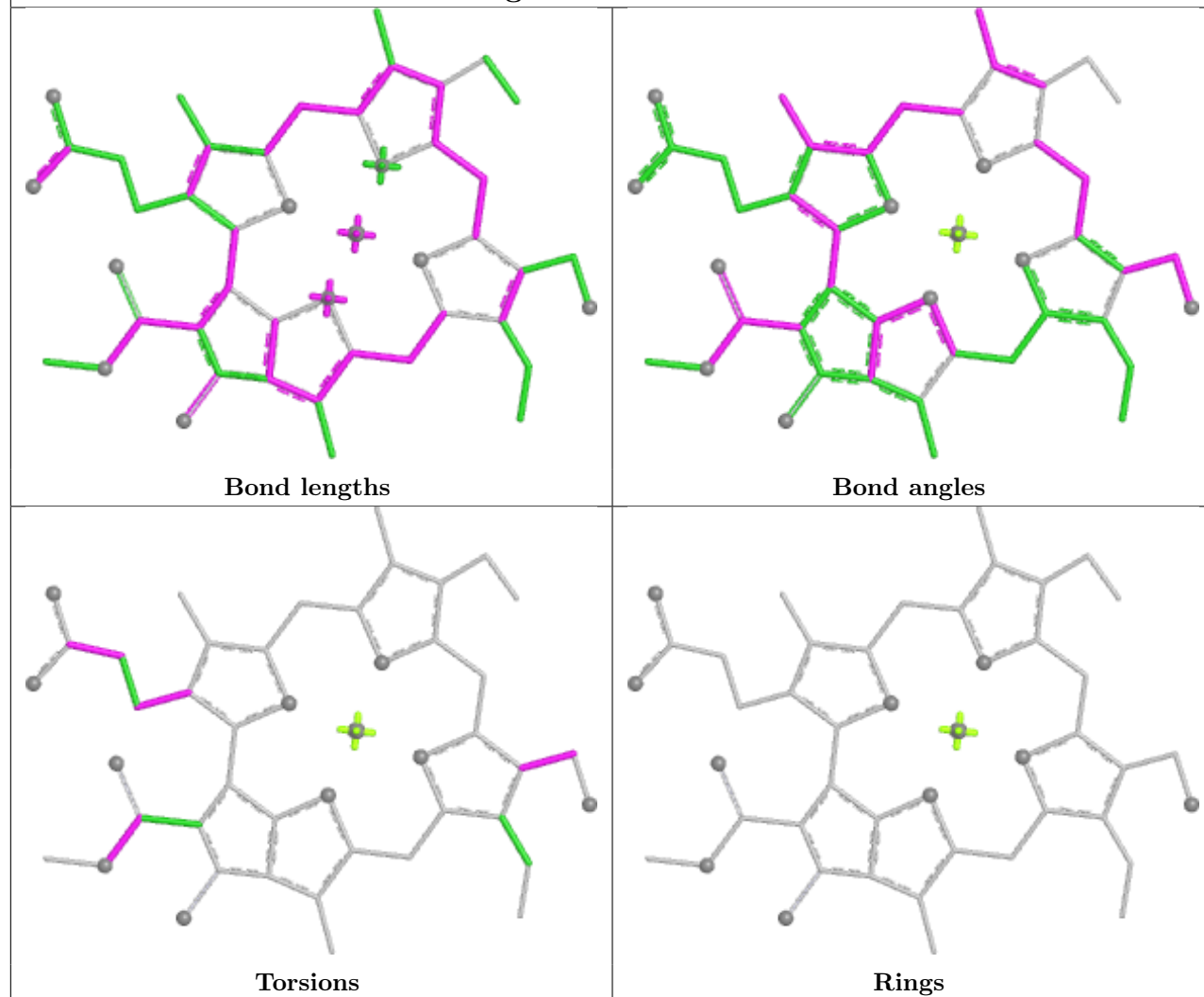
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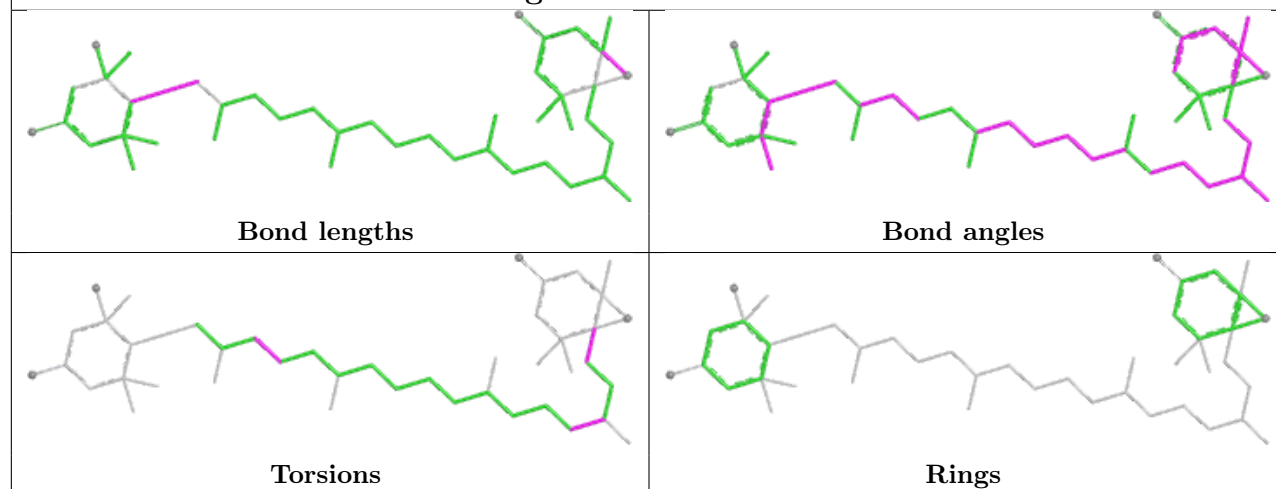
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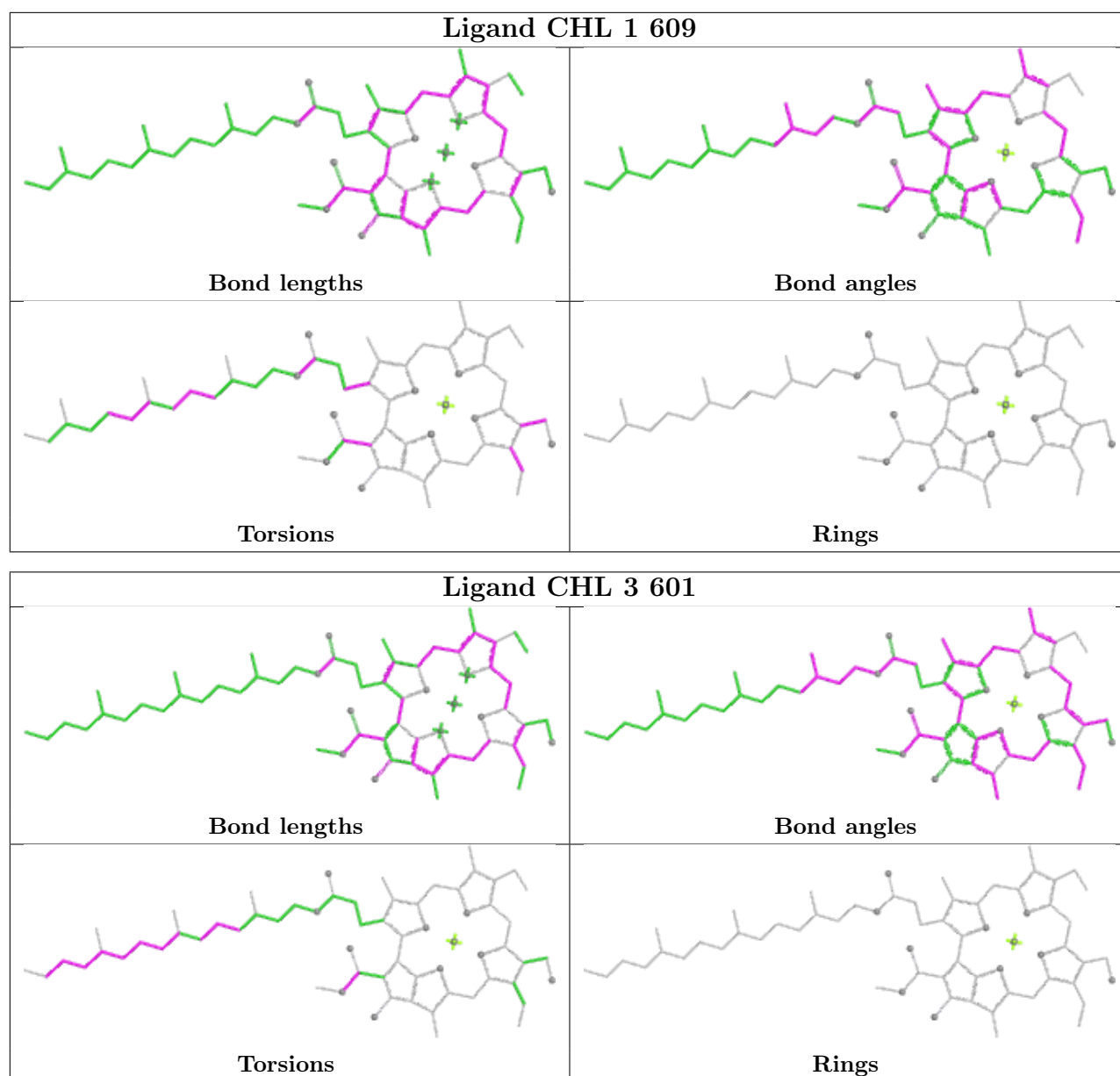


Ligand CHL 1 605

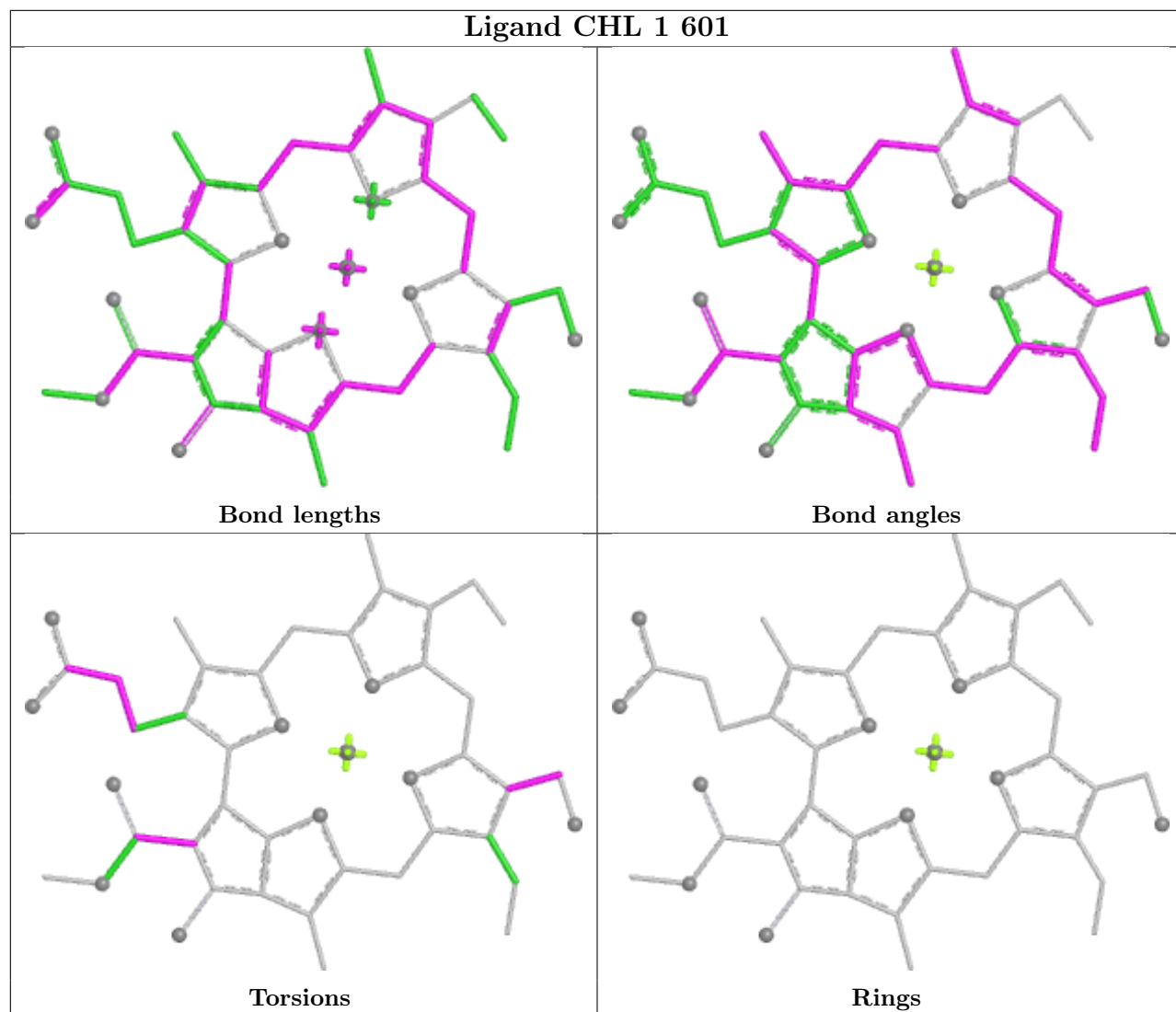


Ligand NEX 2 1623

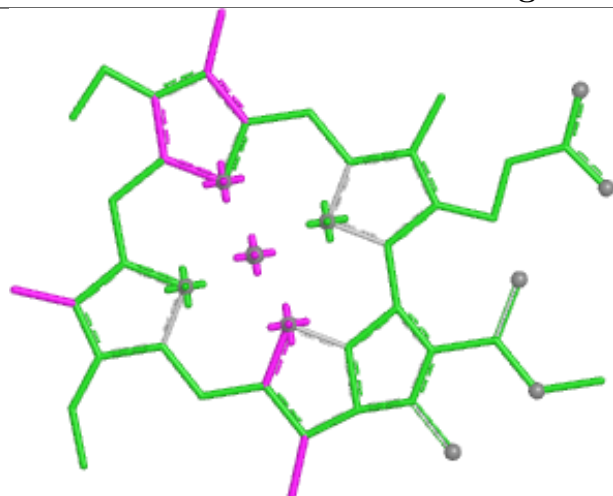




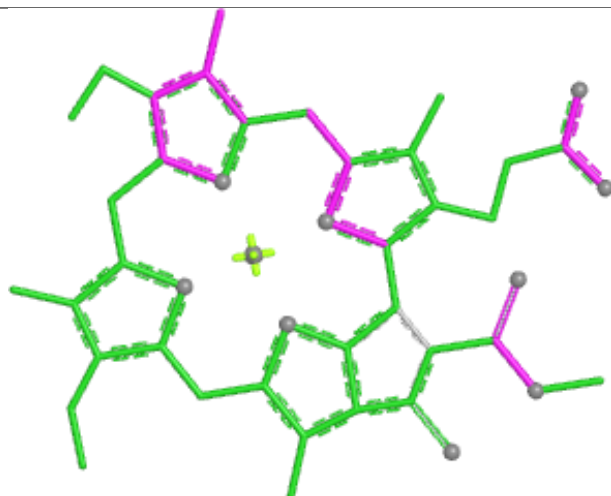
Ligand CHL 1 601



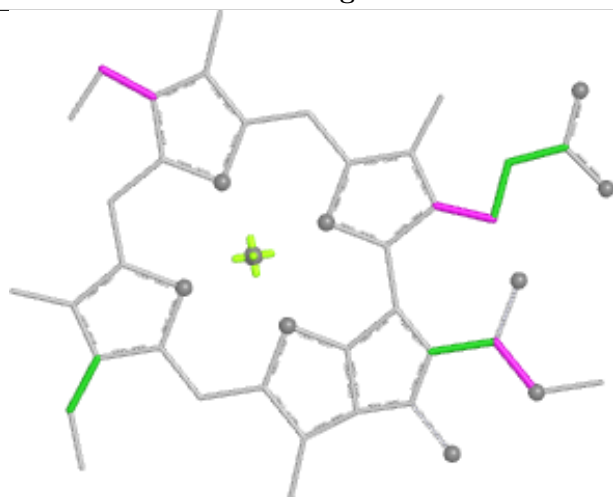
Ligand CLA 1 614



Bond lengths



Bond angles

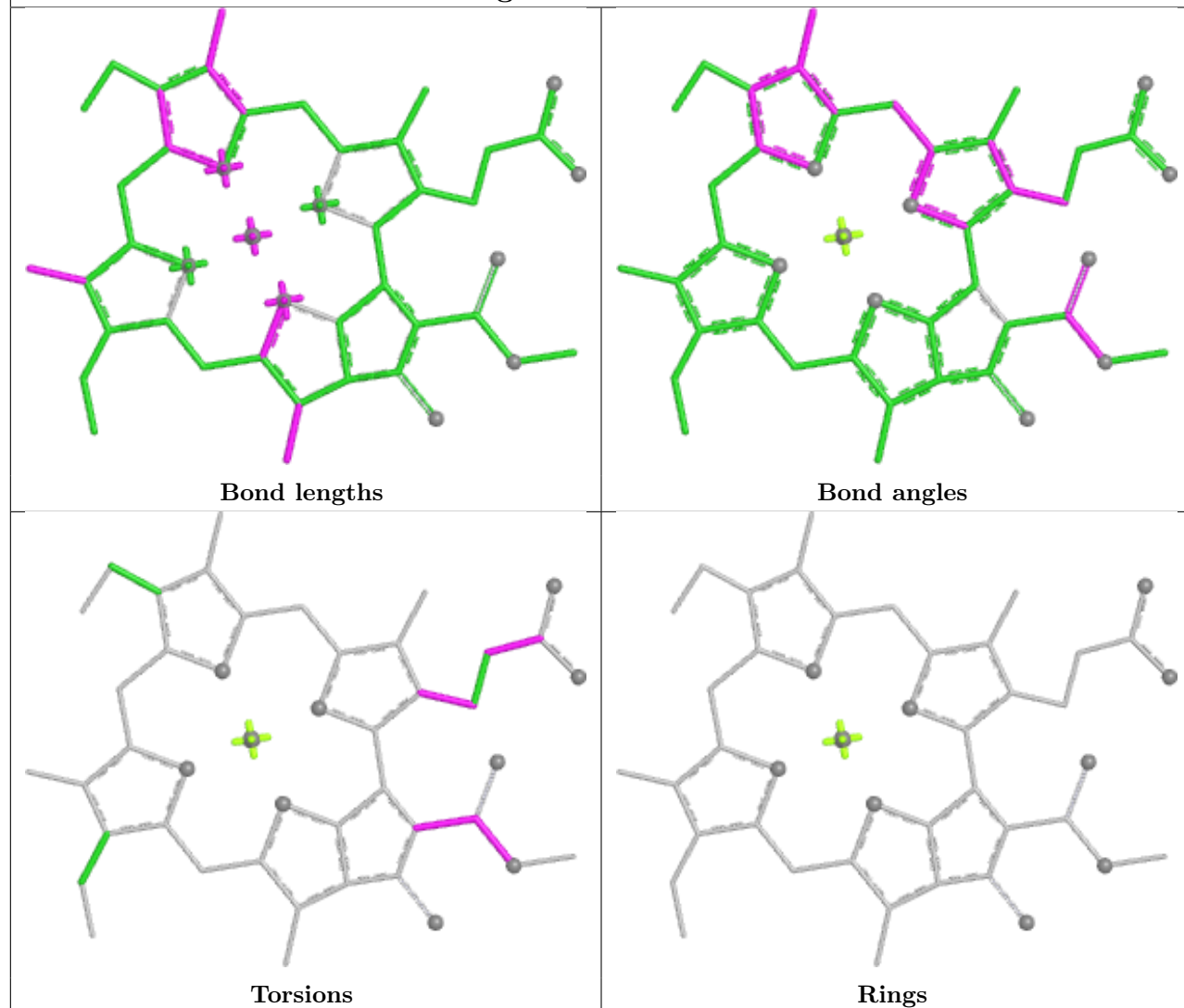


Torsions

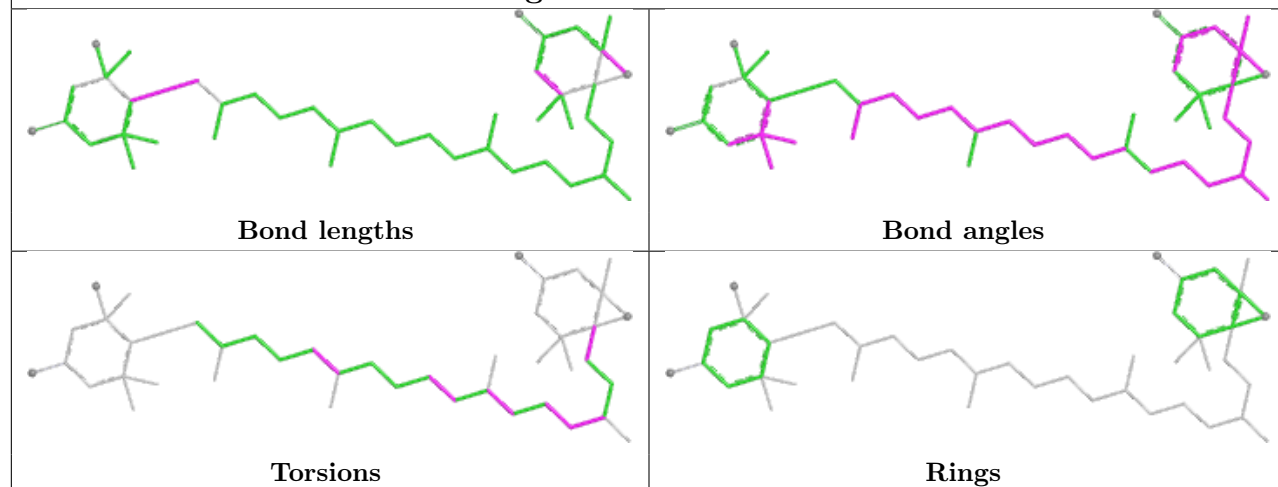


Rings

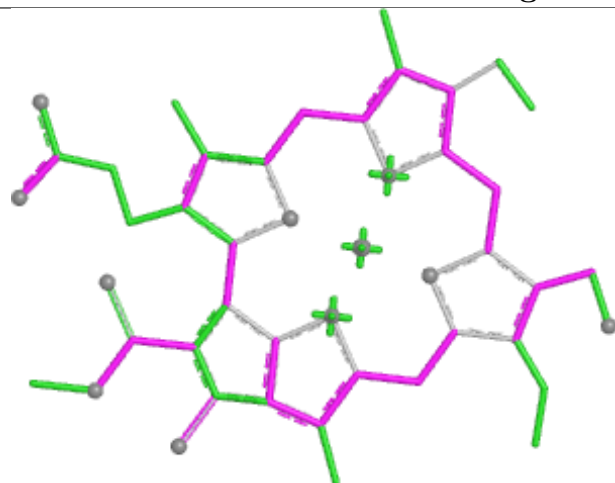
Ligand CLA 4 612



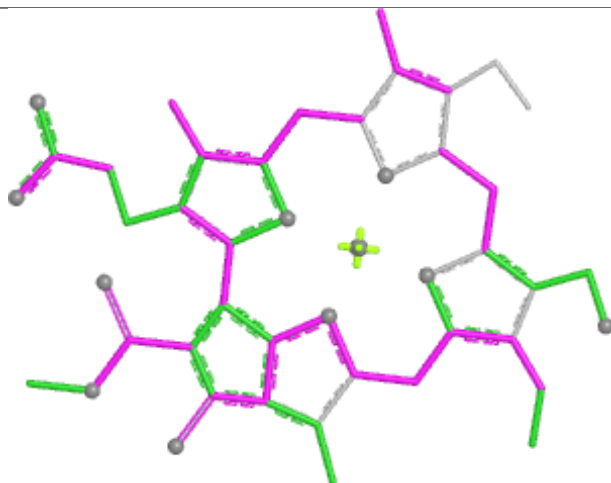
Ligand NEX 1 1623



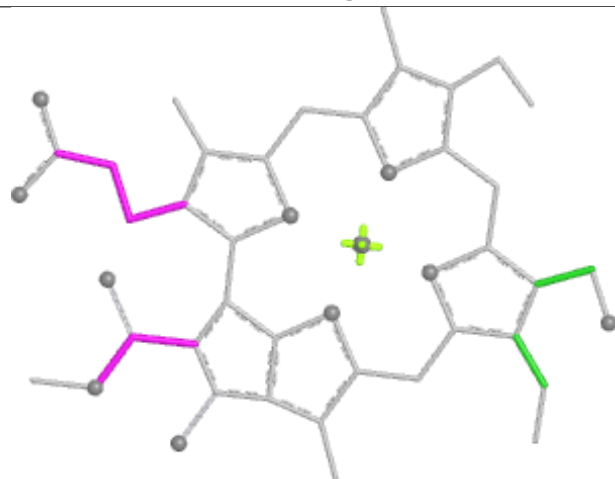
Ligand CHL 3 608



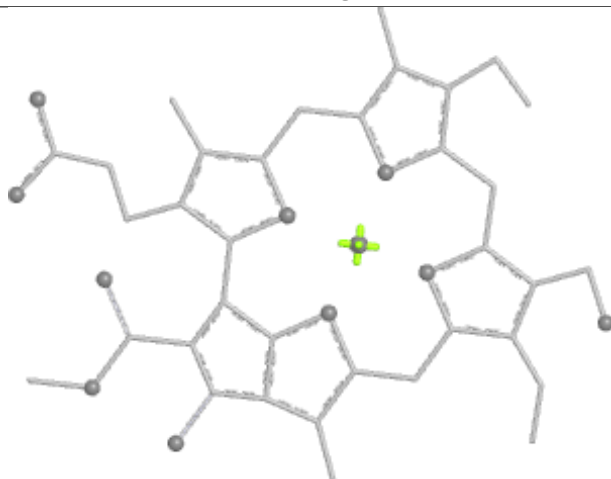
Bond lengths



Bond angles

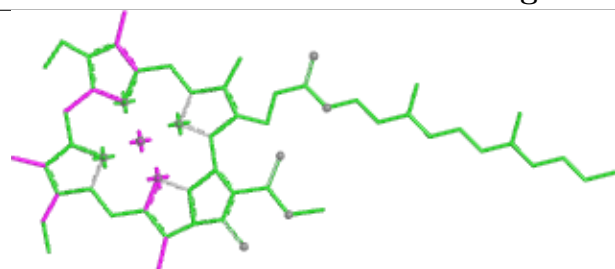


Torsions

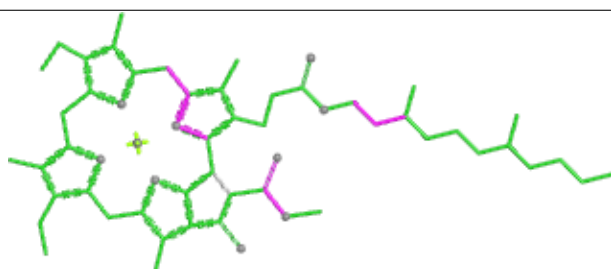


Rings

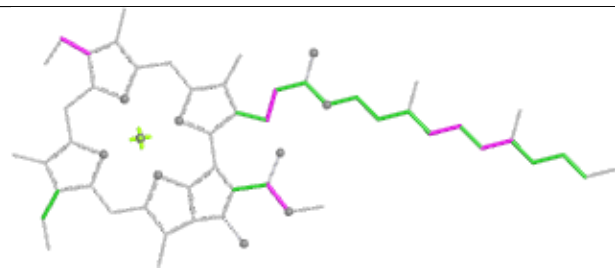
Ligand CLA 3 613



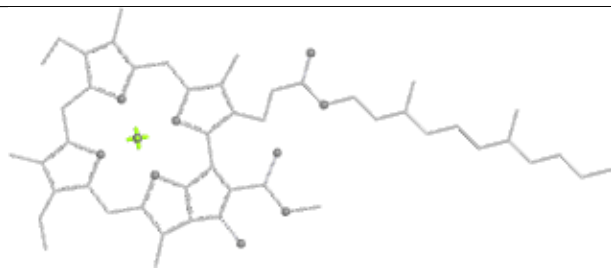
Bond lengths



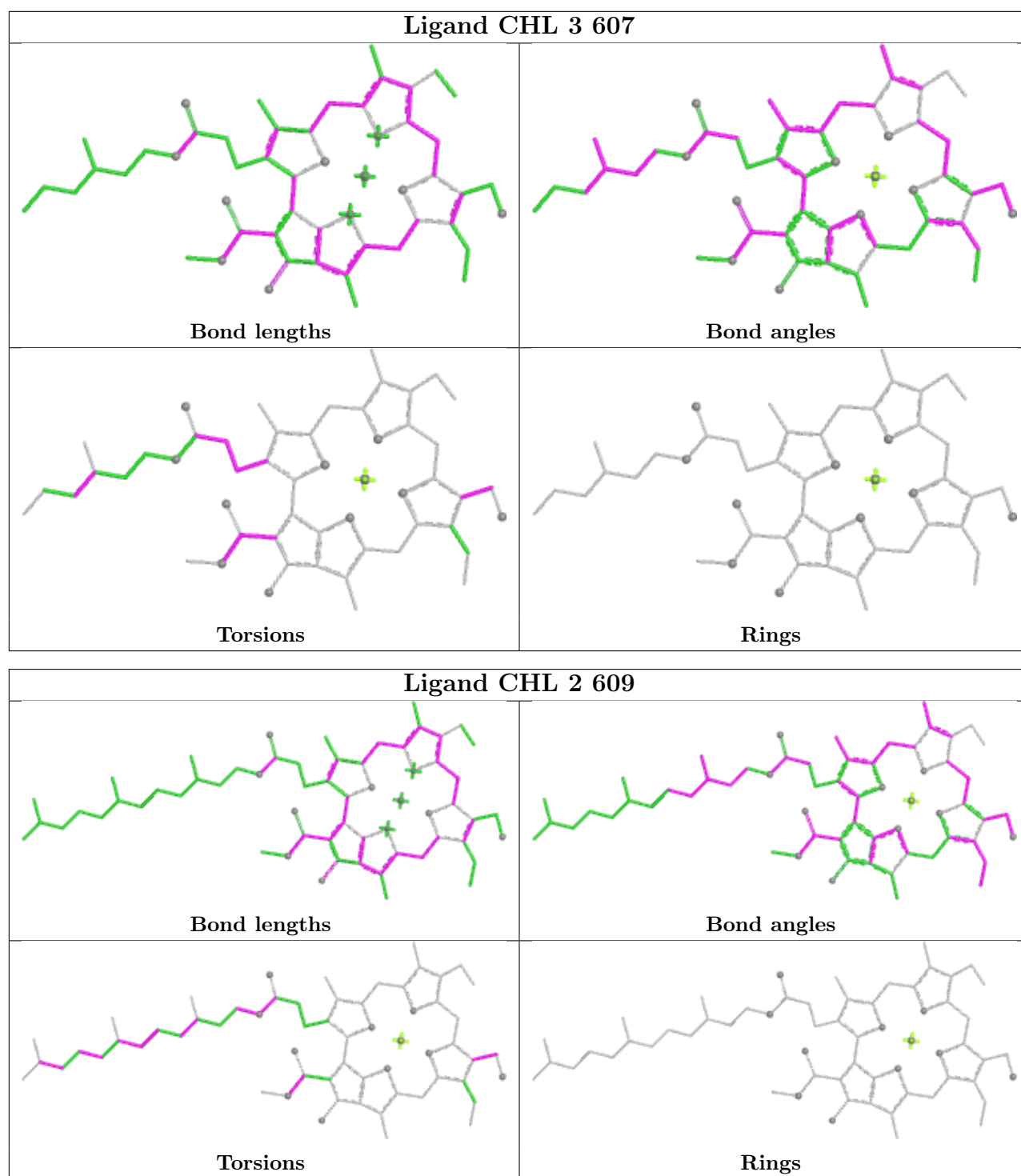
Bond angles



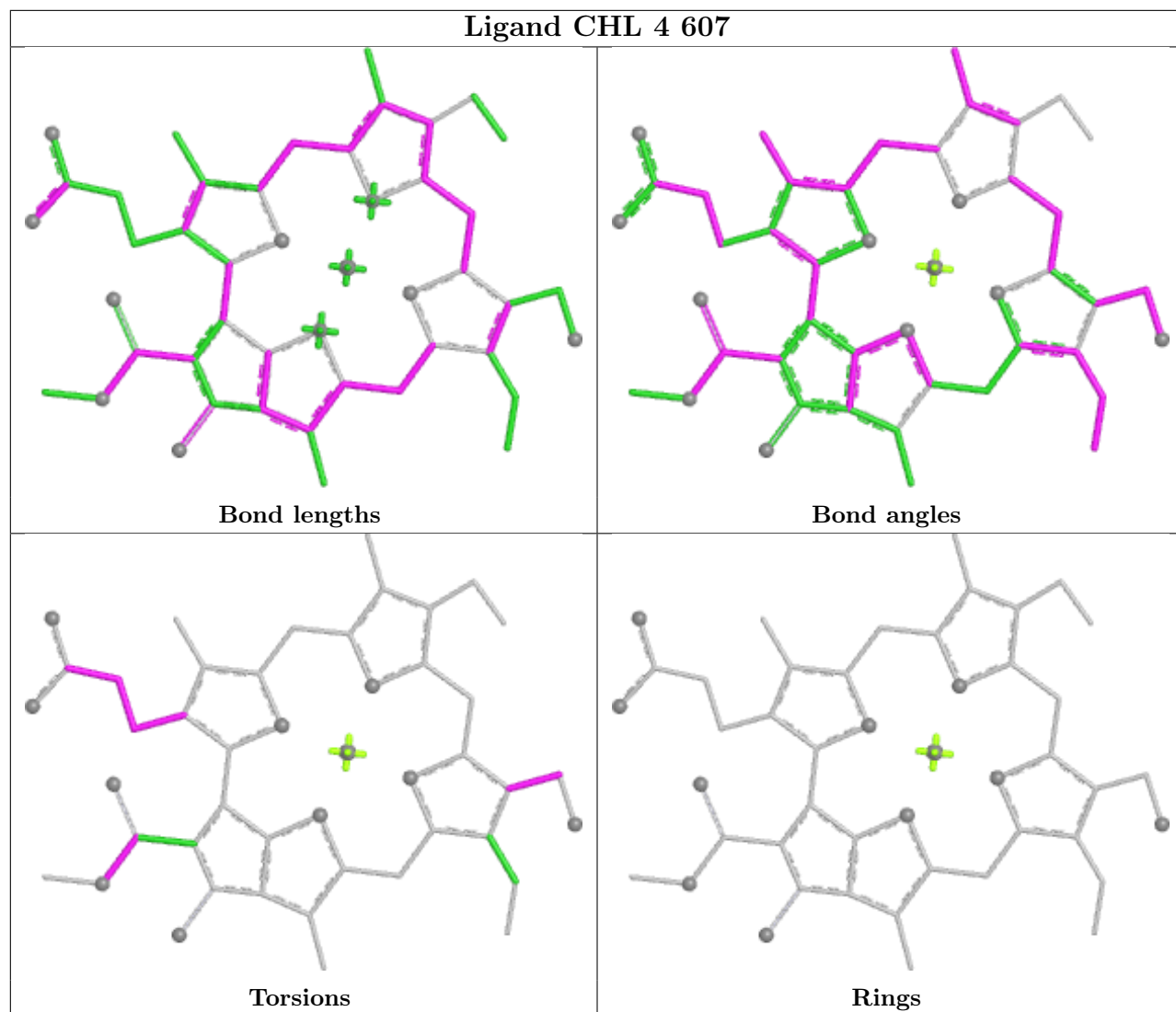
Torsions

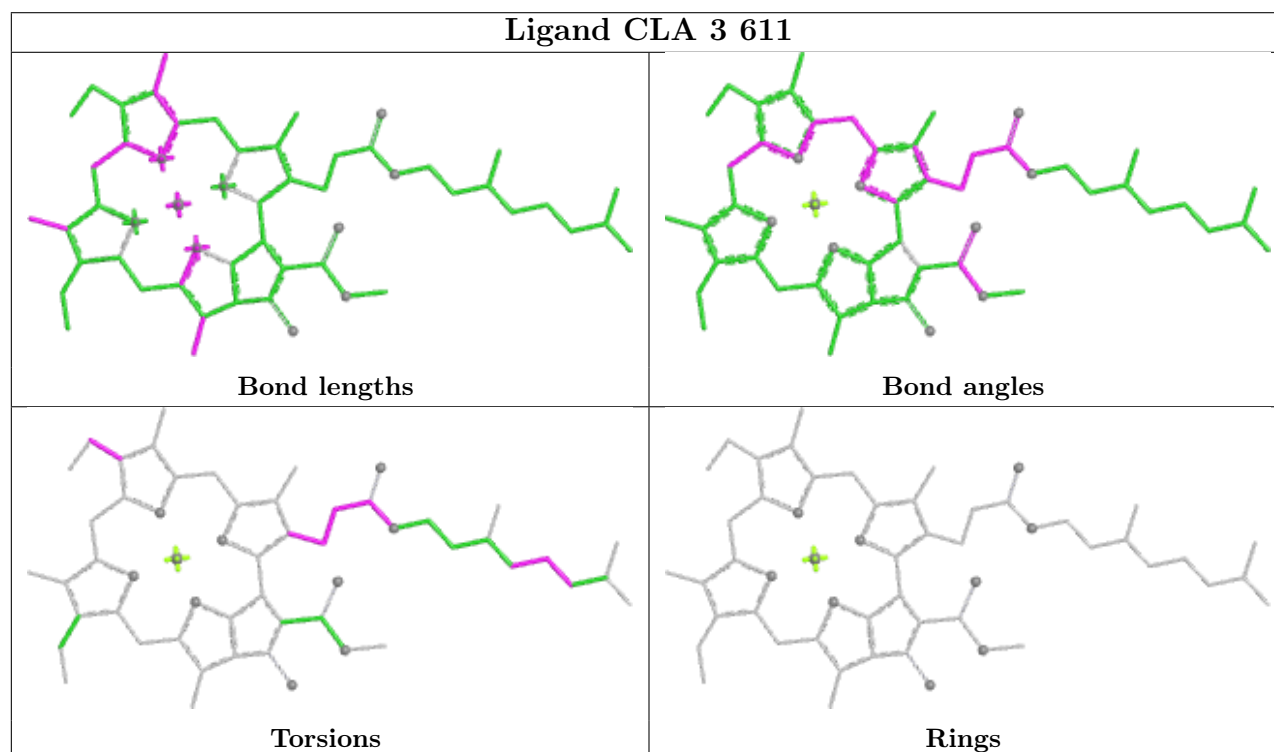
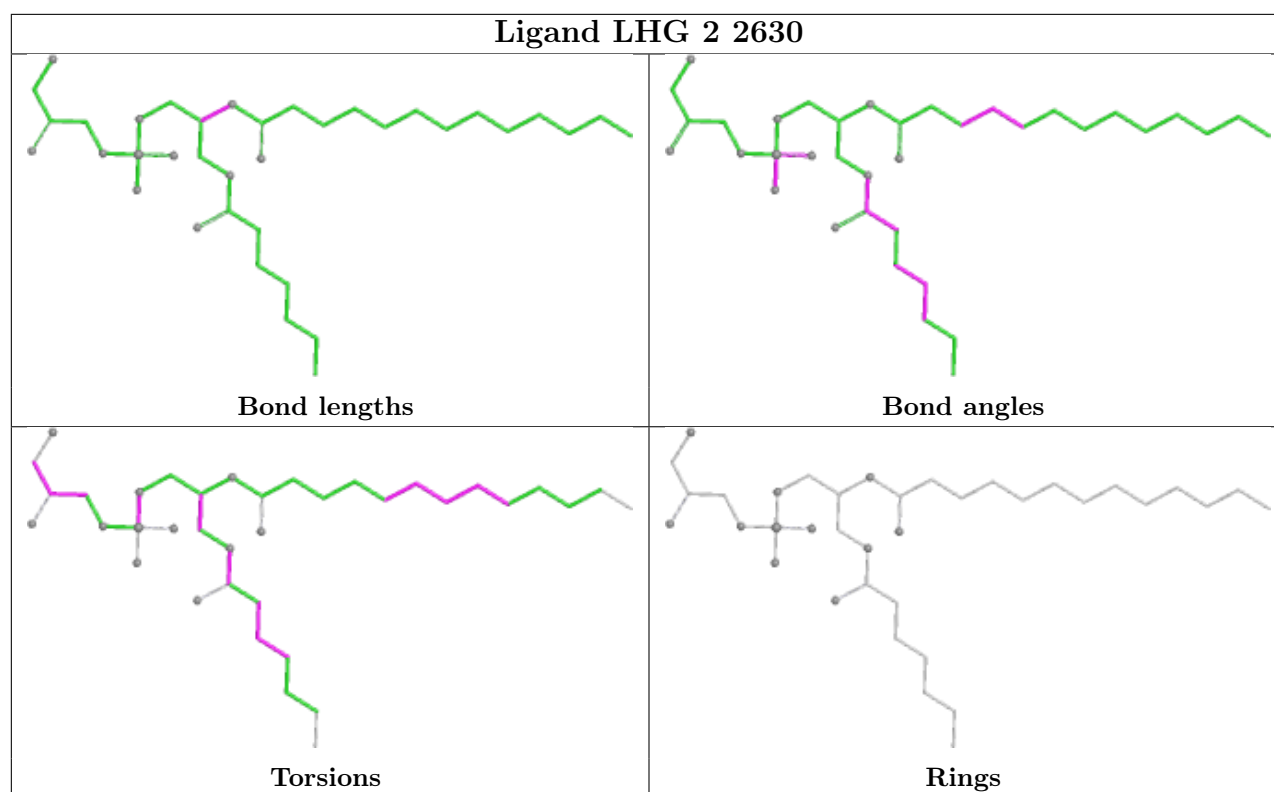


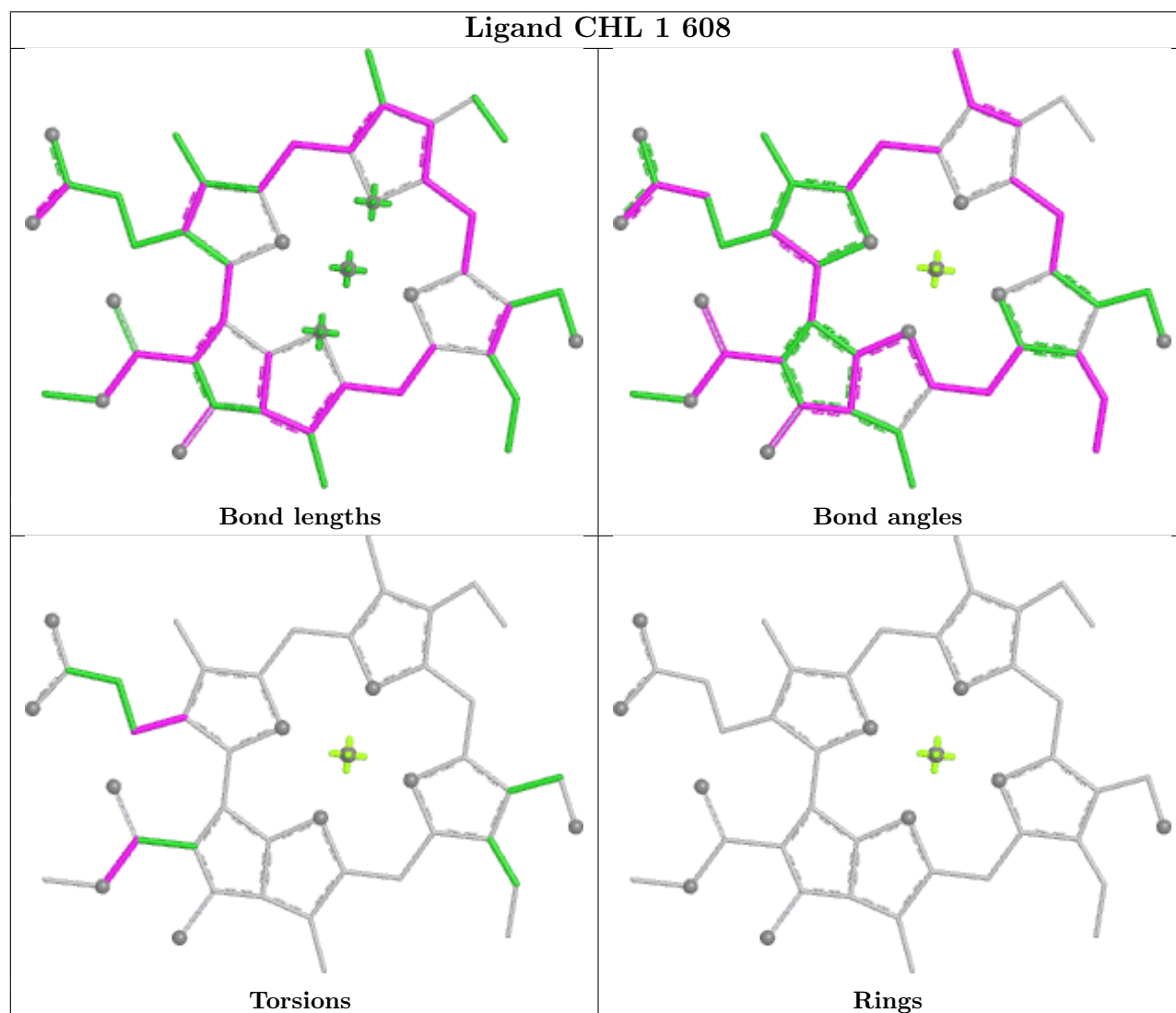
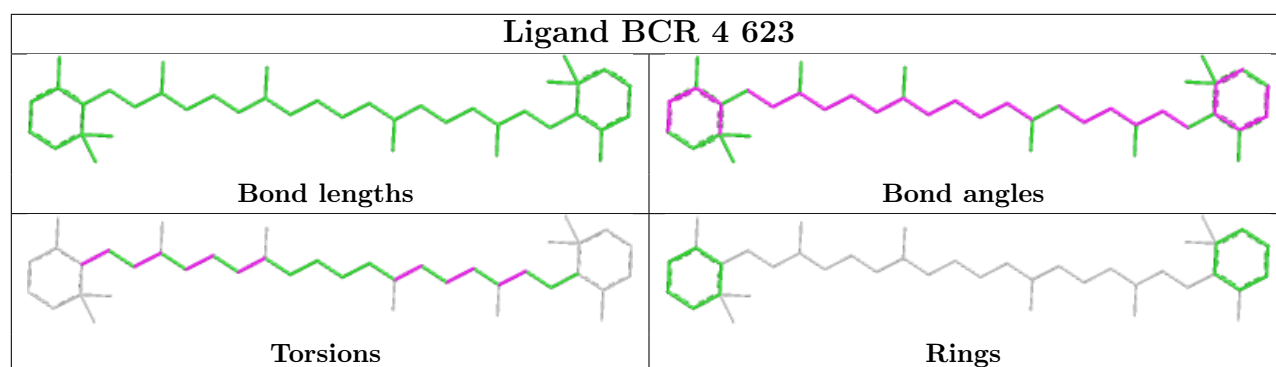
Rings



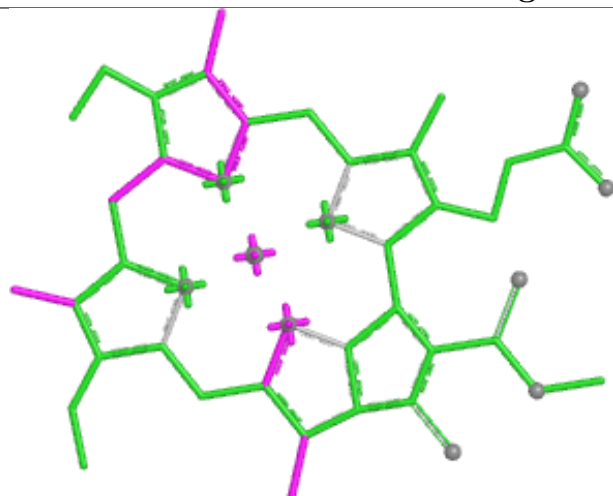
Ligand CHL 4 607



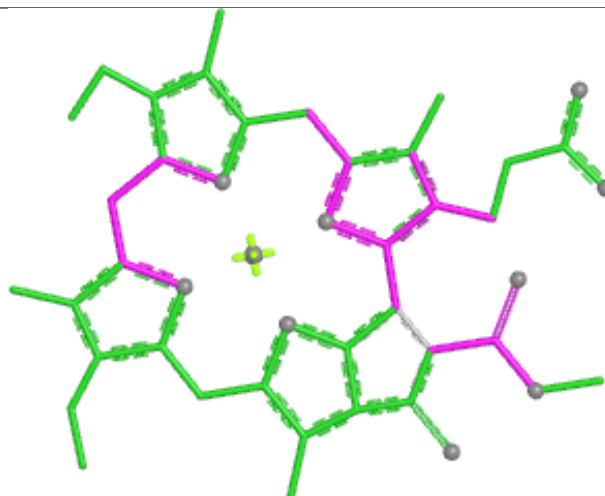




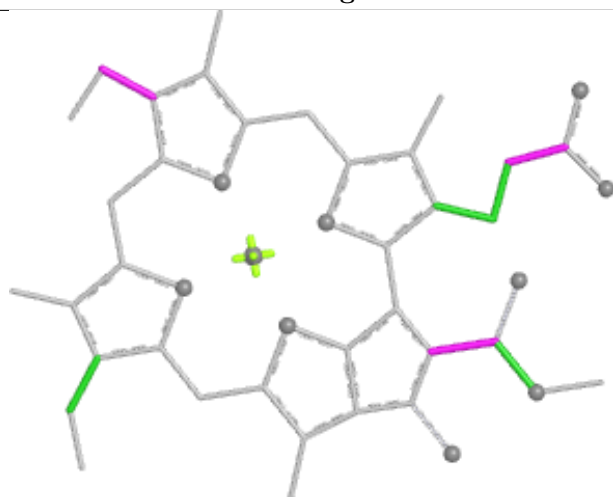
Ligand CLA 1 612



Bond lengths



Bond angles

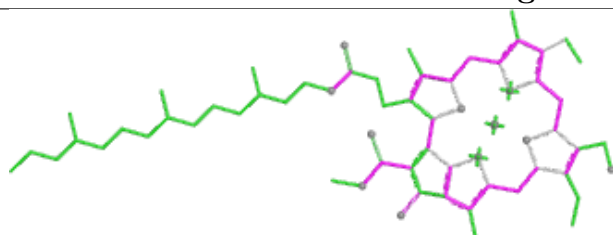


Torsions

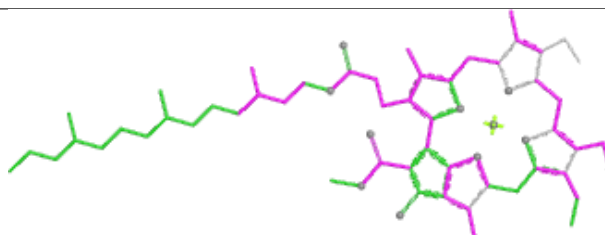


Rings

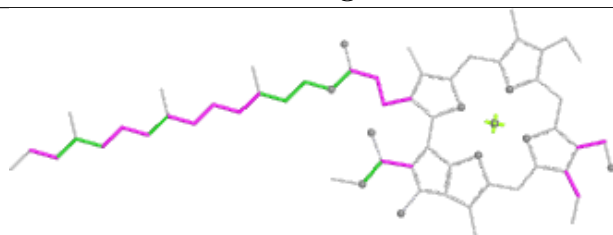
Ligand CHL 1 607



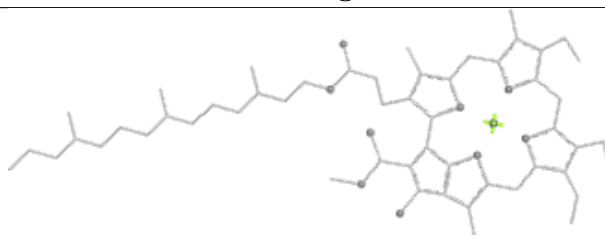
Bond lengths



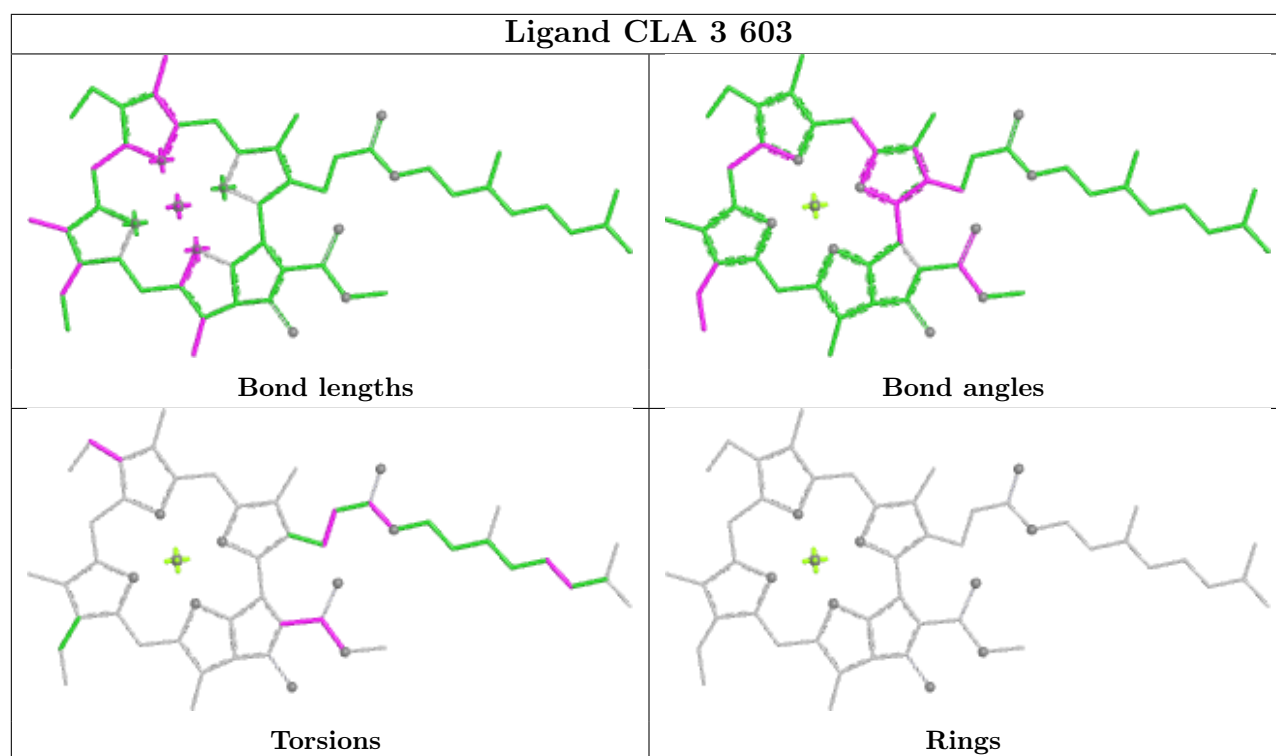
Bond angles



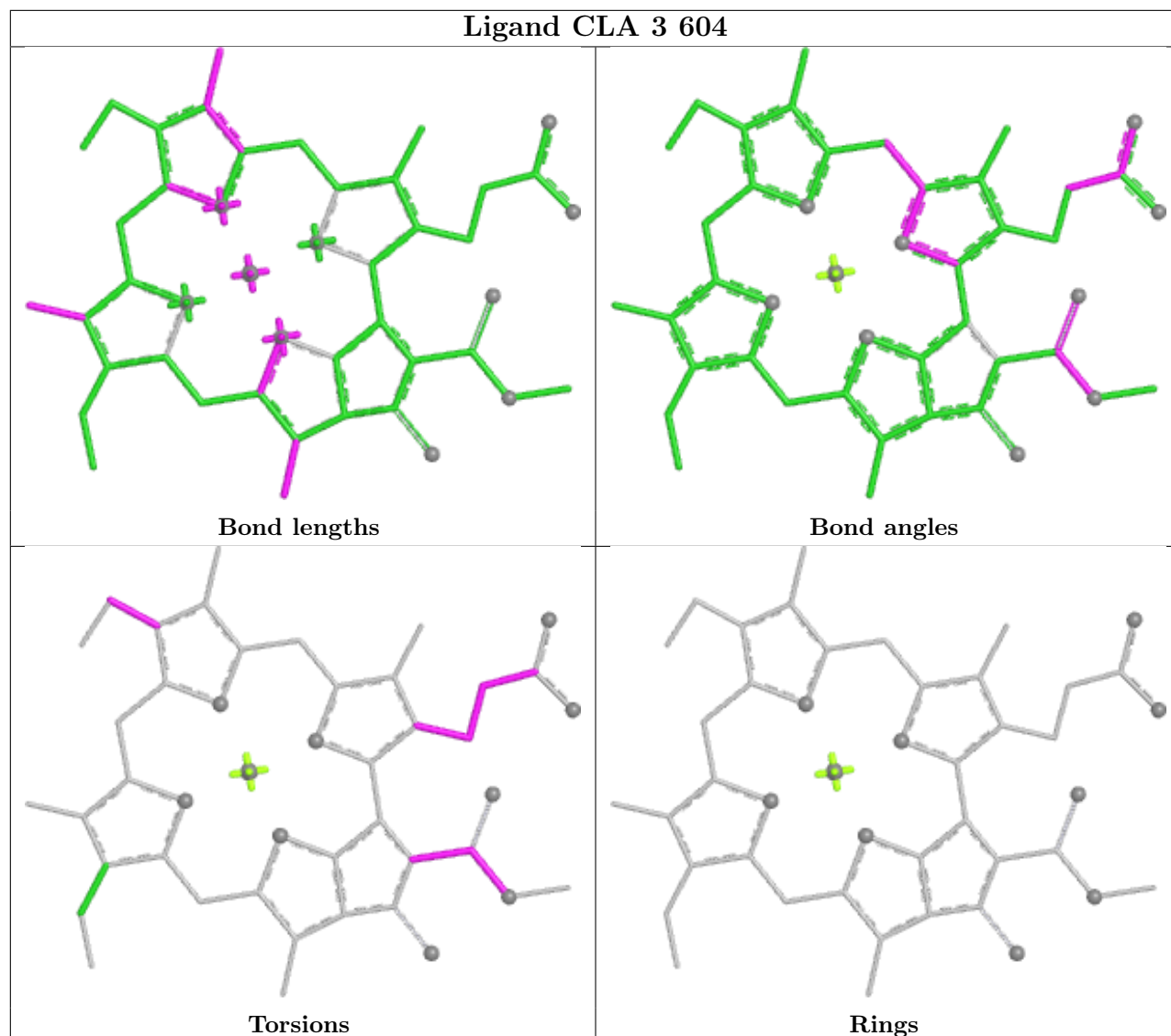
Torsions

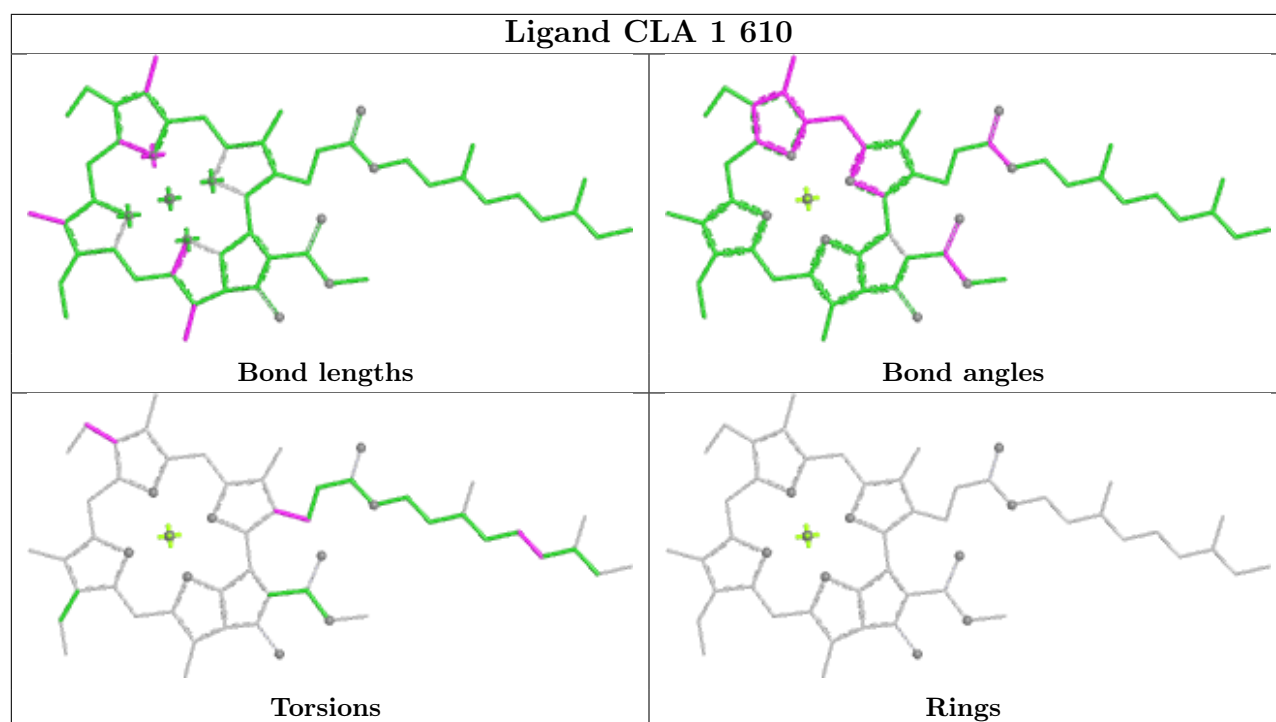


Rings

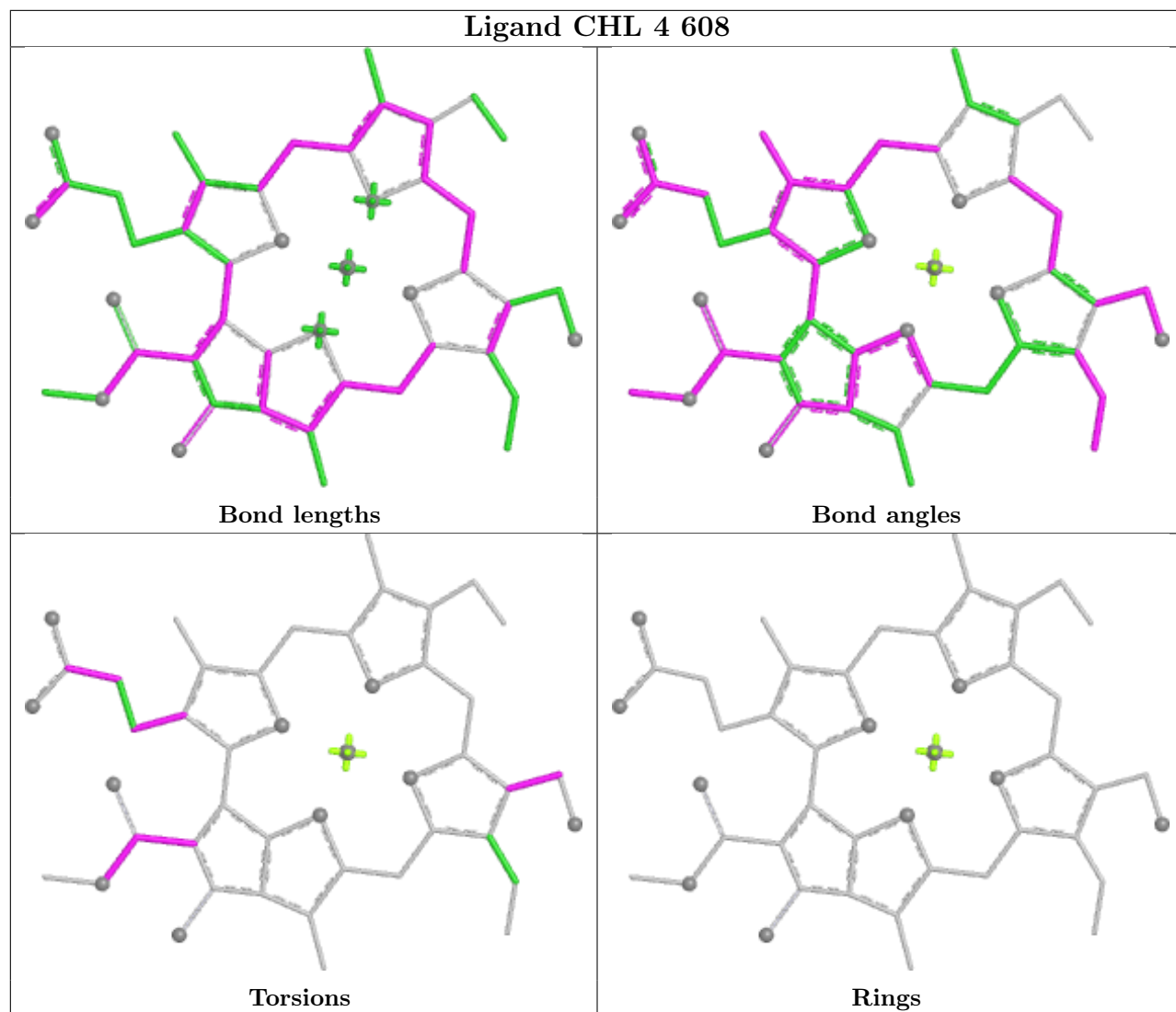


Ligand CLA 3 604

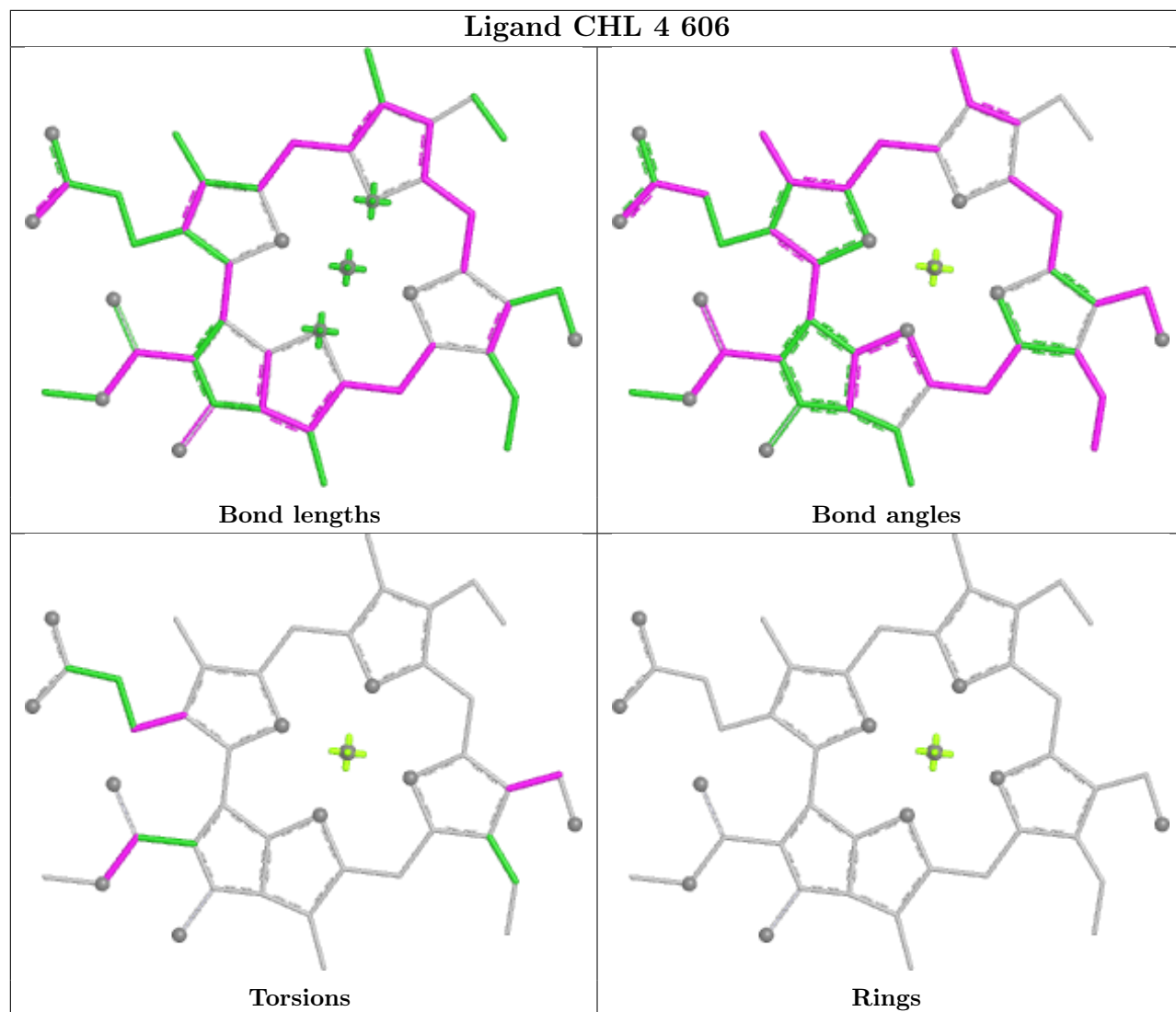




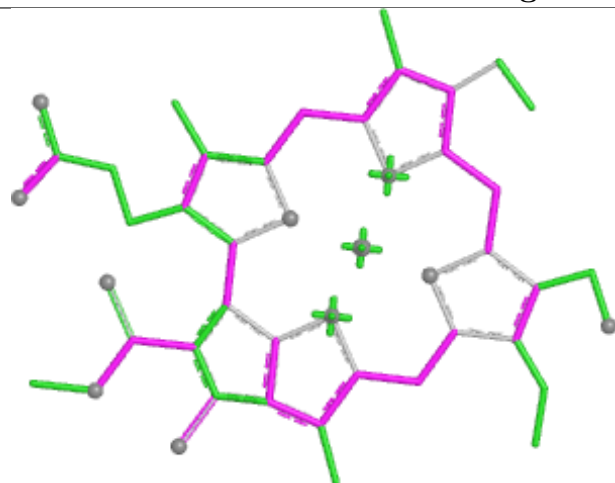
Ligand CHL 4 608



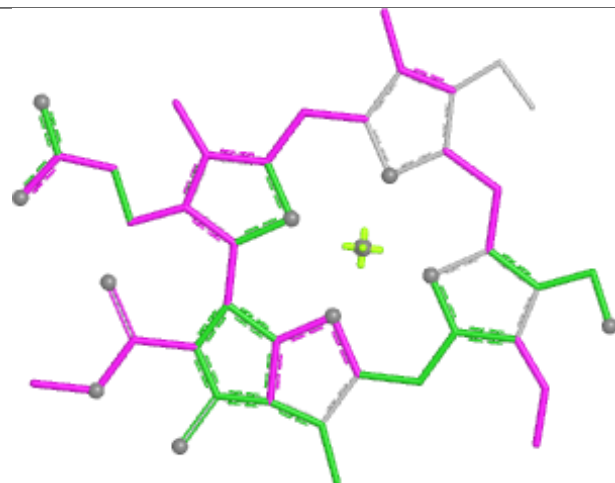
Ligand CHL 4 606



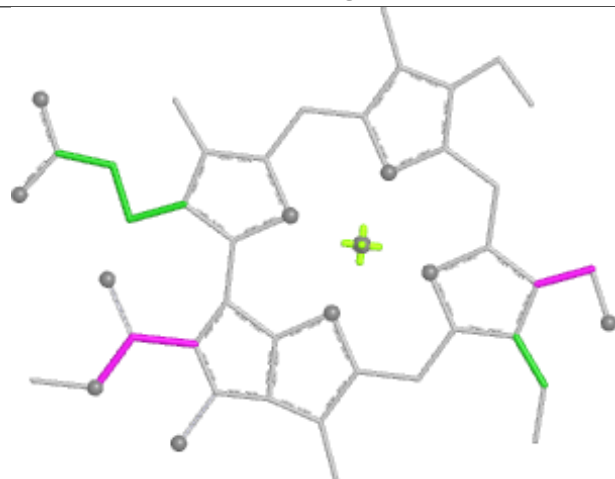
Ligand CHL 2 606



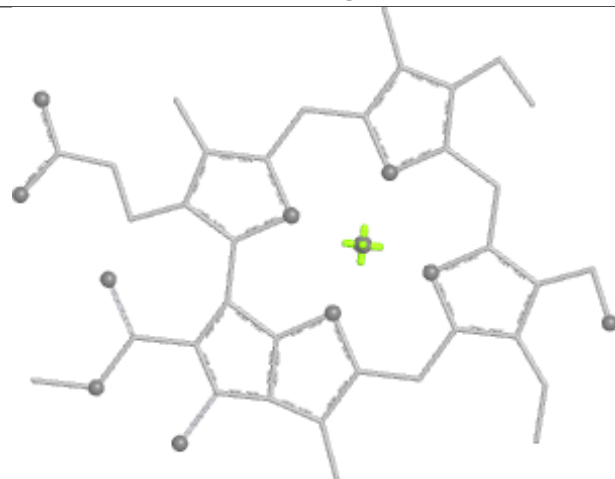
Bond lengths



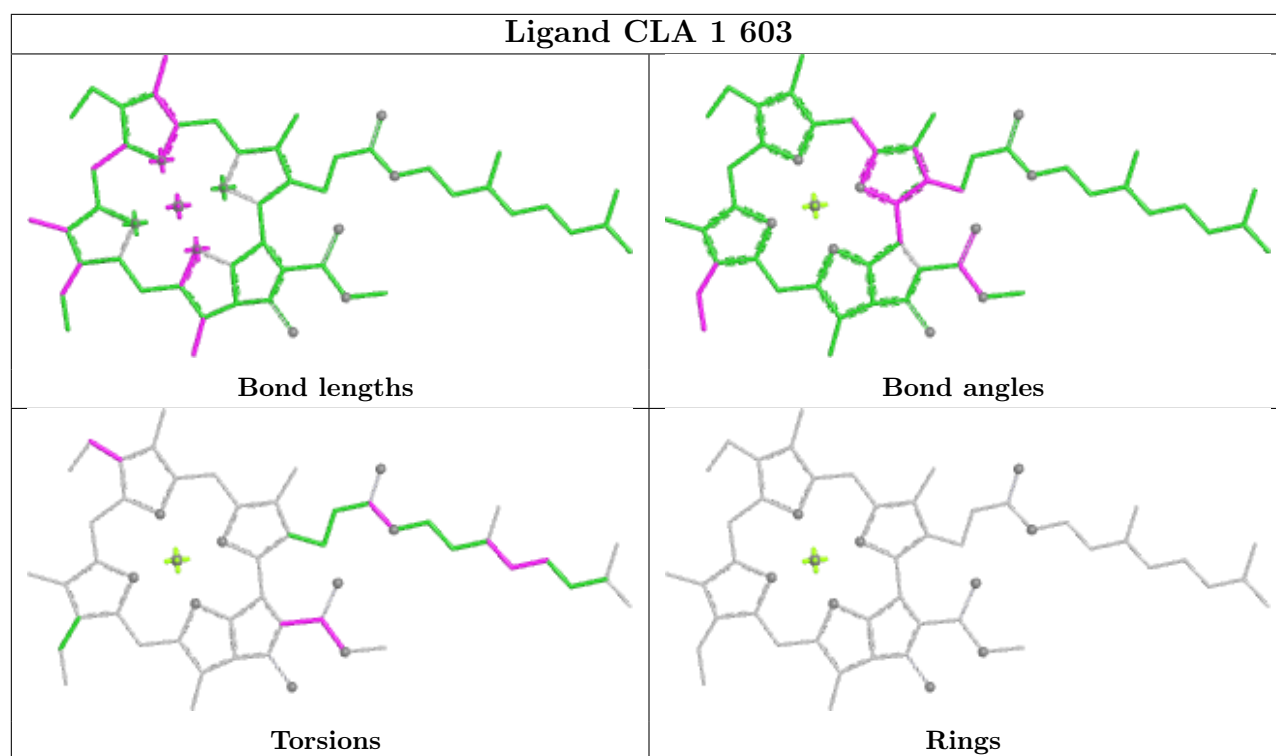
Bond angles



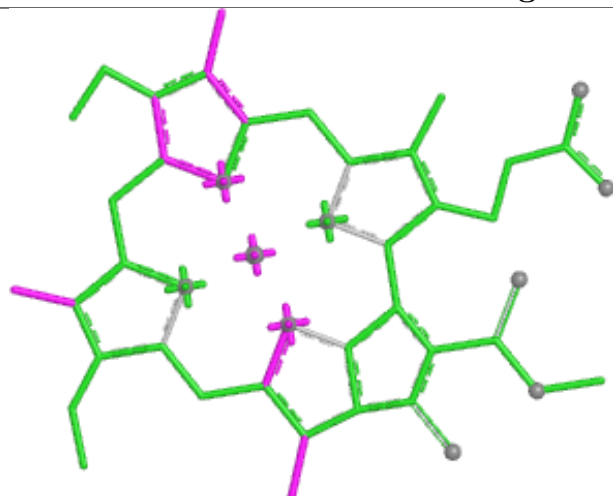
Torsions



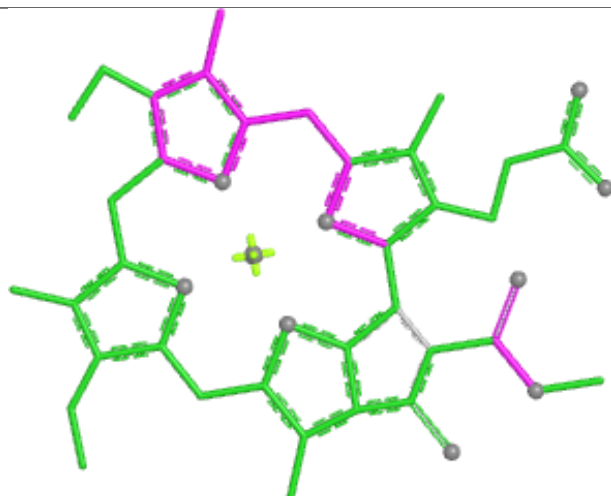
Rings



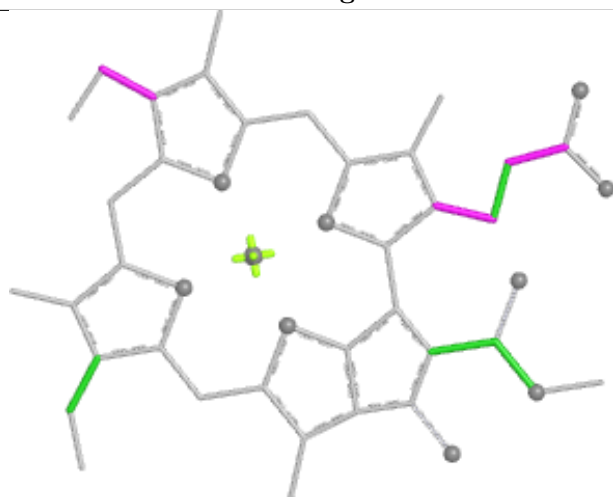
Ligand CLA 4 610



Bond lengths



Bond angles

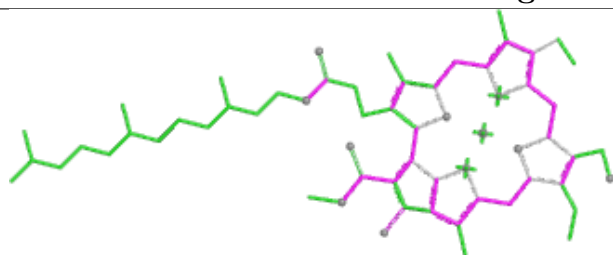


Torsions

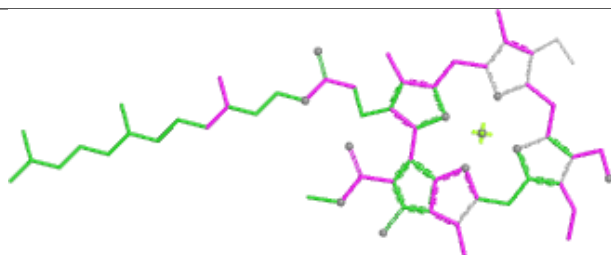


Rings

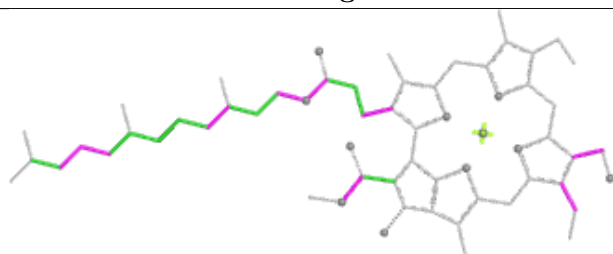
Ligand CHL 3 609



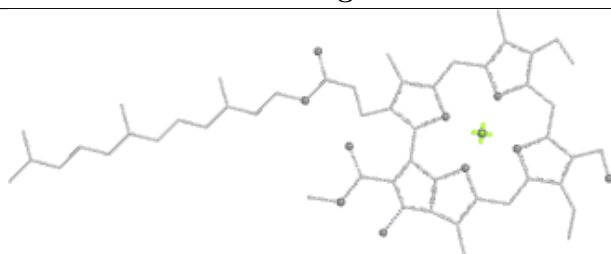
Bond lengths



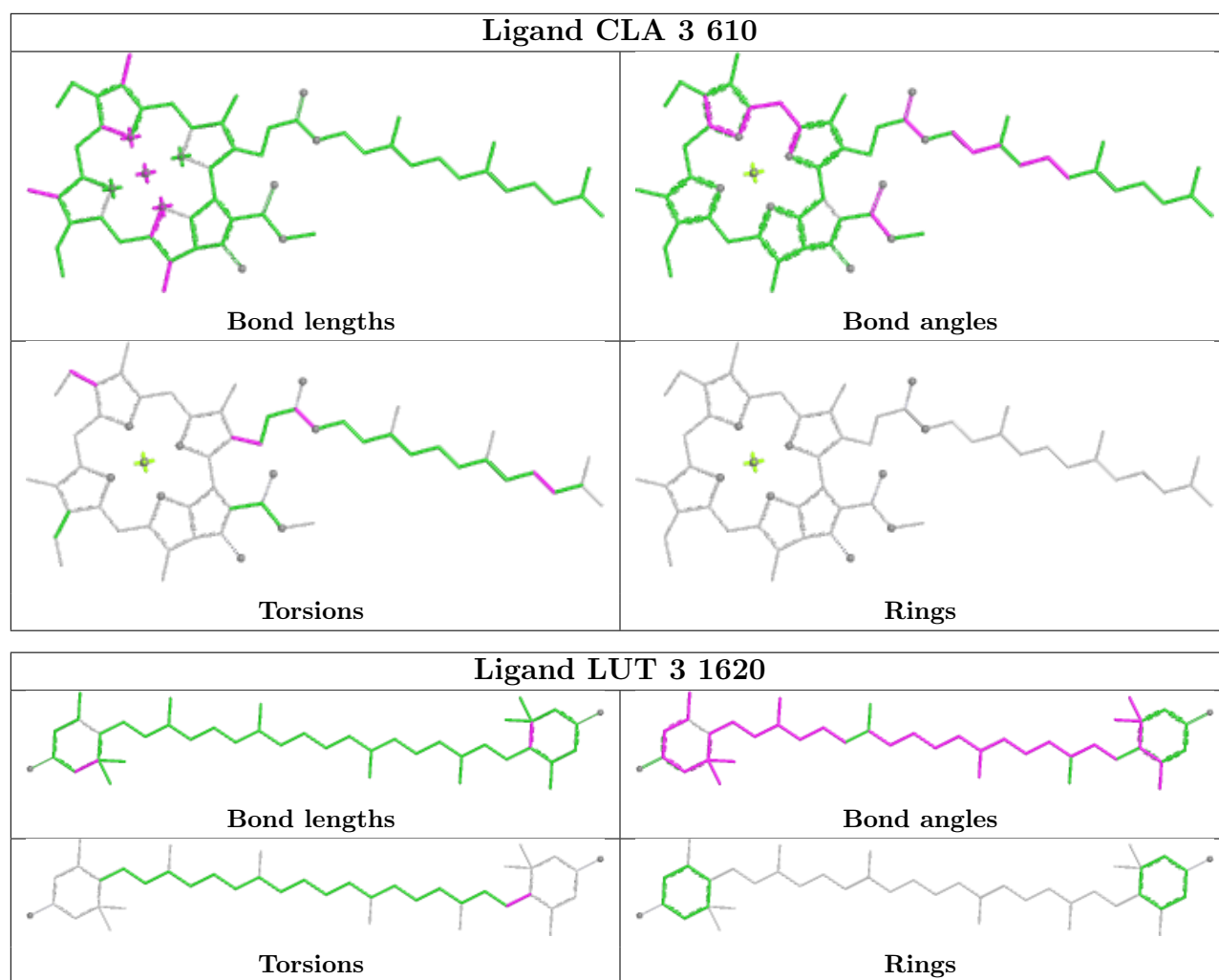
Bond angles



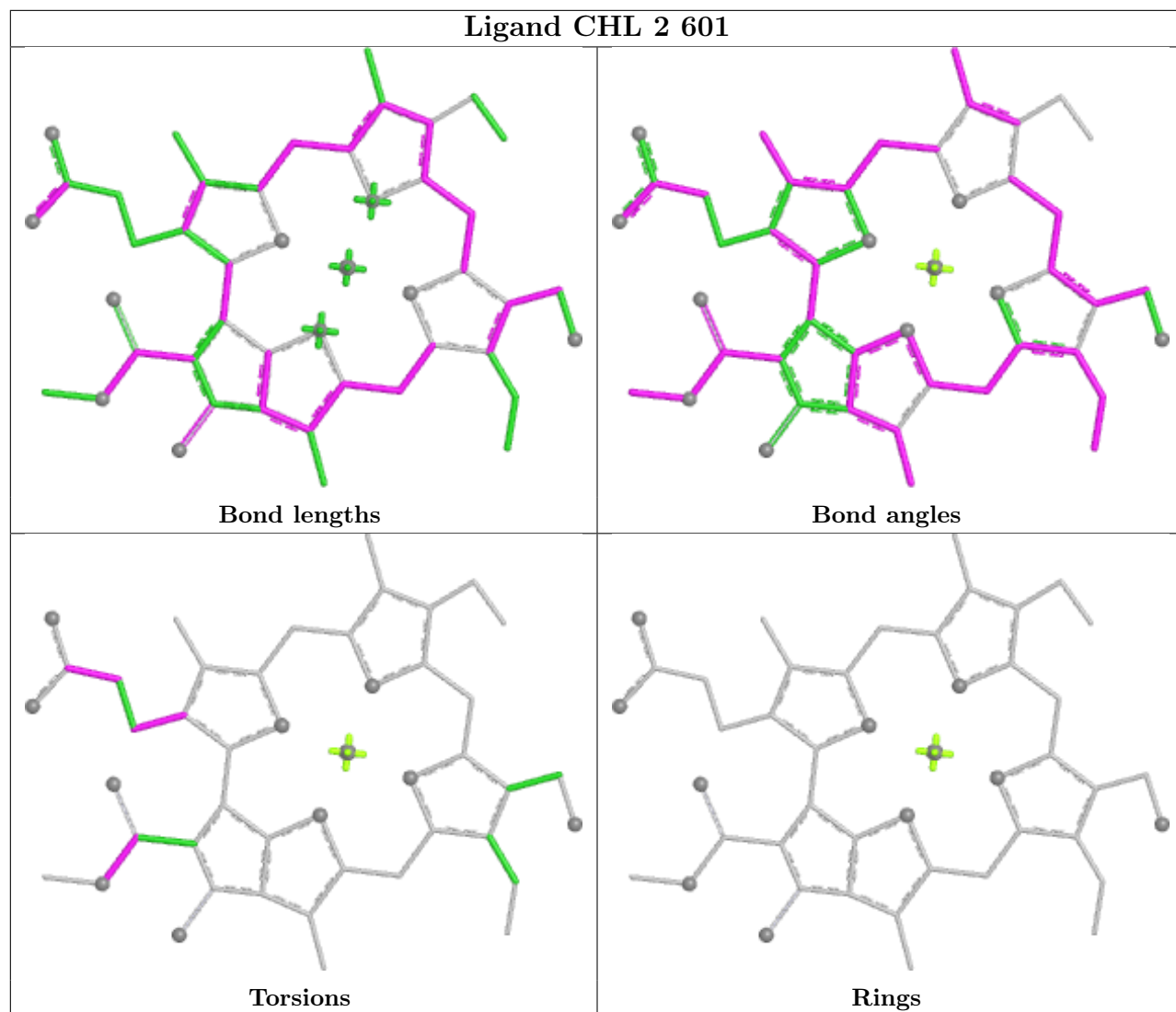
Torsions



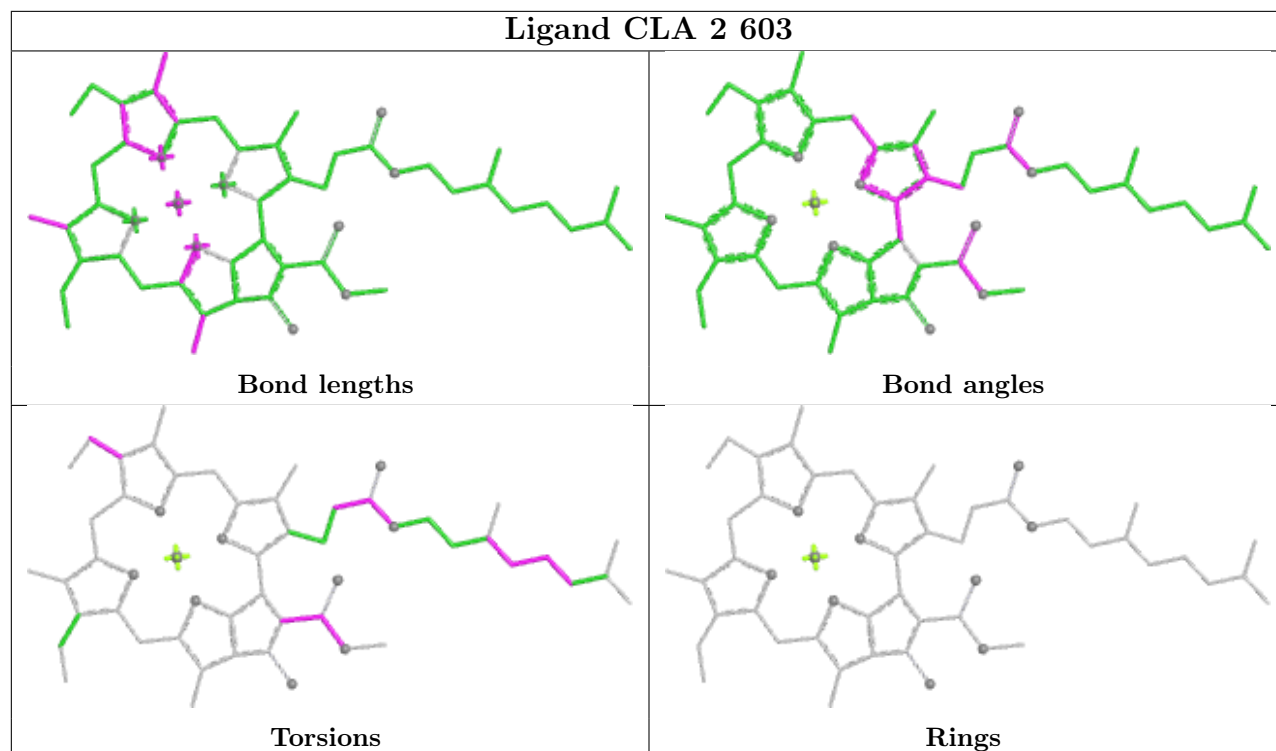
Rings



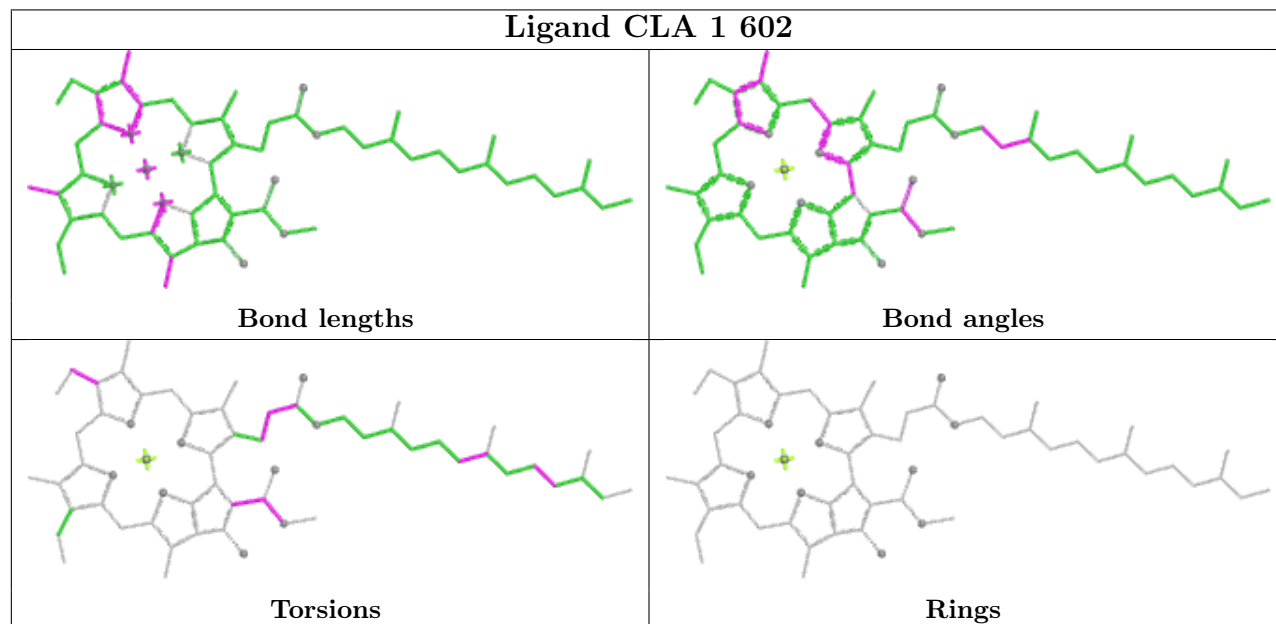
Ligand CHL 2 601



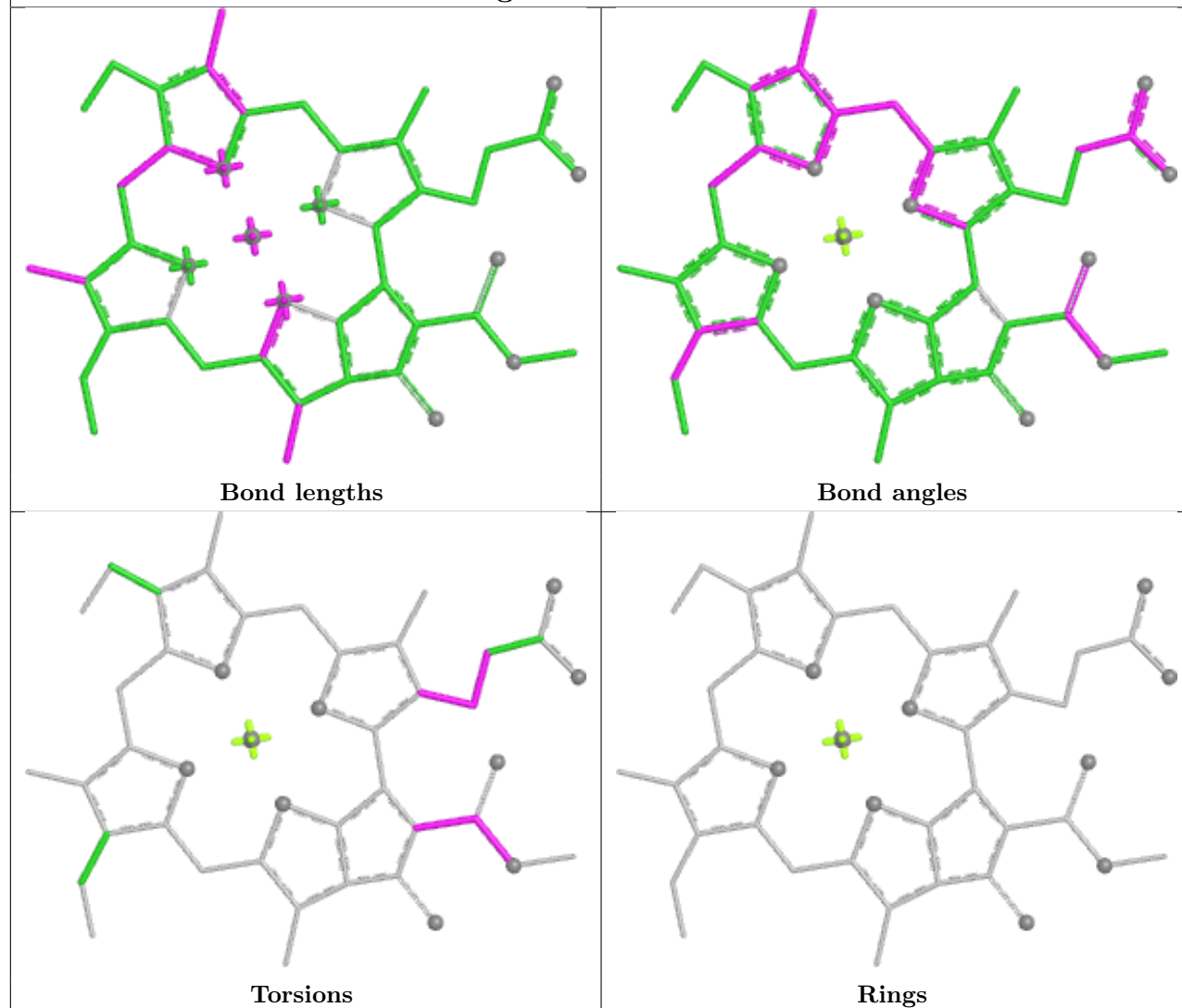
Ligand CLA 2 603



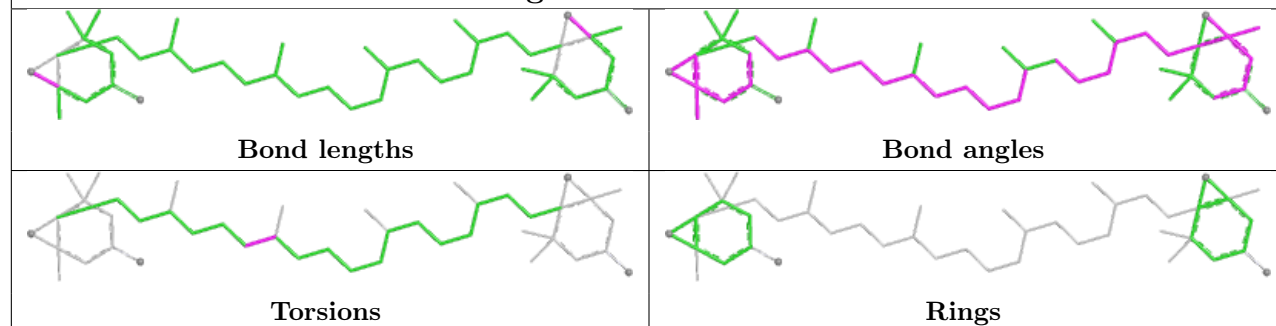
Ligand CLA 1 602

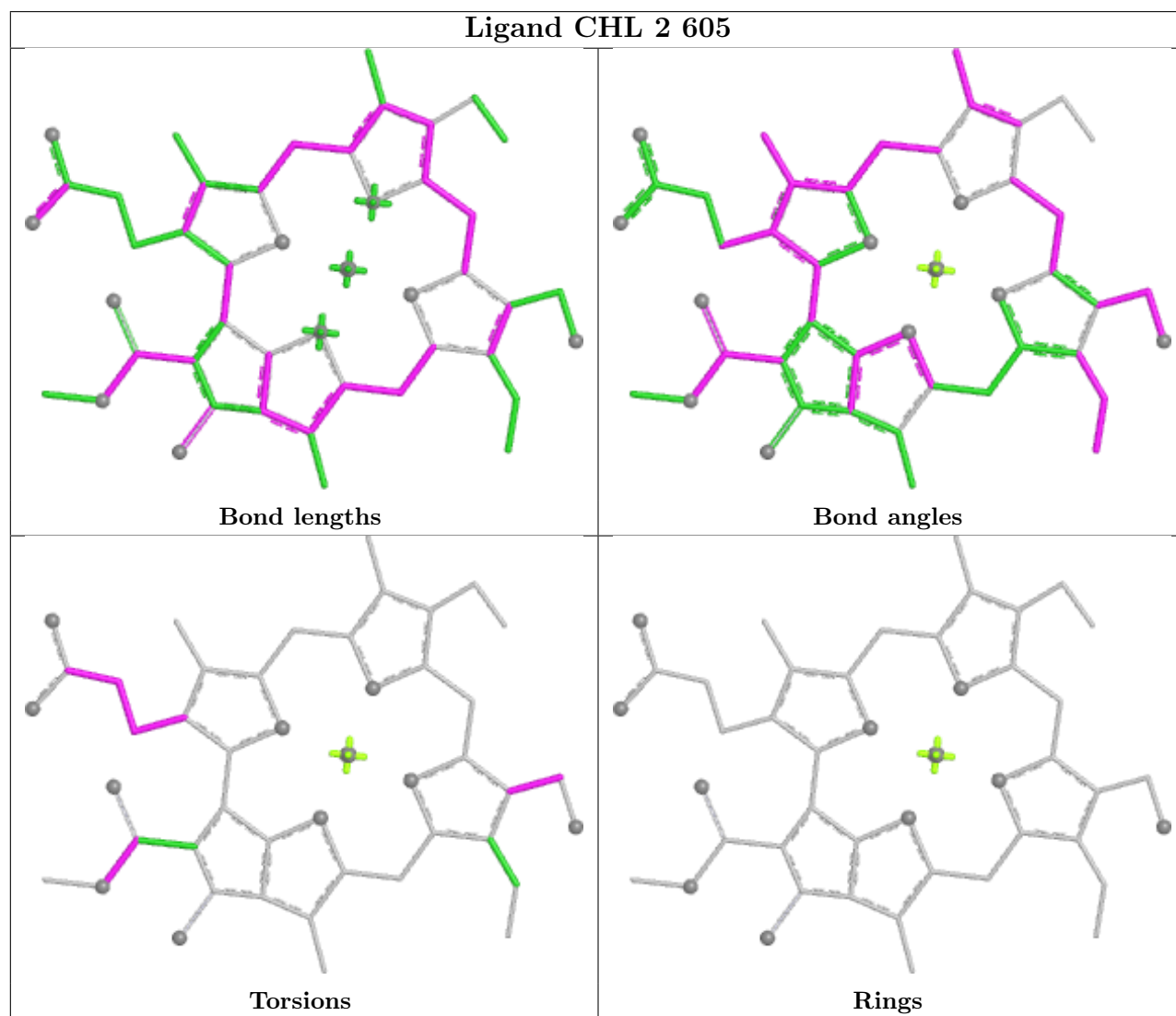
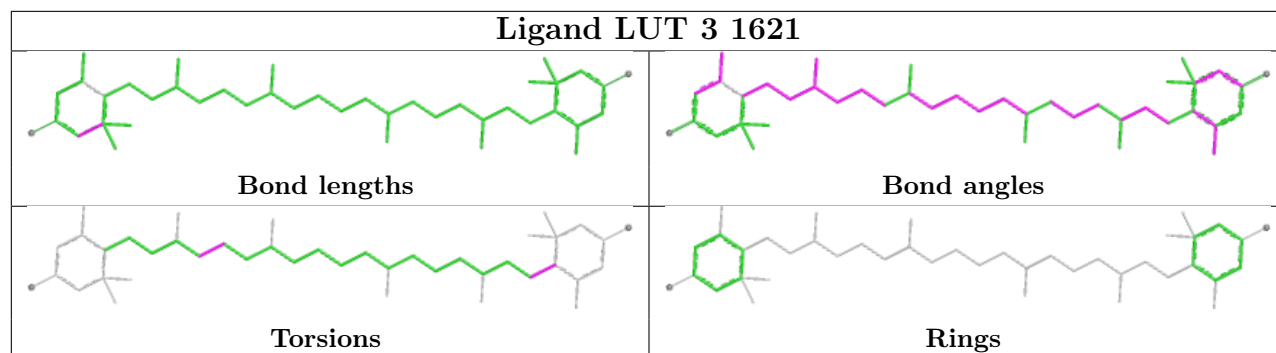


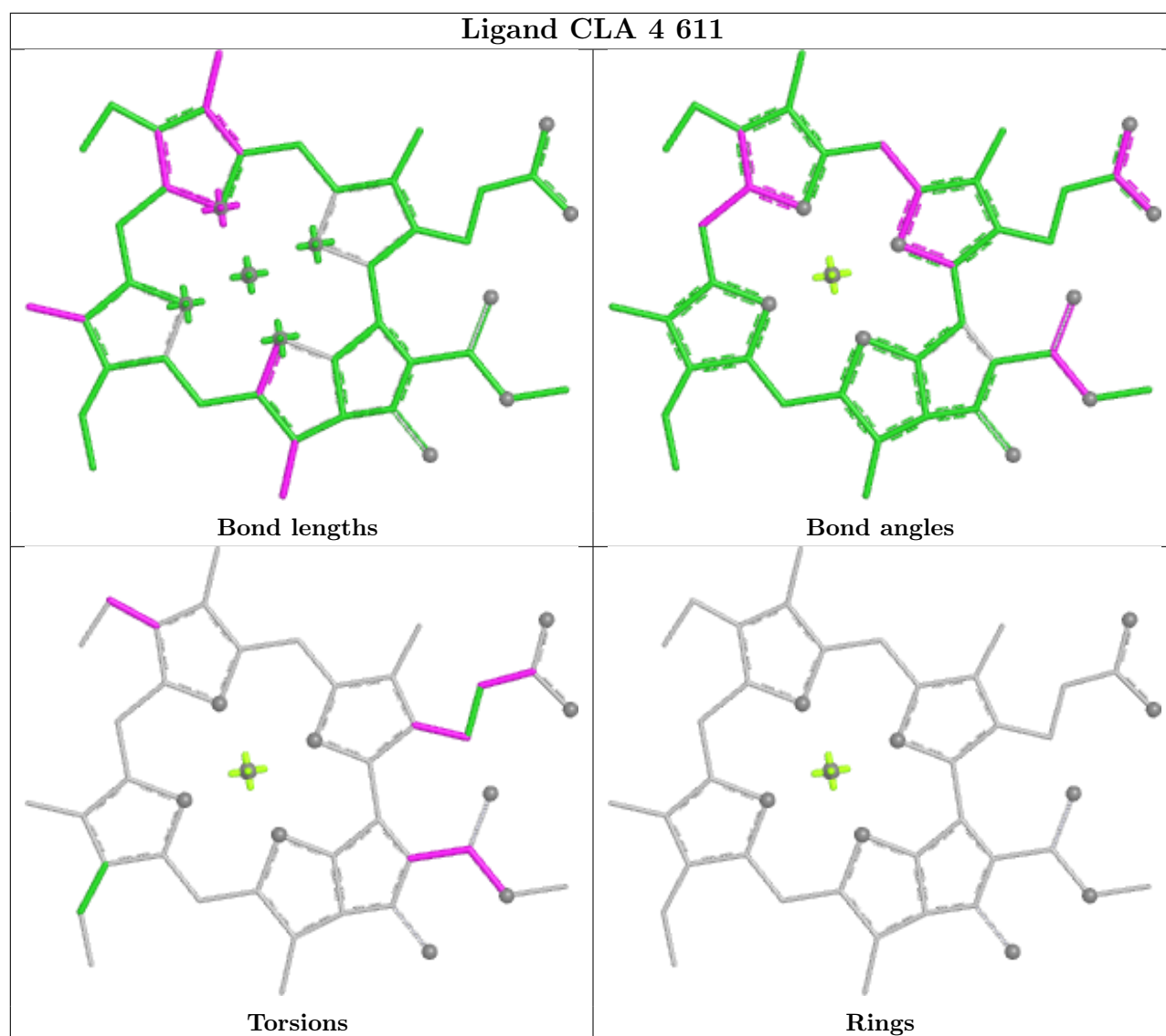
Ligand CLA 2 604



Ligand XAT 1 1622







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

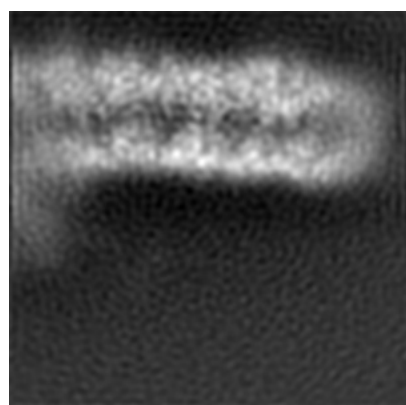
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-6744. These allow visual inspection of the internal detail of the map and identification of artifacts.

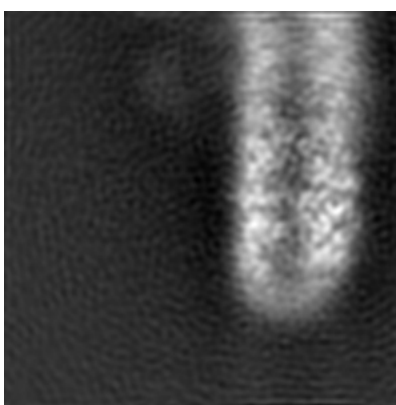
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

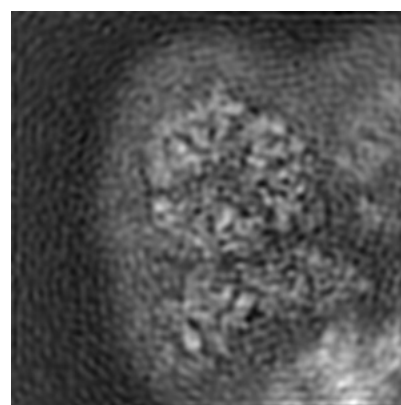
6.1.1 Primary map



X



Y

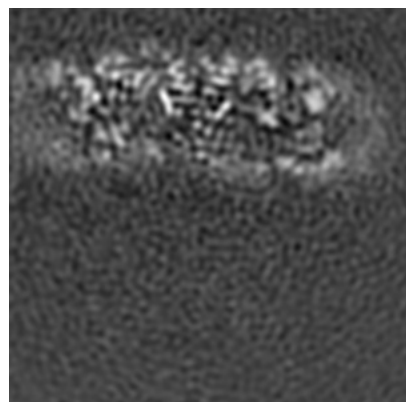


Z

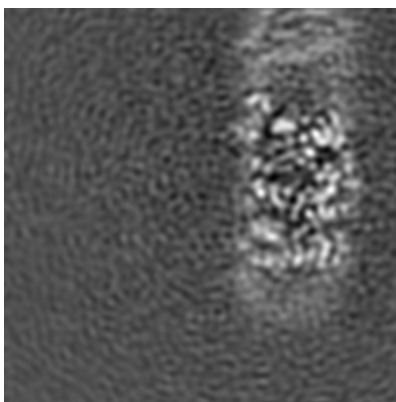
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

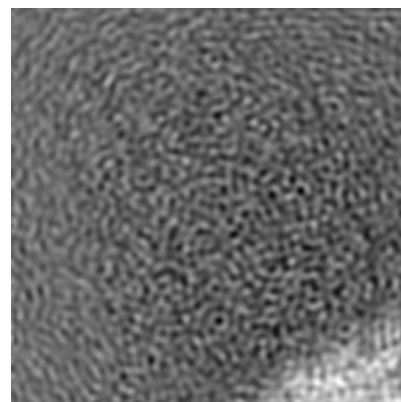
6.2.1 Primary map



X Index: 75



Y Index: 75

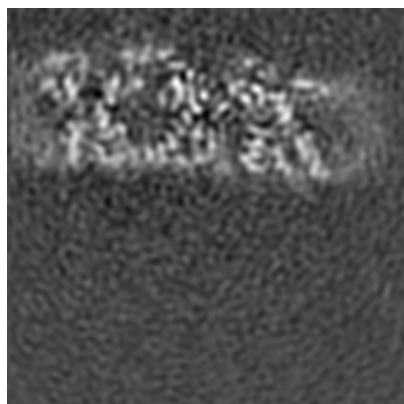


Z Index: 75

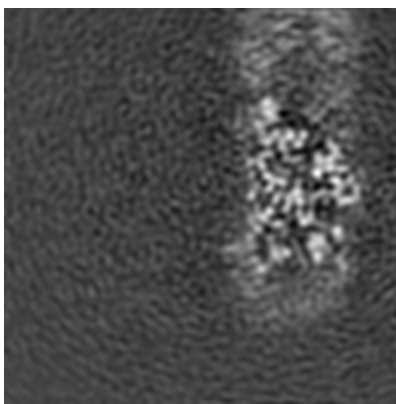
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

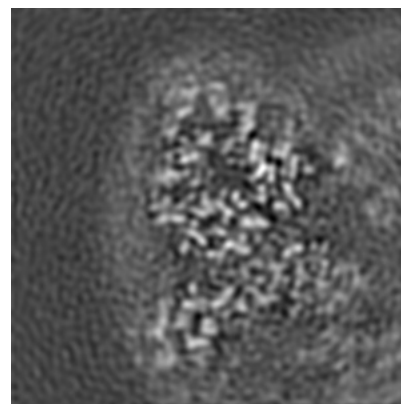
6.3.1 Primary map



X Index: 69



Y Index: 70

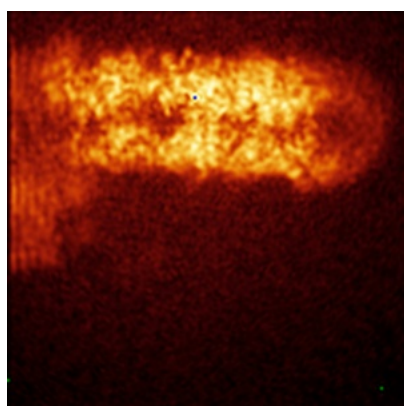


Z Index: 118

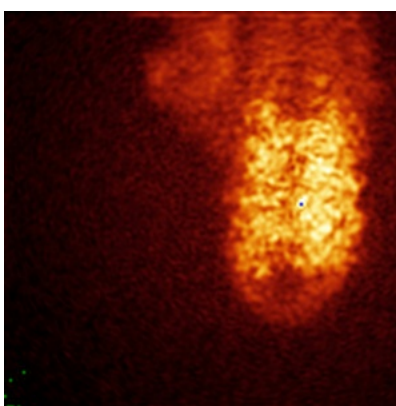
The images above show the largest variance slices of the map in three orthogonal directions.

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

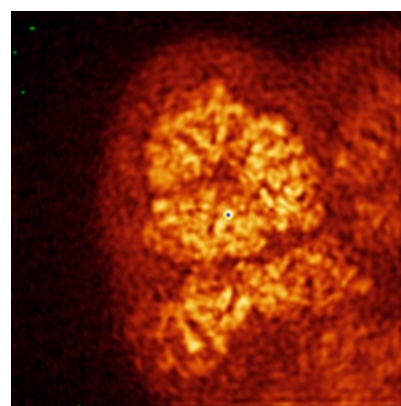
6.4.1 Primary map



X



Y

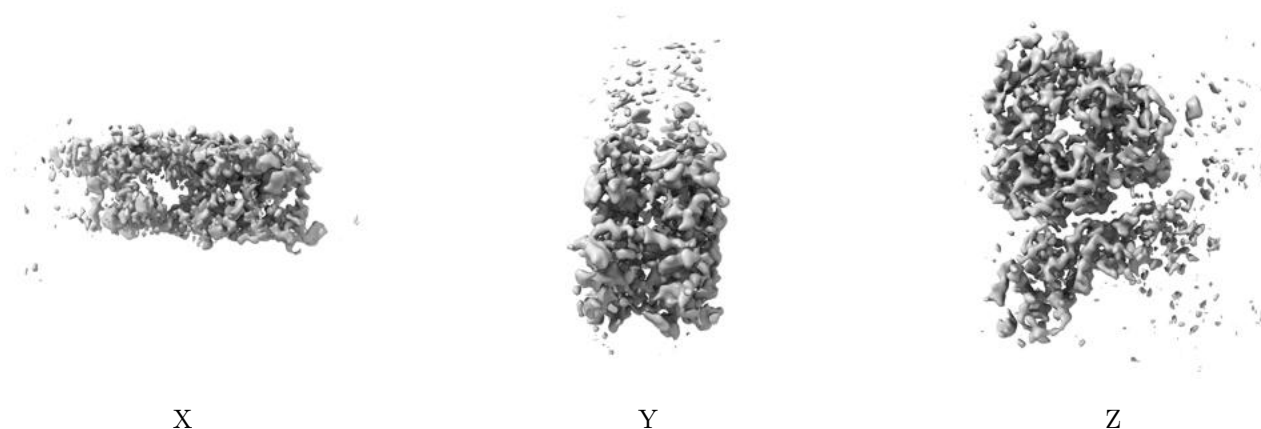


Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.028. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

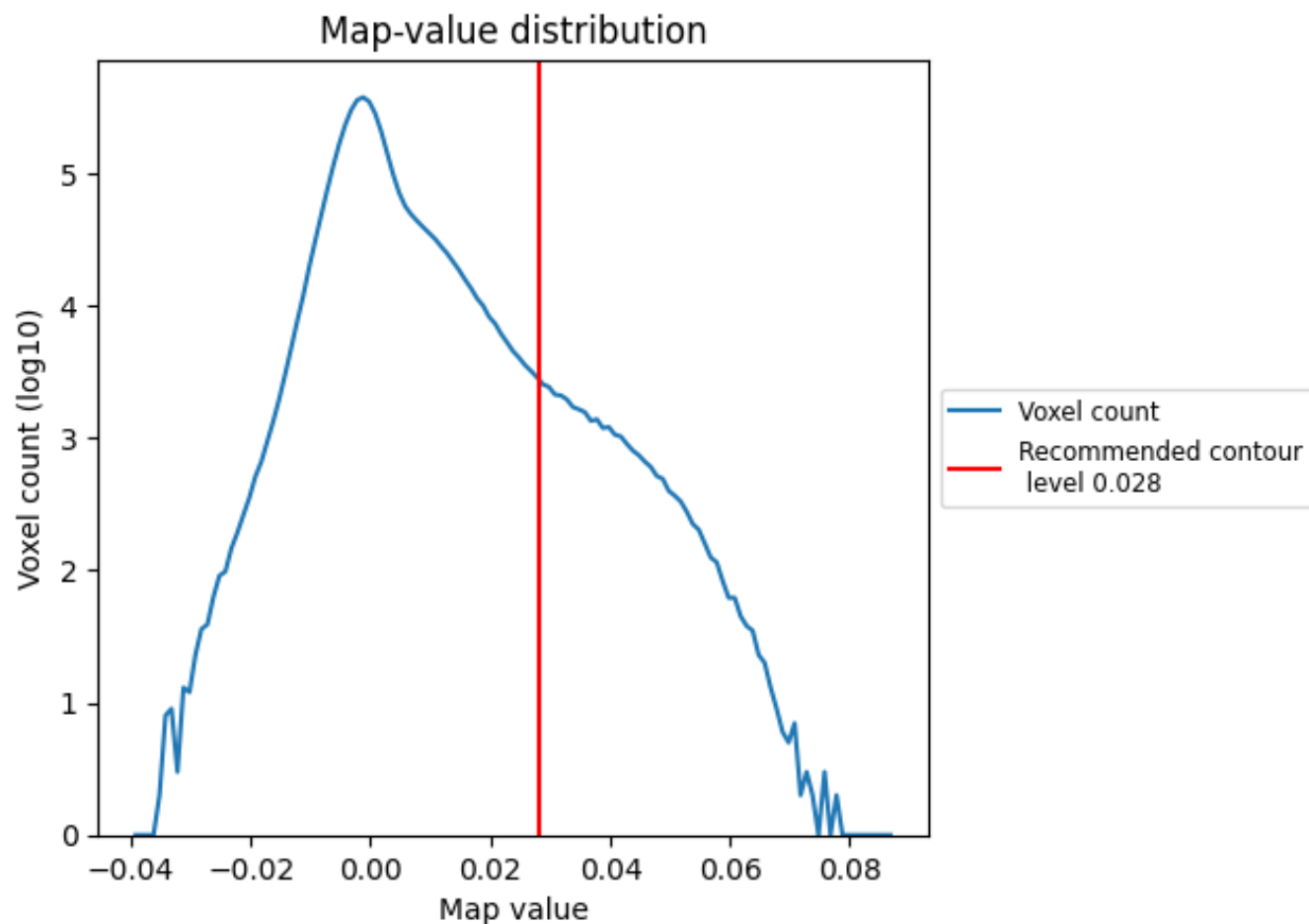
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

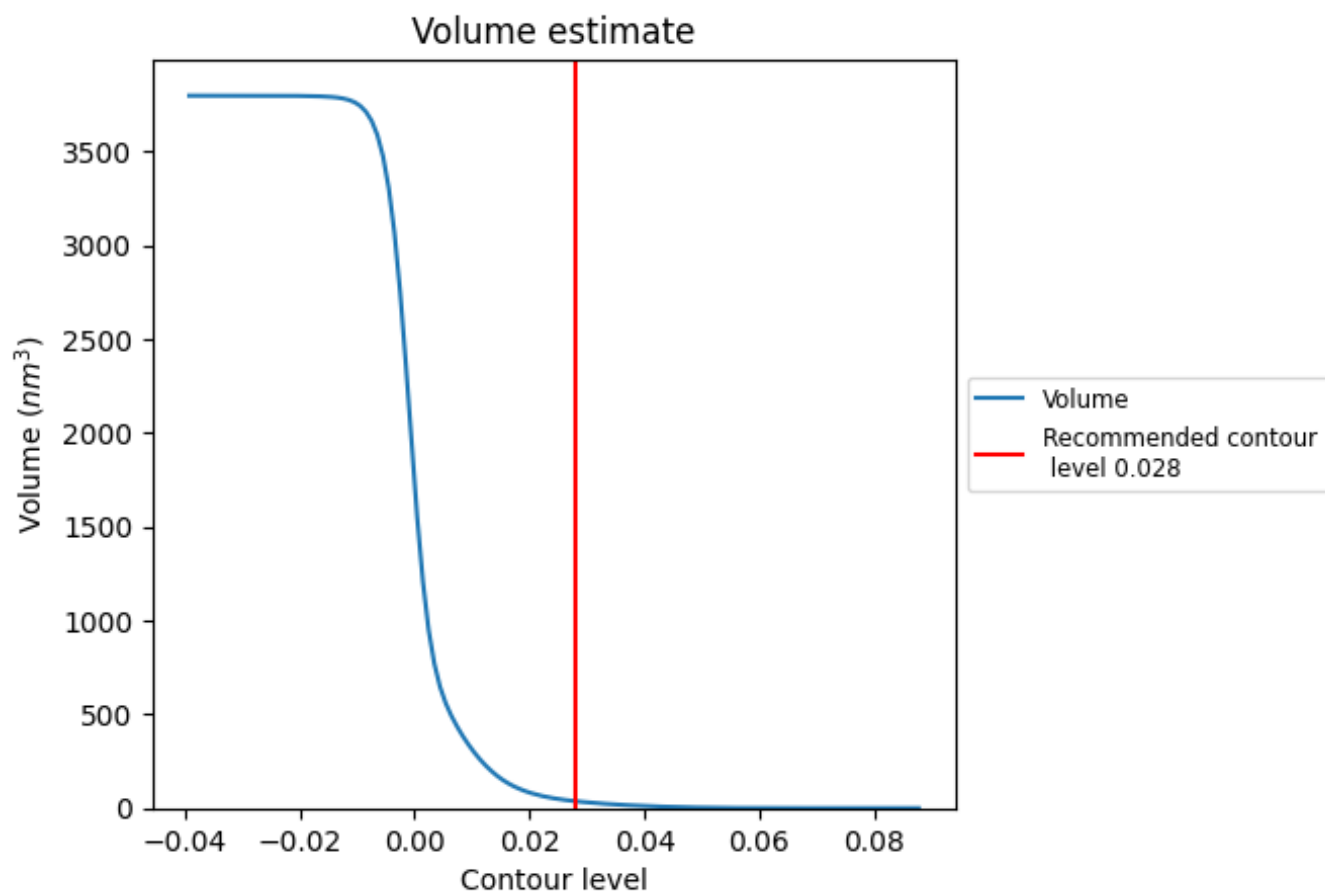
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

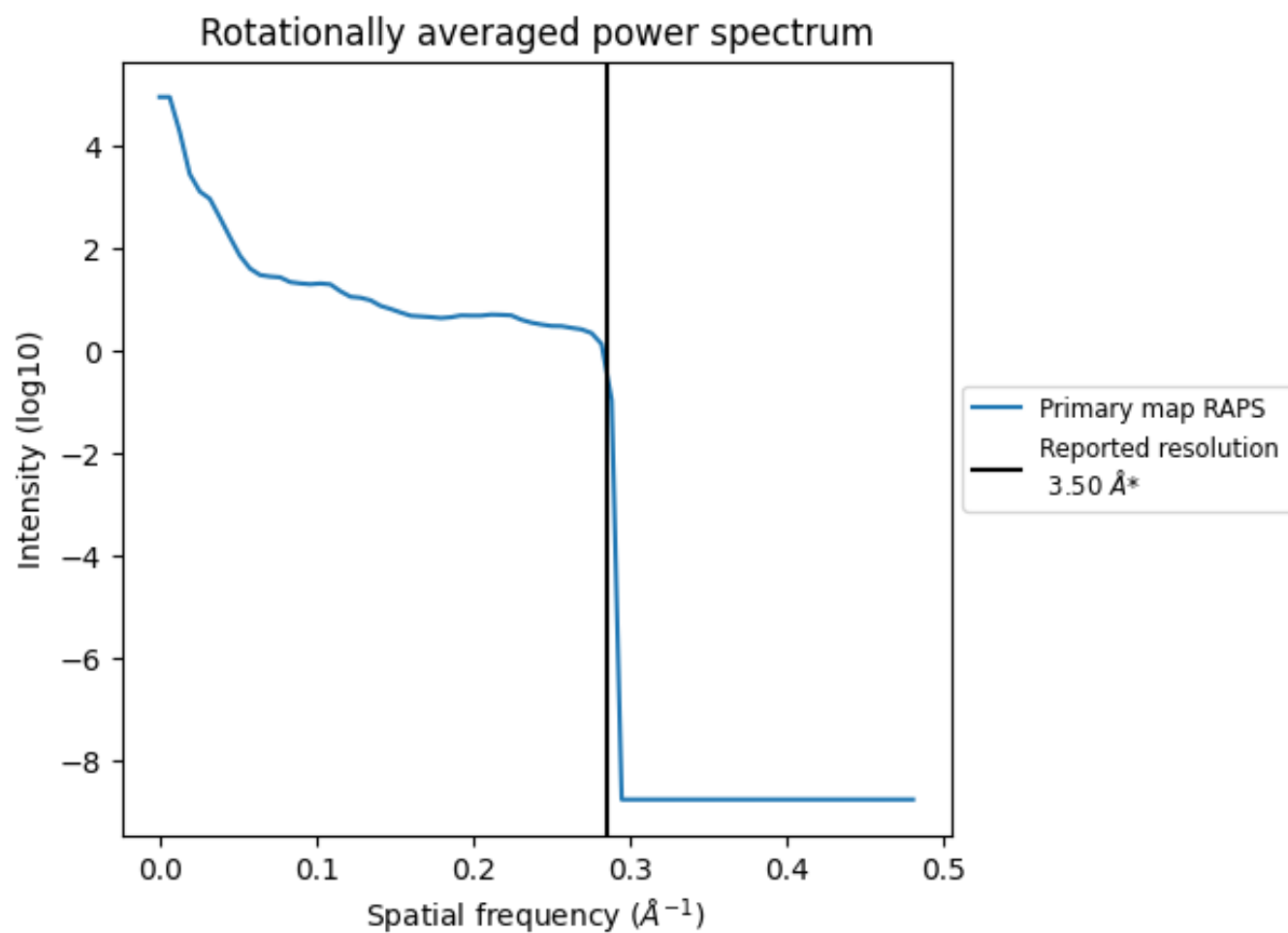
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 37 nm^3 ; this corresponds to an approximate mass of 33 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ



*Reported resolution corresponds to spatial frequency of 0.286 Å⁻¹

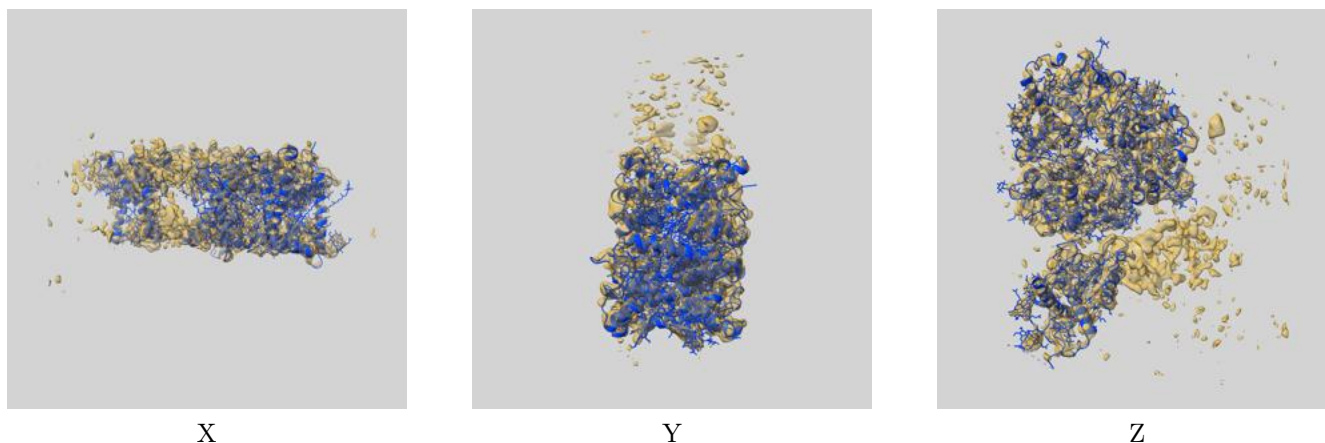
8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit [i](#)

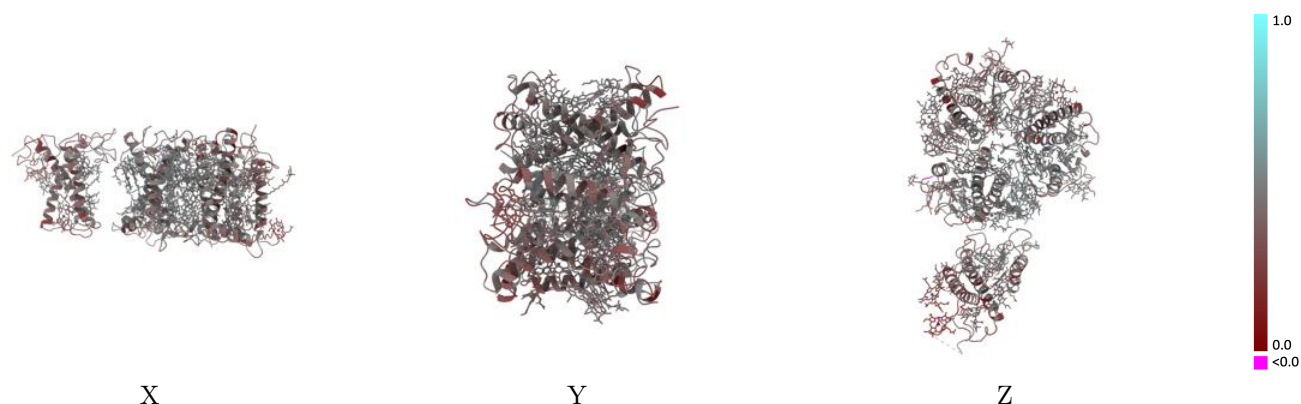
This section contains information regarding the fit between EMDB map EMD-6744 and PDB model 5XNO. Per-residue inclusion information can be found in section [3](#) on page [13](#).

9.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 0.028 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



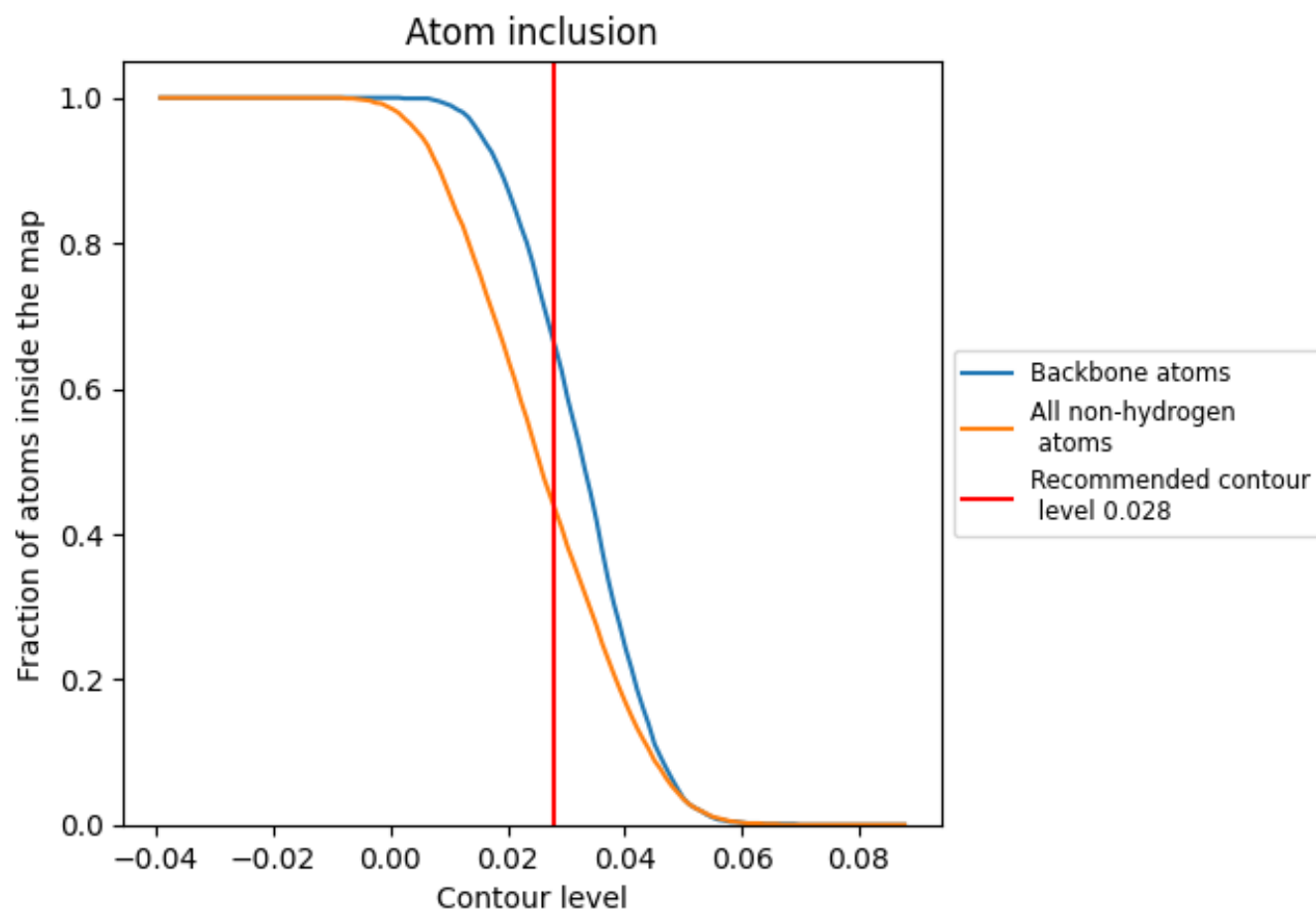
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.028).

9.4 Atom inclusion [i](#)



At the recommended contour level, 66% of all backbone atoms, 44% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.028) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.4360	0.4100
1	0.4500	0.4300
2	0.4070	0.3980
3	0.4720	0.4390
4	0.4080	0.3640

