



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 12:31 AM UTC

PDB ID : 8ABK / pdb_00008abk
EMDB ID : EMD-15324
Title : Complex III₂ from *Yarrowia lipolytica*, decylubiquinol bound, b-position
Authors : Wieferig, J.P.; Kuhlbrandt, W.
Deposited on : 2022-07-04
Resolution : 2.50 Å(reported)

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

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

A user guide is available at

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

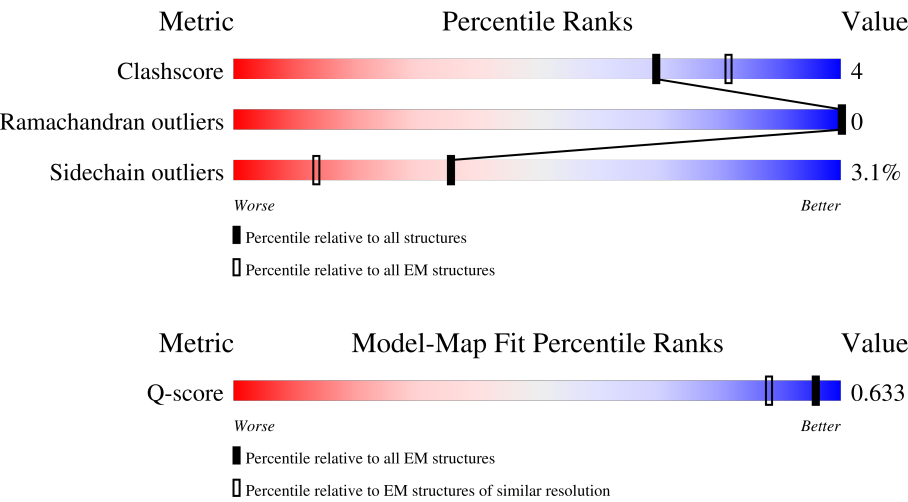
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 i

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 2.50 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.



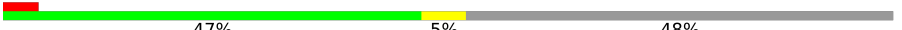
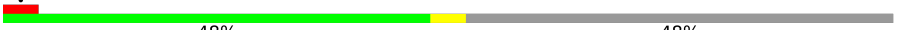
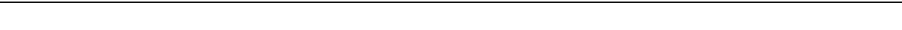
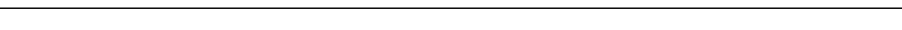
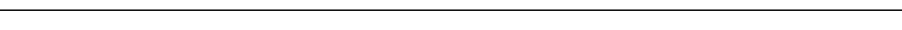
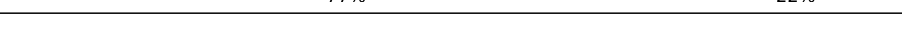
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	7115 (2.00 - 3.00)

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	C	385	92% 7%
1	N	385	93% 6%
2	E	225	22% 74%
2	P	225	18% 60% 20% 17%

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Mol	Chain	Length	Quality of chain
3	G	128	 88% 9% . .
3	R	128	 89% 7% . .
4	F	137	 47% 5% 48%
4	Q	137	 48% . 48%
5	A	474	 85% 7% . 8%
5	L	474	 86% 5% 8%
6	B	417	 88% 8% .
6	M	417	 87% 9% .
7	D	330	 67% 7% 26%
7	O	330	 68% 6% 26%
8	H	93	 85% 5% . 9%
8	S	93	 84% 6% . 9%
9	I	69	 77% . 22%
9	T	69	 77% . 22%
10	J	82	 83% 6% . 9%
10	U	82	 82% 10% 9%

2 Entry composition

There are 19 unique types of molecules in this entry. The entry contains 32573 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 Cytochrome b.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	C	383	Total	C	N	O	S	0	0
			3052	2064	474	496	18		
1	N	383	Total	C	N	O	S	0	0
			3052	2064	474	496	18		

- Molecule 2 is a protein called Cytochrome b-c1 complex subunit Rieske, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	P	186	Total	C	N	O	S	0	0
			1445	920	248	268	9		
2	E	59	Total	C	N	O	S	0	0
			452	288	74	87	3		

- Molecule 3 is a protein called Cytochrome b-c1 complex subunit 7.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	G	124	Total	C	N	O	S	0	0
			994	640	162	190	2		
3	R	124	Total	C	N	O	S	0	0
			994	640	162	190	2		

- Molecule 4 is a protein called YALI0F24673p.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	F	71	Total	C	N	O	S	0	0
			579	361	99	115	4		
4	Q	71	Total	C	N	O	S	0	0
			579	361	99	115	4		

- Molecule 5 is a protein called YALI0A14806p.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	A	438	Total	C	N	O	S	0	0
			3446	2154	603	682	7		
5	L	438	Total	C	N	O	S	0	0
			3446	2154	603	682	7		

- Molecule 6 is a protein called Cytochrome b-c1 complex subunit 2, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	B	402	Total	C	N	O	S	0	0
			3008	1907	516	583	2		
6	M	402	Total	C	N	O	S	0	0
			3008	1907	516	583	2		

- Molecule 7 is a protein called YALI0A17468p.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	D	244	Total	C	N	O	S	0	0
			1893	1210	323	352	8		
7	O	244	Total	C	N	O	S	0	0
			1893	1210	323	352	8		

- Molecule 8 is a protein called Cytochrome b-c1 complex subunit 8.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	H	85	Total	C	N	O	S	0	0
			690	459	118	111	2		
8	S	85	Total	C	N	O	S	0	0
			690	459	118	111	2		

- Molecule 9 is a protein called Complex III subunit 9.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	I	54	Total	C	N	O	S	0	0
			452	297	76	78	1		
9	T	54	Total	C	N	O	S	0	0
			452	297	76	78	1		

- Molecule 10 is a protein called YALI0C12210p.

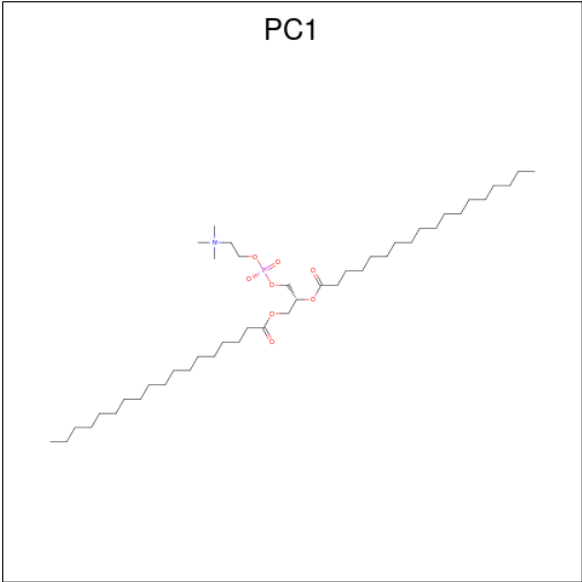
Mol	Chain	Residues	Atoms				AltConf	Trace
10	J	75	Total	C	N	O	0	0
			598	403	99	96		

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Mol	Chain	Residues	Atoms				AltConf	Trace
10	U	75	Total	C	N	O	0	0
			598	403	99	96		

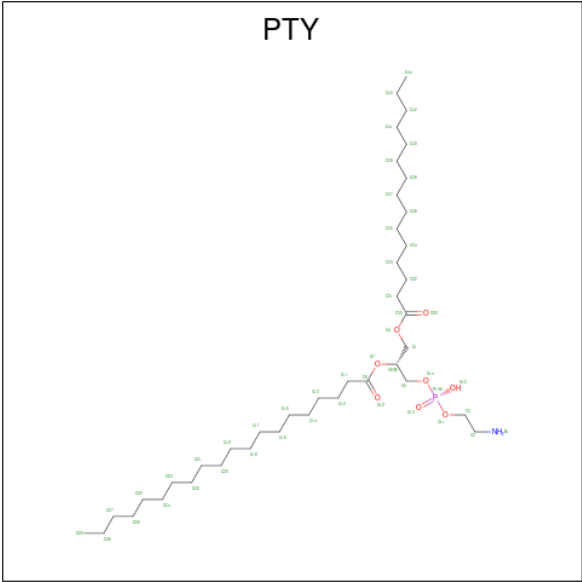
- # HEM

- Molecule 12 is 1,2-DIACYL-SN-GLYCERO-3-PHOSPHOCHOLINE (CCD ID: PC1) (formula: $C_{44}H_{88}NO_8P$).



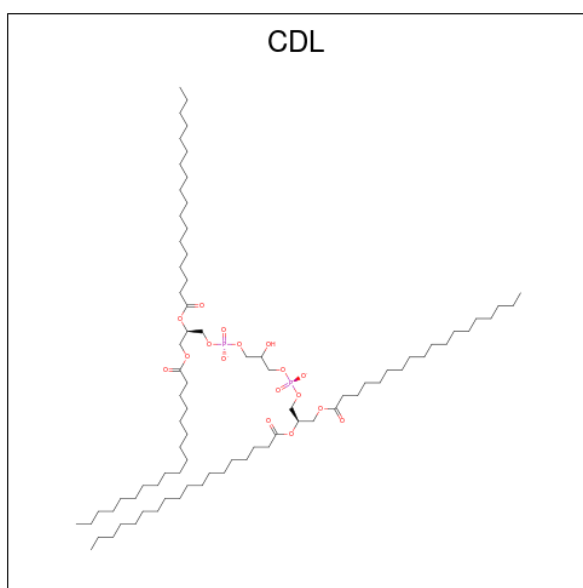
Mol	Chain	Residues	Atoms					AltConf
12	C	1	Total	C	N	O	P	0
			38	28	1	8	1	
12	I	1	Total	C	N	O	P	0
			32	22	1	8	1	
12	N	1	Total	C	N	O	P	0
			38	28	1	8	1	
12	T	1	Total	C	N	O	P	0
			32	22	1	8	1	

- Molecule 13 is PHOSPHATIDYLETHANOLAMINE (CCD ID: PTY) (formula: C₄₀H₈₀NO₈P).



Mol	Chain	Residues	Atoms					AltConf
13	C	1	Total	C	N	O	P	0
			41	31	1	8	1	
13	P	1	Total	C	N	O	P	0
			41	31	1	8	1	
13	N	1	Total	C	N	O	P	0
			41	31	1	8	1	
13	E	1	Total	C	N	O	P	0
			41	31	1	8	1	

- Molecule 14 is CARDIOLIPIN (CCD ID: CDL) (formula: $C_{81}H_{156}O_{17}P_2$).



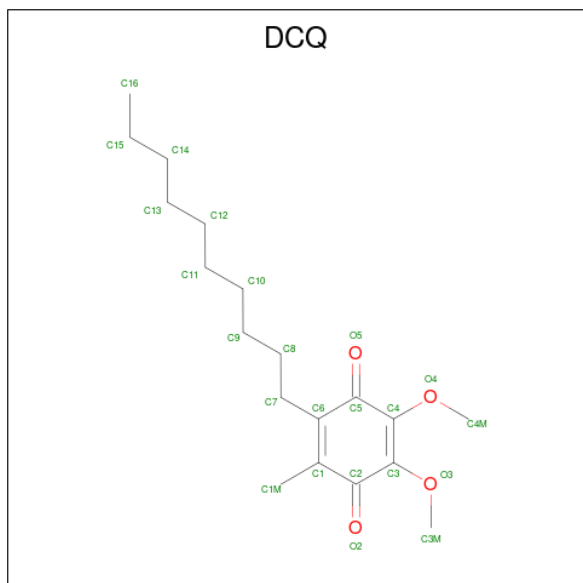
Mol	Chain	Residues	Atoms				AltConf
14	C	1	Total	C	O	P	0
			48	29	17	2	
14	A	1	Total	C	O	P	0
			42	25	15	2	
14	A	1	Total	C	O	P	0
			47	30	15	2	
14	H	1	Total	C	O	P	0
			50	31	17	2	
14	H	1	Total	C	O	P	0
			39	20	17	2	
14	N	1	Total	C	O	P	0
			50	31	17	2	
14	N	1	Total	C	O	P	0
			48	29	17	2	
14	L	1	Total	C	O	P	0
			42	25	15	2	

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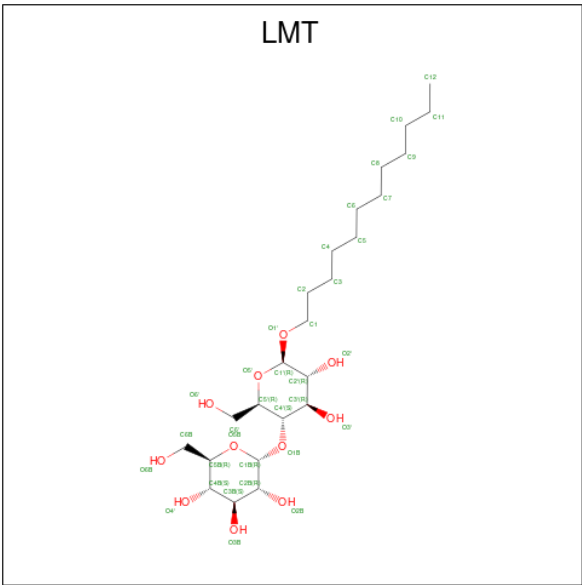
Mol	Chain	Residues	Atoms				AltConf
14	L	1	Total	C	O	P	0
			47	30	15	2	
14	S	1	Total	C	O	P	0
			39	20	17	2	

- Molecule 15 is 2-decyl-5,6-dimethoxy-3-methylcyclohexa-2,5-diene-1,4-dione (CCD ID: DCQ) (formula: $C_{19}H_{30}O_4$) (labeled as "Ligand of Interest" by depositor).



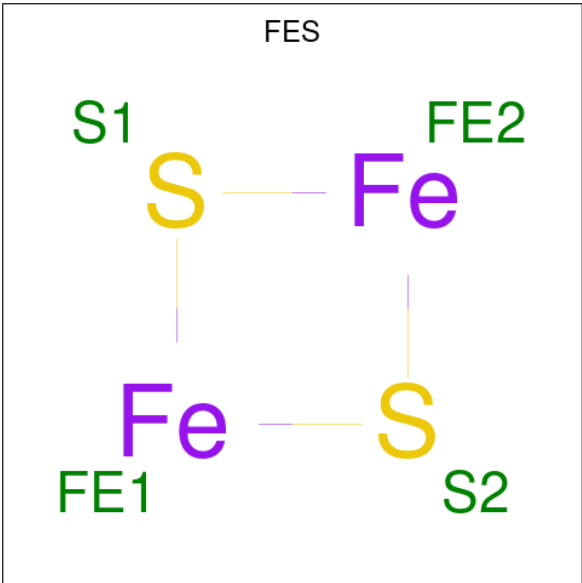
Mol	Chain	Residues	Atoms				AltConf
15	C	1	Total	C	O		0
			23	19	4		
15	N	1	Total	C	O		0
			23	19	4		

- Molecule 16 is DODECYL-BETA-D-MALTOSE (CCD ID: LMT) (formula: $C_{24}H_{46}O_{11}$).



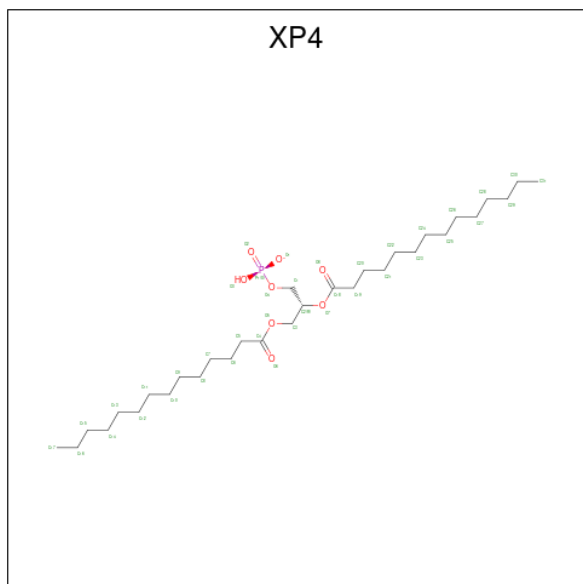
Mol	Chain	Residues	Atoms			AltConf
16	C	1	Total	C	O	0
			35	24	11	
16	J	1	Total	C	O	0
			35	24	11	
16	N	1	Total	C	O	0
			35	24	11	
16	U	1	Total	C	O	0
			35	24	11	

- Molecule 17 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe₂S₂) (labeled as "Ligand of Interest" by depositor).



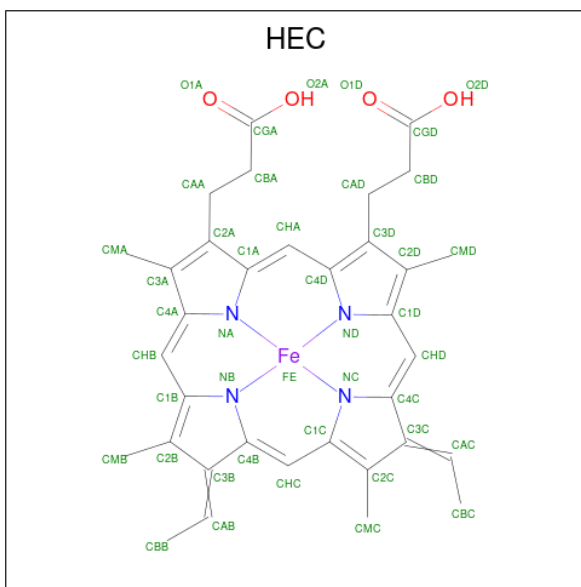
Mol	Chain	Residues	Atoms			AltConf
17	P	1	Total	Fe	S	0
			4	2	2	

- Molecule 18 is 1,2-DIMYRISTOYL-SN-GLYCERO-3-PHOSPHATE (CCD ID: XP4) (formula: $C_{31}H_{60}O_8P$).



Mol	Chain	Residues	Atoms				AltConf
18	A	1	Total	C	O	P	0
			24	15	8	1	
18	L	1	Total	C	O	P	0
			24	15	8	1	

- Molecule 19 is HEME C (CCD ID: HEC) (formula: $C_{34}H_{34}FeN_4O_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
19	D	1	Total 43	C 34	Fe 1	N 4	O 4	0
19	O	1	Total 43	C 34	Fe 1	N 4	O 4	0

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and 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: Cytochrome b

Chain C: 



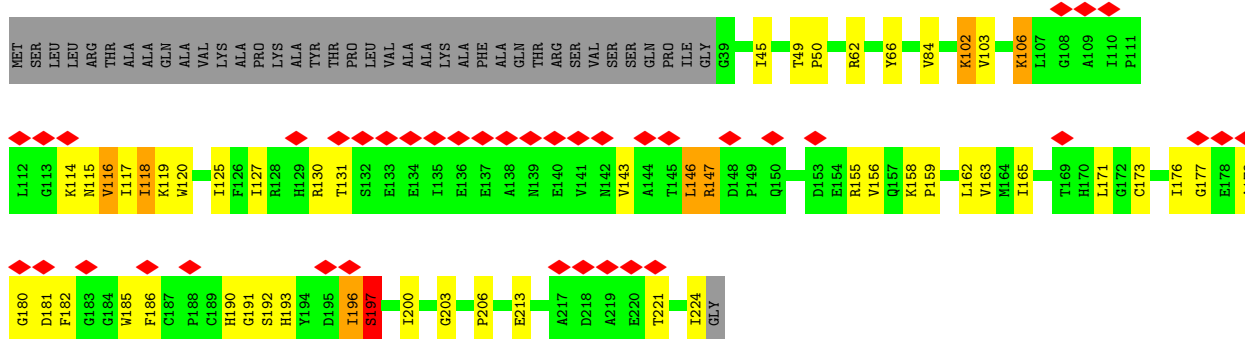
- Molecule 1: Cytochrome b

Chain N: 



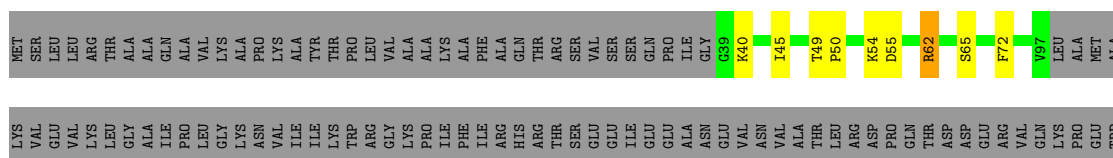
- Molecule 2: Cytochrome b-c1 complex subunit Rieske, mitochondrial

Chain P: 



- Molecule 2: Cytochrome b-c1 complex subunit Rieske, mitochondrial

Chain E: 

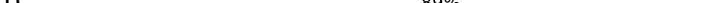


- Molecule 3: Cytochrome b-c1 complex subunit 7

Chain G: 88% 9% ...

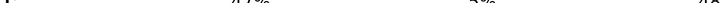


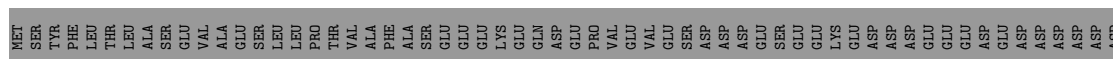
- Molecule 3: Cytochrome b-c1 complex subunit 7

Chain R:  89% 7% ..

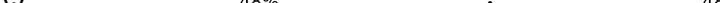


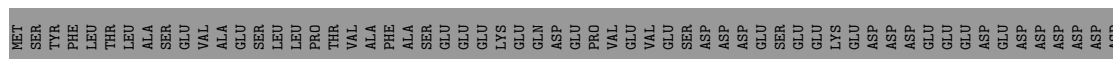
- Molecule 4: YALI0F24673p

Chain F:  47% 5% 48%

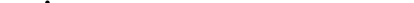


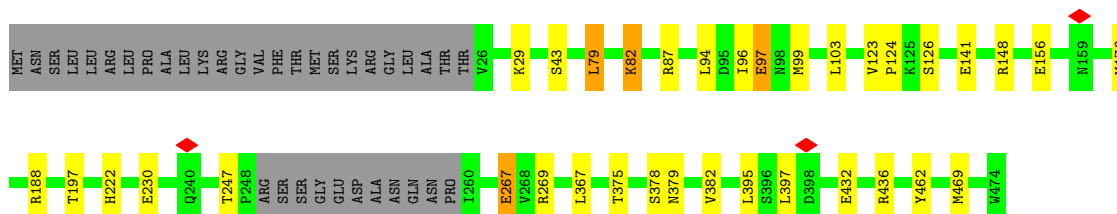
- Molecule 4: YALI0F24673p

Chain Q:  48% 2% 50%

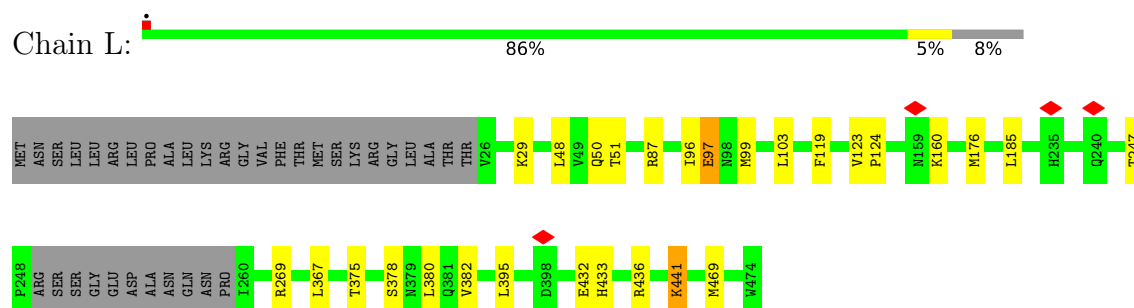


- Molecule 5: YALI0A14806p

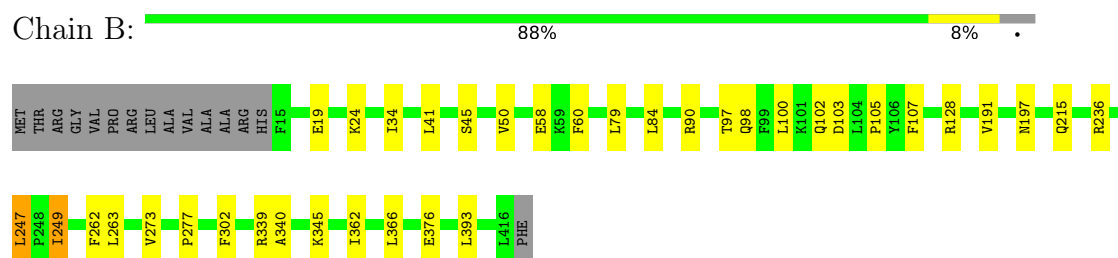
Chain A:  85% 7% 8%



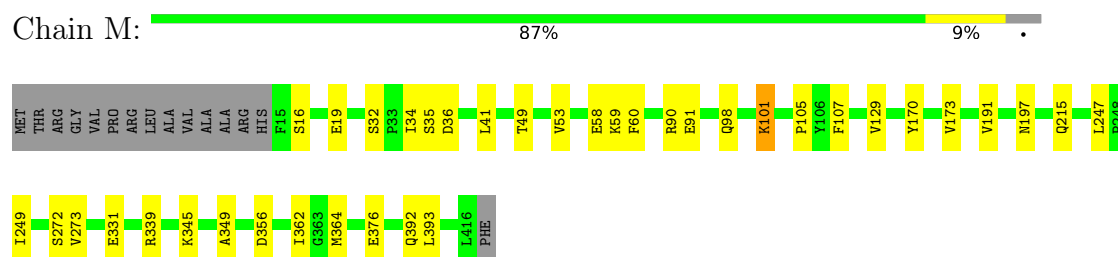
- Molecule 5: YALI0A14806p



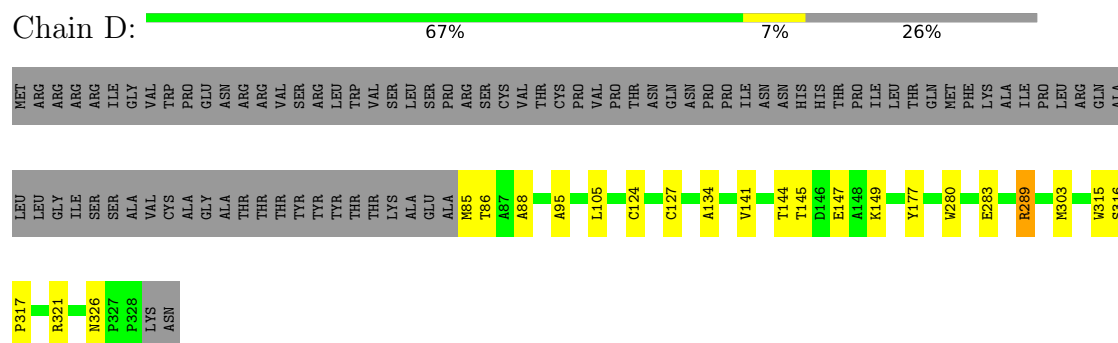
- Molecule 6: Cytochrome b-c1 complex subunit 2, mitochondrial



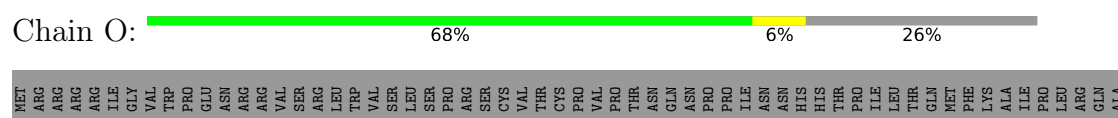
- Molecule 6: Cytochrome b-c1 complex subunit 2, mitochondrial

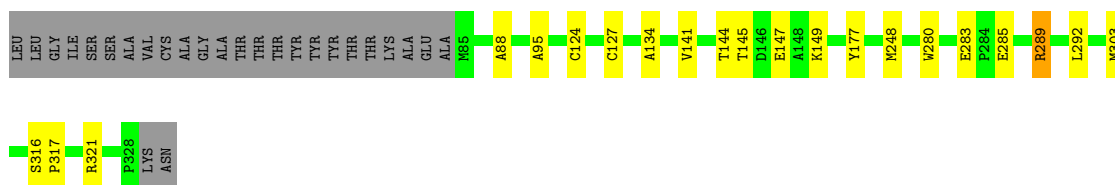


- Molecule 7: YALI0A17468p

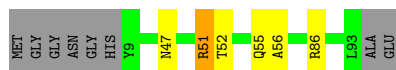


- Molecule 7: YALI0A17468p

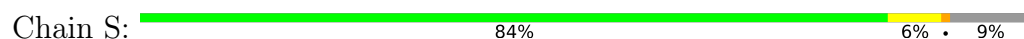




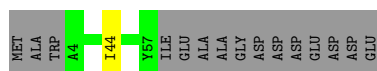
• Molecule 8: Cytochrome b-c1 complex subunit 8



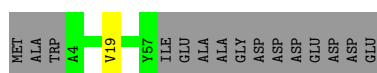
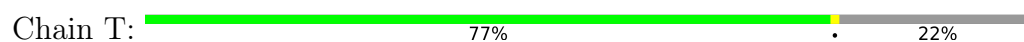
• Molecule 8: Cytochrome b-c1 complex subunit 8



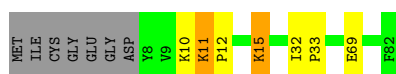
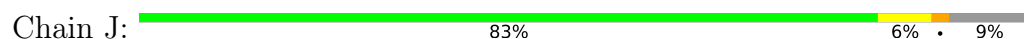
• Molecule 9: Complex III subunit 9



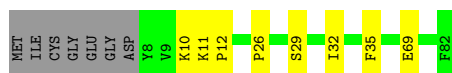
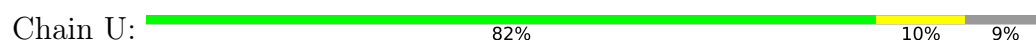
• Molecule 9: Complex III subunit 9



• Molecule 10: YALI0C12210p



• Molecule 10: YALI0C12210p



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	69259	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	55	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.088	Depositor
Minimum map value	-0.025	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.012	Depositor
Map size (Å)	301.32, 301.32, 301.32	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.837, 0.837, 0.837	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: DCQ, FES, PTY, PC1, HEC, HEM, LMT, XP4, CDL

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	C	0.60	0/3153	1.05	2/4305 (0.0%)
1	N	0.60	0/3153	1.04	1/4305 (0.0%)
2	E	0.58	0/461	1.09	1/619 (0.2%)
2	P	0.54	0/1479	1.10	4/2003 (0.2%)
3	G	0.57	0/1012	1.04	1/1373 (0.1%)
3	R	0.57	0/1012	1.06	1/1373 (0.1%)
4	F	0.49	0/595	1.05	0/805
4	Q	0.49	0/595	1.13	2/805 (0.2%)
5	A	0.54	0/3510	1.09	3/4768 (0.1%)
5	L	0.55	0/3510	1.10	3/4768 (0.1%)
6	B	0.55	0/3069	1.09	6/4178 (0.1%)
6	M	0.55	0/3069	1.10	7/4178 (0.2%)
7	D	0.58	0/1950	1.12	5/2656 (0.2%)
7	O	0.57	0/1950	1.10	3/2656 (0.1%)
8	H	0.57	0/717	1.07	1/975 (0.1%)
8	S	0.56	0/717	1.06	2/975 (0.2%)
9	I	0.54	0/465	1.10	0/629
9	T	0.54	0/465	1.09	0/629
10	J	0.54	0/620	1.00	0/846
10	U	0.54	0/620	1.03	0/846
All	All	0.56	0/32122	1.08	42/43692 (0.1%)

There are no bond length outliers.

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	D	283	GLU	CB-CA-C	-11.21	94.80	111.27
7	O	283	GLU	CB-CA-C	-10.88	95.28	111.27
6	M	215	GLN	CB-CA-C	-9.65	92.56	109.64
6	B	215	GLN	CB-CA-C	-8.92	92.67	110.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	90	ARG	CG-CD-NE	-8.66	92.94	112.00
7	D	283	GLU	CB-CG-CD	-8.35	98.41	112.60
6	M	90	ARG	CG-CD-NE	-7.47	95.56	112.00
2	P	62	ARG	CB-CG-CD	-7.11	94.94	111.30
4	Q	93	ASP	CA-CB-CG	6.98	119.58	112.60
4	Q	96	HIS	CB-CA-C	-6.74	99.60	110.79
3	G	76	ARG	NE-CZ-NH1	-6.46	115.04	121.50
3	R	76	ARG	NE-CZ-NH1	-6.40	115.10	121.50
2	P	181	ASP	N-CA-C	-6.33	105.01	114.64
1	N	49	THR	CA-CB-OG1	-6.22	100.27	109.60
5	A	379	ASN	CA-CB-CG	-6.16	106.44	112.60
6	B	107	PHE	CA-CB-CG	-6.08	107.72	113.80
5	A	269	ARG	CG-CD-NE	-6.03	98.73	112.00
6	B	197	ASN	CB-CA-C	6.01	119.51	111.82
2	E	62	ARG	CB-CG-CD	-6.00	97.50	111.30
1	C	49	THR	CA-CB-OG1	-5.93	100.71	109.60
6	M	392	GLN	CB-CG-CD	-5.87	102.62	112.60
6	M	197	ASN	CB-CA-C	5.81	119.53	111.63
6	M	49	THR	N-CA-C	-5.78	105.89	113.12
6	M	107	PHE	CA-CB-CG	-5.71	108.09	113.80
7	O	289	ARG	CG-CD-NE	-5.54	99.81	112.00
8	S	86	ARG	CB-CA-C	-5.50	102.25	110.88
7	D	289	ARG	CG-CD-NE	-5.43	100.05	112.00
2	P	197	SER	N-CA-C	-5.43	106.65	113.28
7	D	321	ARG	CB-CG-CD	-5.37	98.95	111.30
8	H	86	ARG	CB-CA-C	-5.32	102.53	110.88
5	L	441	LYS	N-CA-CB	-5.28	102.77	110.53
7	D	86	THR	CA-CB-OG1	-5.27	101.70	109.60
6	M	215	GLN	CB-CG-CD	-5.26	103.65	112.60
6	B	215	GLN	CB-CG-CD	-5.24	103.70	112.60
2	P	182	PHE	N-CA-C	-5.22	105.58	112.94
5	A	247	THR	CB-CA-C	5.21	116.85	109.08
8	S	71	THR	CA-CB-OG1	-5.20	101.80	109.60
5	L	269	ARG	CG-CD-NE	-5.19	100.58	112.00
1	C	283	ARG	CG-CD-NE	5.18	123.39	112.00
5	L	247	THR	CB-CA-C	5.14	116.74	109.08
7	O	321	ARG	CB-CG-CD	-5.11	99.54	111.30
6	B	128	ARG	CG-CD-NE	-5.05	100.88	112.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	3052	0	3113	22	0
1	N	3052	0	3113	19	0
2	E	452	0	443	4	0
2	P	1445	0	1426	39	0
3	G	994	0	1022	2	0
3	R	994	0	1022	4	0
4	F	579	0	511	3	0
4	Q	579	0	511	2	0
5	A	3446	0	3369	20	0
5	L	3446	0	3369	12	0
6	B	3008	0	2991	18	0
6	M	3008	0	2991	18	0
7	D	1893	0	1834	23	0
7	O	1893	0	1834	21	0
8	H	690	0	673	5	0
8	S	690	0	673	5	0
9	I	452	0	435	1	0
9	T	452	0	435	1	0
10	J	598	0	615	4	0
10	U	598	0	615	5	0
11	C	86	0	60	8	0
11	N	86	0	60	7	0
12	C	38	0	50	2	0
12	I	32	0	38	1	0
12	N	38	0	50	2	0
12	T	32	0	38	4	0
13	C	41	0	58	3	0
13	E	41	0	58	2	0
13	N	41	0	58	9	0
13	P	41	0	58	12	0
14	A	89	0	85	3	0
14	C	48	0	40	1	0
14	H	89	0	66	7	0
14	L	89	0	85	1	0
14	N	98	0	84	5	0
14	S	39	0	22	4	0
15	C	23	0	30	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
15	N	23	0	30	3	0
16	C	35	0	46	1	0
16	J	35	0	46	1	0
16	N	35	0	46	1	0
16	U	35	0	46	1	0
17	P	4	0	0	0	0
18	A	24	0	22	0	0
18	L	24	0	22	1	0
19	D	43	0	32	11	0
19	O	43	0	32	11	0
All	All	32573	0	32257	232	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 4.

All (232) 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)
2:P:84:VAL:CG2	13:P:302:PTY:H441	1.23	1.56
7:D:124:CYS:SG	19:D:401:HEC:HBB3	1.55	1.47
7:O:124:CYS:SG	19:O:401:HEC:HBB3	1.55	1.44
2:P:84:VAL:HG23	13:P:302:PTY:C44	1.62	1.29
2:P:84:VAL:CG2	13:P:302:PTY:C44	2.13	1.27
7:D:124:CYS:SG	19:D:401:HEC:CBB	2.27	1.23
7:O:124:CYS:SG	19:O:401:HEC:CBB	2.25	1.22
7:D:127:CYS:SG	19:D:401:HEC:HBC3	1.85	1.16
5:A:395:LEU:HD23	6:B:34:ILE:HD12	1.30	1.13
7:O:127:CYS:SG	19:O:401:HEC:HBC3	1.89	1.11
7:O:124:CYS:SG	19:O:401:HEC:CAB	2.42	1.06
7:O:127:CYS:SG	19:O:401:HEC:CAC	2.43	1.06
7:O:127:CYS:SG	19:O:401:HEC:CBC	2.44	1.05
2:P:84:VAL:HG22	13:P:302:PTY:H441	1.05	1.05
7:D:124:CYS:SG	19:D:401:HEC:CAB	2.46	1.03
7:D:127:CYS:SG	19:D:401:HEC:CBC	2.46	1.03
7:D:127:CYS:SG	19:D:401:HEC:CAC	2.49	1.01
2:P:84:VAL:HG23	13:P:302:PTY:H441	1.02	0.98
6:M:91:GLU:HG2	6:M:364:MET:HE1	1.50	0.93
1:C:330:ILE:HD12	12:C:503:PC1:H2A1	1.49	0.91
1:N:330:ILE:HD12	12:N:503:PC1:H2A1	1.52	0.91
10:U:11:LYS:HG2	10:U:12:PRO:HD2	1.53	0.88
2:P:84:VAL:HG22	13:P:302:PTY:C44	1.92	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:156:GLU:OE2	5:A:188:ARG:NH1	2.07	0.87
1:C:58:ALA:H	1:C:173:ASN:HD22	1.27	0.82
5:A:395:LEU:CD2	6:B:34:ILE:HD12	2.10	0.80
5:A:395:LEU:HD23	6:B:34:ILE:CD1	2.11	0.79
14:S:101:CDL:HB61	14:S:101:CDL:C72	2.14	0.78
11:N:501:HEM:HBC2	11:N:501:HEM:HHD	1.67	0.77
1:C:344:ILE:HD12	2:P:190:HIS:CD2	2.20	0.76
3:R:17:SER:HB2	3:R:20:LEU:HB2	1.67	0.76
1:C:344:ILE:HD12	2:P:190:HIS:HD2	1.50	0.76
3:G:17:SER:HB2	3:G:20:LEU:HB2	1.67	0.74
11:C:501:HEM:HBC2	11:C:501:HEM:HHD	1.69	0.72
14:A:3001:CDL:OB9	14:A:3001:CDL:HB4	1.90	0.70
2:P:84:VAL:CG2	13:P:302:PTY:C43	2.69	0.70
14:H:702:CDL:CB6	14:H:702:CDL:C72	2.70	0.69
8:H:51:ARG:HH21	14:H:701:CDL:HA22	1.56	0.69
2:P:84:VAL:HG23	13:P:302:PTY:C43	2.21	0.69
7:D:124:CYS:HG	19:D:401:HEC:HBB3	1.57	0.67
14:S:101:CDL:C72	14:S:101:CDL:CB6	2.73	0.66
11:C:501:HEM:HBC2	11:C:501:HEM:CHD	2.24	0.66
6:B:236:ARG:HD3	6:M:170:TYR:CE1	2.31	0.66
7:O:124:CYS:HG	19:O:401:HEC:HBB3	1.58	0.65
11:N:501:HEM:HBC2	11:N:501:HEM:CHD	2.23	0.64
7:O:124:CYS:SG	19:O:401:HEC:C3B	2.85	0.64
8:S:47:ASN:O	8:S:51:ARG:HG2	1.98	0.64
14:H:702:CDL:C72	14:H:702:CDL:HB61	2.28	0.63
1:C:58:ALA:H	1:C:173:ASN:ND2	1.94	0.63
6:B:41:LEU:CD2	6:B:191:VAL:HG22	2.28	0.62
2:P:180:GLY:HA2	2:P:186:PHE:HB2	1.80	0.62
7:D:124:CYS:SG	19:D:401:HEC:C3B	2.87	0.62
2:P:118:ILE:HD13	2:P:127:ILE:HD12	1.83	0.60
6:M:41:LEU:CD2	6:M:191:VAL:HG22	2.31	0.60
11:N:502:HEM:HBA1	15:N:507:DCQ:C3M	2.31	0.60
1:C:265:THR:HG21	2:P:173:CYS:HB3	1.83	0.60
2:P:186:PHE:CE1	2:P:191:GLY:HA2	2.37	0.59
1:C:265:THR:HG21	2:P:173:CYS:CB	2.32	0.59
11:C:502:HEM:HBA1	15:C:506:DCQ:C3M	2.33	0.59
4:Q:135:ASN:HB3	7:O:95:ALA:HB2	1.85	0.59
1:N:76:TRP:CZ3	7:O:289:ARG:HG3	2.38	0.59
5:L:97:GLU:HB3	6:M:339:ARG:HD3	1.85	0.58
7:D:145:THR:O	7:D:149:LYS:HG3	2.04	0.58
1:C:76:TRP:CZ3	7:D:289:ARG:HG3	2.37	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:A:3002:CDL:H132	14:A:3002:CDL:H522	1.85	0.58
15:C:506:DCQ:O5	15:C:506:DCQ:H8	2.04	0.58
14:H:701:CDL:OB3	14:H:702:CDL:O1	2.22	0.57
9:T:19:VAL:HG22	12:T:201:PC1:H332	1.86	0.57
14:N:504:CDL:H132	8:S:56:ALA:HA	1.86	0.57
14:N:504:CDL:HA21	8:S:51:ARG:HH21	1.69	0.57
8:H:47:ASN:O	8:H:51:ARG:HG2	2.05	0.56
5:A:378:SER:HA	5:A:432:GLU:OE1	2.06	0.56
2:P:155:ARG:NH1	2:P:197:SER:O	2.37	0.56
6:B:41:LEU:HD23	6:B:191:VAL:HG22	1.88	0.56
4:F:135:ASN:HB3	7:D:95:ALA:HB2	1.86	0.56
3:G:55:MET:HE1	3:G:102:LEU:HD21	1.88	0.56
7:O:127:CYS:SG	19:O:401:HEC:C3C	2.93	0.55
7:D:127:CYS:SG	19:D:401:HEC:C3C	2.95	0.55
1:C:180:PHE:HE2	1:N:180:PHE:HE2	1.53	0.55
6:B:236:ARG:HD3	6:M:170:TYR:CZ	2.42	0.54
11:C:501:HEM:HHD	11:C:501:HEM:CBC	2.37	0.54
5:A:43:SER:CB	5:A:222:HIS:HD2	2.20	0.54
11:N:501:HEM:HHD	11:N:501:HEM:CBC	2.37	0.54
13:C:504:PTY:H112	13:C:504:PTY:HC12	1.90	0.53
5:A:382:VAL:HG21	5:A:432:GLU:HA	1.91	0.53
5:L:382:VAL:HG21	5:L:432:GLU:HA	1.89	0.53
6:M:41:LEU:HD23	6:M:191:VAL:HG22	1.89	0.53
5:A:96:ILE:HG12	5:A:103:LEU:HD13	1.91	0.53
11:N:502:HEM:HBA1	15:N:507:DCQ:H3MA	1.90	0.53
1:C:234:PHE:CZ	7:D:303:MET:HE2	2.43	0.52
5:A:43:SER:HB3	5:A:222:HIS:HD2	1.72	0.52
6:M:247:LEU:HD13	6:M:393:LEU:HB3	1.91	0.52
6:M:53:VAL:HG12	6:M:173:VAL:HG13	1.91	0.52
6:B:24:LYS:HB3	6:B:366:LEU:HD22	1.92	0.51
8:H:56:ALA:HA	14:H:701:CDL:H132	1.92	0.51
11:C:502:HEM:HBC2	11:C:502:HEM:HMC2	1.93	0.50
5:L:395:LEU:HD12	6:M:34:ILE:CG2	2.41	0.50
6:B:247:LEU:HD13	6:B:393:LEU:HB3	1.92	0.50
1:N:139:MET:HE1	1:N:269:ILE:HA	1.92	0.50
4:F:124:VAL:HG11	7:D:88:ALA:HB2	1.92	0.50
1:N:229:ASP:HB2	13:N:505:PTY:H382	1.91	0.50
2:P:127:ILE:HG12	2:P:163:VAL:HG22	1.93	0.50
6:M:59:LYS:HB3	6:M:129:VAL:HG13	1.93	0.50
4:Q:124:VAL:HG11	7:O:88:ALA:HB2	1.94	0.50
14:A:3001:CDL:H1	14:A:3002:CDL:HA4	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:184:TYR:CD2	11:C:501:HEM:HBC1	2.46	0.49
6:B:249:ILE:HD11	6:B:302:PHE:HB2	1.94	0.49
1:C:344:ILE:CD1	2:P:190:HIS:HD2	2.20	0.49
1:N:234:PHE:CZ	7:O:303:MET:HE2	2.48	0.49
13:N:505:PTY:H112	12:T:201:PC1:H242	1.94	0.49
14:C:505:CDL:HB32	1:N:3:LEU:HD11	1.94	0.49
11:C:502:HEM:HBA1	15:C:506:DCQ:H3MB	1.95	0.49
14:L:3001:CDL:H1	14:L:3002:CDL:HA4	1.95	0.48
2:P:120:TRP:CE3	2:P:125:ILE:HD11	2.49	0.48
14:N:504:CDL:OB3	14:S:101:CDL:O1	2.31	0.48
13:E:401:PTY:H161	13:E:401:PTY:H132	1.51	0.48
2:P:116:VAL:HG23	2:P:127:ILE:O	2.14	0.48
5:L:96:ILE:HG12	5:L:103:LEU:HD13	1.96	0.48
2:P:45:ILE:HG21	5:L:176:MET:HE2	1.96	0.47
13:N:505:PTY:H111	12:T:201:PC1:O31	2.14	0.47
1:C:3:LEU:HD11	14:N:506:CDL:HB32	1.97	0.47
7:D:326:ASN:OD1	2:E:40:LYS:HA	2.15	0.47
1:N:184:TYR:CD2	11:N:501:HEM:HBC1	2.50	0.47
12:N:503:PC1:H31	8:S:55:GLN:HG3	1.96	0.47
13:N:505:PTY:H362	13:N:505:PTY:H331	1.48	0.47
13:P:302:PTY:H132	13:P:302:PTY:H161	1.77	0.47
6:B:236:ARG:HD3	6:M:170:TYR:CD1	2.49	0.47
6:B:84:LEU:HD12	6:B:97:THR:HG22	1.97	0.47
6:M:105:PRO:HB2	10:U:10:LYS:HA	1.96	0.47
1:C:34:SER:HB3	15:C:506:DCQ:H8	1.97	0.46
2:P:177:GLY:C	2:P:179:ALA:H	2.24	0.46
1:N:226:SER:HB3	13:N:505:PTY:H331	1.97	0.46
13:E:401:PTY:H381	13:E:401:PTY:H352	1.37	0.46
7:O:144:THR:OG1	7:O:147:GLU:HG3	2.16	0.46
7:D:316:SER:N	7:D:317:PRO:CD	2.79	0.46
10:U:29:SER:HA	10:U:32:ILE:HG12	1.98	0.46
2:P:102:LYS:H	2:P:102:LYS:HG3	1.52	0.46
5:A:97:GLU:HB3	6:B:339:ARG:HD2	1.97	0.46
14:H:702:CDL:HA62	14:H:702:CDL:H311	1.57	0.46
3:R:55:MET:HE1	3:R:102:LEU:HD21	1.97	0.46
12:C:503:PC1:H31	8:H:55:GLN:HG3	1.98	0.45
7:D:144:THR:OG1	7:D:147:GLU:HG3	2.16	0.45
1:N:49:THR:HG21	1:N:78:ILE:HD13	1.98	0.45
5:A:375:THR:C	5:A:436:ARG:HH12	2.25	0.45
15:N:507:DCQ:H7A	15:N:507:DCQ:H1MB	1.79	0.45
14:S:101:CDL:HA62	14:S:101:CDL:H311	1.53	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:P:302:PTY:H331	13:P:302:PTY:H361	1.51	0.45
6:M:249:ILE:CD1	6:M:393:LEU:HD21	2.47	0.45
12:I:201:PC1:H322	12:I:201:PC1:H351	1.63	0.45
6:M:249:ILE:HD11	6:M:393:LEU:HD21	1.99	0.45
7:O:316:SER:N	7:O:317:PRO:CD	2.80	0.45
5:A:82:LYS:HD3	5:A:82:LYS:HA	1.57	0.45
1:C:79:ARG:HD3	1:C:79:ARG:C	2.42	0.44
1:C:12:MET:HB3	16:C:507:LMT:H41	2.00	0.44
5:A:367:LEU:HD13	5:A:469:MET:SD	2.58	0.44
7:O:292:LEU:HD21	16:U:101:LMT:H6'2	2.00	0.44
1:C:230:LEU:HD11	13:C:504:PTY:H332	1.99	0.44
5:A:395:LEU:HD22	6:B:98:GLN:HG2	1.99	0.44
5:L:119:PHE:HE1	6:M:349:ALA:HB2	1.83	0.44
2:P:196:ILE:H	2:P:196:ILE:HG12	1.55	0.44
13:N:505:PTY:H372	13:N:505:PTY:H402	1.73	0.44
5:L:367:LEU:HD13	5:L:469:MET:SD	2.58	0.44
2:P:130:ARG:HD2	2:P:162:LEU:HB2	1.99	0.44
10:J:11:LYS:HD2	10:J:12:PRO:HD2	2.00	0.44
5:L:375:THR:C	5:L:436:ARG:HH12	2.24	0.44
6:M:36:ASP:OD1	6:M:98:GLN:HG3	2.16	0.44
2:P:66:TYR:CE1	12:T:201:PC1:H133	2.53	0.44
1:N:16:TYR:CE1	16:N:508:LMT:H42	2.52	0.44
11:N:502:HEM:HBC2	11:N:502:HEM:HMC2	2.00	0.44
5:L:395:LEU:HD12	6:M:34:ILE:HG22	1.99	0.44
7:O:145:THR:HG22	7:O:149:LYS:HD2	1.99	0.44
13:P:302:PTY:H141	13:P:302:PTY:H311	1.98	0.43
6:B:105:PRO:HB2	10:J:10:LYS:HA	1.99	0.43
1:N:227:PHE:HZ	13:N:505:PTY:HC6	1.83	0.43
5:A:99:MET:SD	5:A:126:SER:HB3	2.57	0.43
13:C:504:PTY:H132	2:E:72:PHE:CE2	2.53	0.43
2:P:192:SER:HA	2:P:203:GLY:HA3	2.00	0.43
5:L:123:VAL:N	5:L:124:PRO:CD	2.82	0.43
5:A:123:VAL:N	5:A:124:PRO:CD	2.82	0.43
6:B:100:LEU:HB2	6:B:103:ASP:OD2	2.17	0.43
1:N:79:ARG:HD3	1:N:79:ARG:C	2.42	0.43
1:C:49:THR:HG21	1:C:78:ILE:HD13	2.01	0.43
5:A:79:LEU:HD12	5:A:79:LEU:HA	1.88	0.43
6:B:263:LEU:HA	6:B:277:PRO:HG2	2.00	0.43
10:J:15:LYS:HB2	10:J:15:LYS:HE3	1.77	0.43
5:L:160:LYS:HA	5:L:160:LYS:HD2	1.92	0.43
2:P:106:LYS:HA	2:P:221:THR:HG22	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:O:134:ALA:HA	7:O:177:TYR:HA	2.01	0.43
2:P:177:GLY:HA2	2:P:185:TRP:CD1	2.54	0.42
2:P:102:LYS:HA	2:P:224:ILE:O	2.19	0.42
3:R:20:LEU:HD23	3:R:20:LEU:HA	1.89	0.42
2:P:193:HIS:O	2:P:200:ILE:HD12	2.19	0.42
3:R:71:VAL:HG11	8:S:21:GLN:HG2	2.02	0.42
1:C:345:GLU:OE2	7:D:85:MET:N	2.53	0.42
7:D:134:ALA:HA	7:D:177:TYR:HA	2.01	0.42
1:N:108:THR:HA	14:N:506:CDL:OA4	2.20	0.42
2:P:147:ARG:HD2	2:P:206:PRO:HA	2.02	0.42
10:J:32:ILE:HB	10:J:33:PRO:HD3	2.01	0.42
1:N:226:SER:HB3	13:N:505:PTY:H321	2.02	0.42
5:A:148:ARG:HH11	5:A:148:ARG:HD3	1.65	0.42
7:D:127:CYS:HG	19:D:401:HEC:HBC3	1.77	0.42
7:O:141:VAL:HG11	7:O:280:TRP:CE2	2.55	0.42
1:C:224:TYR:HB3	7:D:315:TRP:CZ2	2.55	0.41
2:P:146:LEU:O	2:P:147:ARG:C	2.63	0.41
2:P:171:LEU:HD12	2:P:190:HIS:CE1	2.54	0.41
13:P:302:PTY:H412	13:P:302:PTY:H443	1.63	0.41
7:O:248:MET:HB2	19:O:401:HEC:C1D	2.50	0.41
5:A:176:MET:HE2	2:E:45:ILE:HG21	2.02	0.41
1:N:131:GLY:HA3	1:N:183:HIS:CE1	2.56	0.41
2:P:102:LYS:HG3	2:P:102:LYS:HZ3	1.63	0.41
19:O:401:HEC:HHC	19:O:401:HEC:HAB	1.77	0.41
18:L:3003:XP4:H8	10:U:35:PHE:HE1	1.85	0.41
7:D:141:VAL:HG11	7:D:280:TRP:CE2	2.56	0.41
19:D:401:HEC:HAB	19:D:401:HEC:HHC	1.83	0.41
5:L:380:LEU:HD11	10:U:26:PRO:HG2	2.02	0.41
2:P:49:THR:N	2:P:50:PRO:CD	2.84	0.41
2:P:125:ILE:HD13	2:P:165:ILE:HA	2.03	0.41
2:P:156:VAL:HG21	2:P:159:PRO:HA	2.02	0.41
1:C:99:ARG:NH2	11:C:502:HEM:HBD1	2.36	0.41
1:C:208:ASN:HB2	1:C:209:PRO:HD2	2.03	0.41
5:A:267:GLU:OE1	5:A:462:TYR:HB2	2.20	0.41
8:H:52:THR:HG23	14:H:701:CDL:H112	2.03	0.41
1:N:227:PHE:CZ	13:N:505:PTY:H322	2.56	0.41
4:F:89:GLY:HA3	4:F:90:PRO:HD3	1.97	0.41
7:D:105:LEU:HB3	9:I:44:ILE:HD12	2.03	0.41
1:N:346:PRO:HA	1:N:347:PRO:HA	1.88	0.41
6:M:35:SER:OG	6:M:101:LYS:HG2	2.21	0.41
16:J:101:LMT:H52	16:J:101:LMT:H21	1.75	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:76:TRP:CG	7:O:285:GLU:HG3	2.56	0.40
2:P:103:VAL:HG22	2:P:120:TRP:CD1	2.56	0.40
2:E:49:THR:N	2:E:50:PRO:CD	2.84	0.40
6:B:262:PHE:HB2	6:B:340:ALA:HB2	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	C	381/385 (99%)	374 (98%)	7 (2%)	0	100	100
1	N	381/385 (99%)	374 (98%)	7 (2%)	0	100	100
2	E	57/225 (25%)	56 (98%)	1 (2%)	0	100	100
2	P	184/225 (82%)	172 (94%)	12 (6%)	0	100	100
3	G	122/128 (95%)	122 (100%)	0	0	100	100
3	R	122/128 (95%)	122 (100%)	0	0	100	100
4	F	69/137 (50%)	67 (97%)	2 (3%)	0	100	100
4	Q	69/137 (50%)	67 (97%)	2 (3%)	0	100	100
5	A	434/474 (92%)	424 (98%)	10 (2%)	0	100	100
5	L	434/474 (92%)	423 (98%)	11 (2%)	0	100	100
6	B	400/417 (96%)	384 (96%)	16 (4%)	0	100	100
6	M	400/417 (96%)	386 (96%)	14 (4%)	0	100	100
7	D	242/330 (73%)	238 (98%)	4 (2%)	0	100	100
7	O	242/330 (73%)	238 (98%)	4 (2%)	0	100	100
8	H	83/93 (89%)	82 (99%)	1 (1%)	0	100	100
8	S	83/93 (89%)	82 (99%)	1 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
9	I	52/69 (75%)	51 (98%)	1 (2%)	0	100	100
9	T	52/69 (75%)	51 (98%)	1 (2%)	0	100	100
10	J	73/82 (89%)	71 (97%)	2 (3%)	0	100	100
10	U	73/82 (89%)	71 (97%)	2 (3%)	0	100	100
All	All	3953/4680 (84%)	3855 (98%)	98 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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	C	331/333 (99%)	326 (98%)	5 (2%)	57	80
1	N	331/333 (99%)	325 (98%)	6 (2%)	51	77
2	E	48/182 (26%)	44 (92%)	4 (8%)	10	22
2	P	154/182 (85%)	137 (89%)	17 (11%)	6	13
3	G	113/117 (97%)	105 (93%)	8 (7%)	13	29
3	R	113/117 (97%)	108 (96%)	5 (4%)	25	50
4	F	61/123 (50%)	58 (95%)	3 (5%)	22	45
4	Q	61/123 (50%)	60 (98%)	1 (2%)	55	79
5	A	377/407 (93%)	366 (97%)	11 (3%)	37	65
5	L	377/407 (93%)	366 (97%)	11 (3%)	37	65
6	B	311/322 (97%)	298 (96%)	13 (4%)	26	52
6	M	311/322 (97%)	298 (96%)	13 (4%)	26	52
7	D	192/268 (72%)	192 (100%)	0	100	100
7	O	192/268 (72%)	192 (100%)	0	100	100
8	H	67/71 (94%)	66 (98%)	1 (2%)	57	80
8	S	67/71 (94%)	66 (98%)	1 (2%)	57	80
9	I	46/57 (81%)	46 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
9	T	46/57 (81%)	46 (100%)	0	100	100
10	J	63/68 (93%)	60 (95%)	3 (5%)	23	46
10	U	63/68 (93%)	62 (98%)	1 (2%)	55	79
All	All	3324/3896 (85%)	3221 (97%)	103 (3%)	36	62

All (103) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	146	VAL
1	C	178	ARG
1	C	197	HIS
1	C	288	LYS
1	C	324	LEU
2	P	102	LYS
2	P	106	LYS
2	P	114	LYS
2	P	115	ASN
2	P	116	VAL
2	P	117	ILE
2	P	118	ILE
2	P	119	LYS
2	P	131	THR
2	P	143	VAL
2	P	146	LEU
2	P	147	ARG
2	P	158	LYS
2	P	176	ILE
2	P	196	ILE
2	P	197	SER
2	P	213	GLU
3	G	3	SER
3	G	6	SER
3	G	17	SER
3	G	19	LEU
3	G	22	LYS
3	G	23	ILE
3	G	50	GLU
3	G	123	GLU
4	F	92	HIS
4	F	96	HIS
4	F	100	GLU

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Mol	Chain	Res	Type
5	A	29	LYS
5	A	79	LEU
5	A	82	LYS
5	A	87	ARG
5	A	94	LEU
5	A	97	GLU
5	A	141	GLU
5	A	197	THR
5	A	230	GLU
5	A	267	GLU
5	A	397	LEU
6	B	19	GLU
6	B	45	SER
6	B	50	VAL
6	B	58	GLU
6	B	60	PHE
6	B	79	LEU
6	B	102	GLN
6	B	247	LEU
6	B	249	ILE
6	B	273	VAL
6	B	345	LYS
6	B	362	ILE
6	B	376	GLU
8	H	51	ARG
10	J	11	LYS
10	J	15	LYS
10	J	69	GLU
1	N	178	ARG
1	N	197	HIS
1	N	255	ASP
1	N	270	VAL
1	N	288	LYS
1	N	324	LEU
2	E	54	LYS
2	E	55	ASP
2	E	62	ARG
2	E	65	SER
3	R	17	SER
3	R	19	LEU
3	R	22	LYS
3	R	23	ILE

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Mol	Chain	Res	Type
3	R	32	VAL
4	Q	100	GLU
5	L	29	LYS
5	L	48	LEU
5	L	50	GLN
5	L	51	THR
5	L	87	ARG
5	L	97	GLU
5	L	99	MET
5	L	185	LEU
5	L	378	SER
5	L	433	HIS
5	L	441	LYS
6	M	16	SER
6	M	19	GLU
6	M	32	SER
6	M	58	GLU
6	M	60	PHE
6	M	101	LYS
6	M	272	SER
6	M	273	VAL
6	M	331	GLU
6	M	345	LYS
6	M	356	ASP
6	M	362	ILE
6	M	376	GLU
8	S	51	ARG
10	U	69	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	173	ASN
1	C	177	GLN
1	C	202	HIS
1	C	332	ASN
2	P	90	ASN
2	P	157	GLN
4	F	92	HIS
4	F	97	HIS
4	F	109	GLN
4	F	132	HIS

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Mol	Chain	Res	Type
5	A	222	HIS
5	A	316	HIS
5	A	410	GLN
6	B	28	GLN
6	B	63	GLN
6	B	122	GLN
6	B	189	GLN
6	B	215	GLN
6	B	408	HIS
7	D	193	GLN
7	D	208	HIS
10	J	70	HIS
1	N	177	GLN
1	N	202	HIS
1	N	332	ASN
1	N	343	HIS
2	E	90	ASN
3	R	84	HIS
4	Q	132	HIS
5	L	410	GLN
6	M	28	GLN
6	M	63	GLN
6	M	122	GLN
6	M	184	GLN
6	M	215	GLN
6	M	408	HIS
7	O	90	HIS
7	O	193	GLN
7	O	208	HIS
10	U	70	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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](#)

33 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$
11	HEM	N	502	1	50,50,50	1.55	9 (18%)	67,82,82	2.54	27 (40%)
19	HEC	D	401	7	46,50,50	2.46	22 (47%)	58,82,82	2.31	26 (44%)
18	XP4	L	3003	-	23,23,39	1.45	2 (8%)	26,28,44	2.07	7 (26%)
15	DCQ	C	506	-	23,23,23	0.84	1 (4%)	27,29,29	2.06	5 (18%)
18	XP4	A	3003	-	23,23,39	1.46	2 (8%)	26,28,44	2.18	6 (23%)
16	LMT	C	507	-	36,36,36	0.45	0	47,47,47	1.27	6 (12%)
16	LMT	U	101	-	36,36,36	0.43	0	47,47,47	0.94	4 (8%)
14	CDL	N	506	-	47,47,99	0.45	0	53,59,111	0.75	2 (3%)
16	LMT	J	101	-	36,36,36	0.53	0	47,47,47	1.41	4 (8%)
14	CDL	H	701	-	49,49,99	0.37	0	55,61,111	0.80	2 (3%)
13	PTY	N	505	-	40,40,49	0.33	0	43,45,54	0.73	2 (4%)
14	CDL	S	101	-	38,38,99	0.41	0	44,50,111	1.30	8 (18%)
14	CDL	H	702	-	38,38,99	0.40	0	44,50,111	1.17	3 (6%)
11	HEM	N	501	1	50,50,50	1.56	10 (20%)	67,82,82	1.99	20 (29%)
12	PC1	N	503	-	37,37,53	0.76	1 (2%)	43,45,61	0.98	4 (9%)
14	CDL	A	3002	-	46,46,99	0.36	0	51,56,111	0.95	3 (5%)
14	CDL	N	504	-	49,49,99	0.35	0	55,61,111	0.69	1 (1%)
13	PTY	C	504	-	40,40,49	0.34	0	43,45,54	0.76	0
14	CDL	L	3001	-	41,41,99	0.42	0	45,51,111	0.85	2 (4%)
14	CDL	L	3002	-	46,46,99	0.34	0	51,56,111	0.92	3 (5%)
11	HEM	C	501	1	50,50,50	1.47	7 (14%)	67,82,82	2.02	20 (29%)
15	DCQ	N	507	-	23,23,23	0.79	1 (4%)	27,29,29	1.53	5 (18%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
11	HEM	C	502	1	50,50,50	1.57	9 (18%)	67,82,82	2.49	26 (38%)
12	PC1	I	201	-	31,31,53	0.37	0	37,39,61	0.75	1 (2%)
12	PC1	T	201	-	31,31,53	0.37	0	37,39,61	0.51	0
14	CDL	A	3001	-	41,41,99	0.48	0	45,51,111	0.73	1 (2%)
13	PTY	E	401	-	40,40,49	0.48	0	43,45,54	0.51	0
12	PC1	C	503	-	37,37,53	0.71	1 (2%)	43,45,61	1.05	4 (9%)
17	FES	P	301	2	0,4,4	-	-	-	-	-
14	CDL	C	505	-	47,47,99	0.43	0	53,59,111	0.78	2 (3%)
16	LMT	N	508	-	36,36,36	0.40	0	47,47,47	1.22	6 (12%)
19	HEC	O	401	7	46,50,50	2.45	21 (45%)	58,82,82	2.30	26 (44%)
13	PTY	P	302	-	40,40,49	0.46	0	43,45,54	0.71	1 (2%)

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
11	HEM	N	502	1	-	4/14/54/54	-
19	HEC	D	401	7	-	4/14/54/54	-
18	XP4	L	3003	-	-	1/24/24/41	-
15	DCQ	C	506	-	-	8/14/38/38	0/1/1/1
18	XP4	A	3003	-	-	0/24/24/41	-
16	LMT	C	507	-	-	14/21/61/61	0/2/2/2
16	LMT	U	101	-	-	10/21/61/61	0/2/2/2
14	CDL	N	506	-	-	27/57/57/110	-
16	LMT	J	101	-	-	10/21/61/61	0/2/2/2
14	CDL	H	701	-	-	29/59/59/110	-
13	PTY	N	505	-	-	29/44/44/53	-
14	CDL	S	101	-	-	27/48/48/110	-
14	CDL	H	702	-	-	25/48/48/110	-
11	HEM	N	501	1	-	5/14/54/54	-
12	PC1	N	503	-	-	12/41/41/57	-
14	CDL	A	3002	-	-	30/54/54/110	-
14	CDL	N	504	-	-	33/59/59/110	-
13	PTY	C	504	-	-	30/44/44/53	-
14	CDL	L	3001	-	-	21/48/48/110	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
14	CDL	L	3002	-	-	29/54/54/110	-
11	HEM	C	501	1	-	5/14/54/54	-
15	DCQ	N	507	-	-	8/14/38/38	0/1/1/1
11	HEM	C	502	1	-	4/14/54/54	-
12	PC1	I	201	-	-	14/35/35/57	-
12	PC1	T	201	-	-	11/35/35/57	-
14	CDL	A	3001	-	-	14/48/48/110	-
13	PTY	E	401	-	-	24/44/44/53	-
12	PC1	C	503	-	-	11/41/41/57	-
17	FES	P	301	2	-	-	0/1/1/1
14	CDL	C	505	-	-	27/57/57/110	-
16	LMT	N	508	-	-	12/21/61/61	0/2/2/2
19	HEC	O	401	7	-	3/14/54/54	-
13	PTY	P	302	-	-	23/44/44/53	-

All (86) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
18	A	3003	XP4	O7-C18	5.26	1.46	1.35
19	D	401	HEC	C4B-NB	-5.20	1.29	1.39
18	L	3003	XP4	O7-C18	5.14	1.46	1.35
19	O	401	HEC	C4B-NB	-5.14	1.29	1.39
11	C	501	HEM	FE-NB	5.13	2.10	1.94
11	N	501	HEM	FE-NB	5.00	2.10	1.94
19	O	401	HEC	C1B-NB	-4.79	1.30	1.39
11	C	502	HEM	FE-NB	4.77	2.09	1.94
19	D	401	HEC	C1B-NB	-4.36	1.31	1.39
19	O	401	HEC	C4A-NA	-4.23	1.31	1.39
19	O	401	HEC	CHD-C4C	4.19	1.46	1.38
19	D	401	HEC	C4D-ND	-4.13	1.31	1.39
19	O	401	HEC	CHA-C1A	4.10	1.46	1.38
19	D	401	HEC	C4A-NA	-4.00	1.32	1.39
11	N	501	HEM	C1B-NB	-3.95	1.33	1.40
19	D	401	HEC	CHA-C1A	3.90	1.46	1.38
11	N	502	HEM	FE-NB	3.83	2.06	1.94
19	O	401	HEC	C2A-C3A	3.81	1.45	1.36
19	D	401	HEC	C2A-C3A	3.81	1.45	1.36
11	N	502	HEM	C1B-NB	-3.76	1.33	1.40
19	D	401	HEC	CAC-C3C	3.74	1.47	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
19	O	401	HEC	CAC-C3C	3.68	1.47	1.35
19	D	401	HEC	CHD-C4C	3.66	1.45	1.38
18	L	3003	XP4	O5-C4	3.63	1.43	1.33
18	A	3003	XP4	O5-C4	3.62	1.43	1.33
19	D	401	HEC	C4C-NC	-3.61	1.32	1.39
19	D	401	HEC	CHB-C4A	3.59	1.45	1.38
19	O	401	HEC	C4D-ND	-3.58	1.32	1.39
11	C	502	HEM	C3C-C4C	-3.53	1.39	1.46
11	N	501	HEM	FE-NC	3.49	2.06	1.95
19	D	401	HEC	C1A-NA	-3.49	1.33	1.39
11	C	502	HEM	C1B-NB	-3.32	1.34	1.40
19	D	401	HEC	CHC-C4B	3.31	1.44	1.38
19	O	401	HEC	C1A-NA	-3.29	1.33	1.39
11	C	501	HEM	FE-NC	3.27	2.06	1.95
11	N	502	HEM	CMA-C3A	3.26	1.57	1.50
11	N	502	HEM	C4D-ND	-3.26	1.34	1.40
19	O	401	HEC	CHB-C4A	3.26	1.44	1.38
11	C	502	HEM	FE-NC	3.26	2.05	1.95
19	D	401	HEC	CHA-C4D	3.26	1.46	1.39
11	N	502	HEM	C4B-NB	-3.24	1.32	1.38
19	D	401	HEC	C1D-ND	-3.23	1.33	1.39
19	O	401	HEC	C4C-NC	-3.21	1.33	1.39
11	C	501	HEM	C1B-NB	-3.17	1.34	1.40
11	C	502	HEM	C1C-C2C	-3.16	1.39	1.45
19	O	401	HEC	CHA-C4D	3.13	1.46	1.39
19	D	401	HEC	C1C-NC	-3.13	1.33	1.39
19	O	401	HEC	CHC-C4B	3.08	1.44	1.38
11	N	502	HEM	C1C-C2C	-3.02	1.39	1.45
11	C	502	HEM	C4B-NB	-3.01	1.32	1.38
11	N	501	HEM	C4B-NB	-2.92	1.33	1.38
11	N	502	HEM	FE-NC	2.91	2.04	1.95
19	O	401	HEC	CHD-C1D	2.86	1.45	1.39
11	C	501	HEM	C4B-NB	-2.81	1.33	1.38
19	O	401	HEC	C1C-NC	-2.79	1.34	1.39
19	O	401	HEC	CAB-C3B	2.79	1.44	1.35
11	C	502	HEM	C4D-ND	-2.74	1.35	1.40
19	O	401	HEC	C1D-ND	-2.71	1.34	1.39
19	D	401	HEC	CAB-C3B	2.68	1.43	1.35
19	O	401	HEC	C3D-C2D	2.68	1.45	1.38
11	N	501	HEM	C1C-NC	-2.67	1.34	1.39
19	O	401	HEC	C3B-C4B	-2.56	1.41	1.46
19	D	401	HEC	CHD-C1D	2.55	1.45	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	C	502	HEM	CMA-C3A	2.51	1.55	1.50
19	D	401	HEC	C3B-C4B	-2.47	1.41	1.46
15	C	506	DCQ	C4-C3	2.47	1.45	1.36
11	N	501	HEM	C4D-ND	-2.46	1.36	1.40
11	C	501	HEM	C1C-NC	-2.45	1.35	1.39
19	D	401	HEC	CHC-C1C	2.37	1.44	1.39
11	C	501	HEM	O2D-CGD	-2.36	1.23	1.30
19	D	401	HEC	C3D-C2D	2.35	1.45	1.38
19	O	401	HEC	CHC-C1C	2.32	1.44	1.39
11	N	501	HEM	C1D-C2D	2.30	1.49	1.44
11	N	501	HEM	O2D-CGD	-2.30	1.23	1.30
12	N	503	PC1	C22-C21	2.29	1.57	1.50
11	N	502	HEM	C3C-C4C	-2.20	1.42	1.46
11	N	502	HEM	FE-ND	-2.19	1.88	1.94
19	O	401	HEC	CHB-C1B	2.15	1.44	1.39
19	D	401	HEC	C1C-C2C	2.14	1.48	1.43
11	N	501	HEM	C4A-NA	-2.12	1.35	1.39
11	N	501	HEM	C1C-C2C	-2.11	1.41	1.45
11	C	502	HEM	CMD-C2D	2.10	1.55	1.50
11	C	501	HEM	C1D-C2D	2.08	1.48	1.44
12	C	503	PC1	C22-C21	2.08	1.56	1.50
15	N	507	DCQ	C4-C5	-2.04	1.42	1.48
19	D	401	HEC	CHB-C1B	2.00	1.43	1.39

All (227) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	C	502	HEM	CHD-C4C-NC	9.24	134.52	124.45
11	N	502	HEM	CHD-C4C-NC	7.87	133.02	124.45
18	A	3003	XP4	O7-C18-C19	7.62	124.68	111.09
11	N	502	HEM	CHC-C4B-NB	7.59	132.59	124.42
18	L	3003	XP4	O7-C18-C19	6.96	123.50	111.09
11	C	502	HEM	CHC-C4B-NB	6.83	131.77	124.42
19	D	401	HEC	CAA-CBA-CGA	-5.82	98.24	113.67
15	C	506	DCQ	C1-C6-C5	5.72	125.18	119.62
16	J	101	LMT	C1-O1'-C1'	-5.66	104.01	113.68
11	N	502	HEM	CHD-C1D-ND	5.64	130.49	124.42
19	O	401	HEC	CAA-CBA-CGA	-5.32	99.56	113.67
11	C	501	HEM	CHD-C1D-ND	5.26	130.08	124.42
15	C	506	DCQ	C7-C6-C5	5.20	125.56	116.92
19	D	401	HEC	C2A-C1A-NA	5.16	115.30	110.32
18	A	3003	XP4	O7-C18-O8	-4.95	113.43	122.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	C	501	HEM	CHC-C4B-NB	4.94	129.73	124.42
19	D	401	HEC	C2B-C1B-NB	4.92	118.03	110.14
11	N	501	HEM	CHC-C4B-NB	4.83	129.63	124.42
19	O	401	HEC	C2B-C1B-NB	4.77	117.80	110.14
11	N	502	HEM	C4C-CHD-C1D	-4.75	115.93	126.02
11	N	502	HEM	CHD-C1D-C2D	-4.72	117.57	125.03
19	O	401	HEC	CBC-CAC-C3C	-4.66	118.11	127.43
19	O	401	HEC	C2A-C1A-NA	4.66	114.82	110.32
11	N	501	HEM	CHD-C1D-ND	4.55	129.32	124.42
18	L	3003	XP4	O7-C18-O8	-4.50	114.30	122.99
11	C	502	HEM	C1B-NB-C4B	4.48	110.52	105.21
11	N	502	HEM	CHA-C4D-ND	4.44	129.86	124.37
11	C	502	HEM	CHA-C4D-ND	4.42	129.84	124.37
11	C	501	HEM	CBA-CAA-C2A	-4.39	100.40	112.53
11	C	502	HEM	C4C-CHD-C1D	-4.39	116.70	126.02
11	C	502	HEM	CHD-C1D-ND	4.37	129.13	124.42
11	N	501	HEM	CBA-CAA-C2A	-4.28	100.70	112.53
11	N	502	HEM	O2A-CGA-O1A	-4.27	112.36	123.33
19	D	401	HEC	C1D-C2D-C3D	-4.25	101.95	106.82
11	N	502	HEM	C1B-NB-C4B	4.24	110.22	105.21
16	J	101	LMT	O5'-C1'-O1'	-4.22	100.08	110.04
15	C	506	DCQ	C1M-C1-C6	-4.06	117.78	124.45
14	L	3001	CDL	OB6-CB4-CB3	-3.95	96.04	109.70
11	C	501	HEM	CHA-C4D-ND	3.93	129.23	124.37
19	O	401	HEC	C1D-C2D-C3D	-3.88	102.37	106.82
11	N	501	HEM	CHA-C4D-ND	3.87	129.16	124.37
19	D	401	HEC	C1A-C2A-C3A	-3.86	102.02	107.11
11	C	502	HEM	C4C-NC-C1C	3.80	112.02	105.82
11	C	502	HEM	CHD-C1D-C2D	-3.79	119.05	125.03
19	D	401	HEC	CMD-C2D-C1D	3.72	131.09	125.42
11	C	501	HEM	C1B-NB-C4B	3.72	109.61	105.21
11	C	501	HEM	CHA-C1A-NA	3.70	130.56	123.86
14	S	101	CDL	CB4-OB6-CB5	3.65	124.28	117.85
15	N	507	DCQ	O4-C4-C5	-3.64	104.34	116.64
11	C	501	HEM	C1A-CHA-C4D	-3.61	117.75	126.25
19	O	401	HEC	C2C-C1C-NC	3.61	115.93	110.14
19	O	401	HEC	C1A-C2A-C3A	-3.57	102.41	107.11
15	C	506	DCQ	C9-C8-C7	-3.54	100.13	113.13
11	N	501	HEM	CBC-CAC-C3C	-3.53	109.90	127.53
11	C	501	HEM	CBC-CAC-C3C	-3.52	109.95	127.53
11	N	502	HEM	CHB-C1B-NB	3.49	128.69	124.37
19	D	401	HEC	CMB-C2B-C3B	3.48	134.74	126.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	C	502	HEM	C4C-C3C-C2C	3.47	109.83	106.81
19	O	401	HEC	CMB-C2B-C3B	3.45	134.66	126.55
19	D	401	HEC	CAA-C2A-C1A	3.43	131.83	124.85
11	C	502	HEM	C3B-C4B-NB	-3.43	107.00	109.47
11	N	502	HEM	C1C-CHC-C4B	-3.42	118.74	126.02
14	H	702	CDL	CB4-OB6-CB5	3.30	123.68	117.85
14	S	101	CDL	OA5-PA1-OA3	-3.30	95.84	108.94
12	C	503	PC1	O22-C21-C22	3.28	136.62	123.78
11	N	501	HEM	C1B-NB-C4B	3.28	109.09	105.21
19	O	401	HEC	C3D-C4D-ND	3.25	113.75	110.15
19	O	401	HEC	CHC-C4B-C3B	-3.19	119.83	125.21
16	N	508	LMT	C1B-O5B-C5B	3.17	119.92	113.72
11	C	501	HEM	CHD-C1D-C2D	-3.15	120.05	125.03
11	C	502	HEM	O2A-CGA-O1A	-3.15	115.23	123.33
12	C	503	PC1	O21-C21-C22	-3.15	104.68	111.48
14	H	702	CDL	OB6-CB4-CB6	3.14	119.62	108.34
11	N	502	HEM	C1A-CHA-C4D	-3.14	118.87	126.25
11	C	502	HEM	CHB-C1B-NB	3.13	128.24	124.37
19	D	401	HEC	CBD-CAD-C3D	-3.12	103.91	112.53
11	C	501	HEM	C2A-C1A-NA	-3.10	106.71	110.15
15	N	507	DCQ	O4-C4-C3	3.08	135.30	123.64
16	C	507	LMT	O5B-C5B-C4B	3.08	115.24	109.70
11	N	502	HEM	O2D-CGD-CBD	3.06	123.68	114.00
11	N	502	HEM	O2A-CGA-CBA	3.06	123.68	114.00
11	N	501	HEM	CHC-C1C-NC	3.06	127.78	124.45
11	N	501	HEM	CHA-C4D-C3D	-3.06	119.59	125.23
19	D	401	HEC	C3D-C4D-ND	3.04	113.53	110.15
11	C	502	HEM	CHA-C4D-C3D	-3.04	119.63	125.23
18	A	3003	XP4	O5-C4-O6	-3.04	116.03	123.63
11	C	501	HEM	CAC-C3C-C4C	3.01	132.02	124.82
14	S	101	CDL	OB6-CB4-CB6	3.01	119.13	108.34
11	C	502	HEM	C1D-C2D-C3D	-2.98	103.84	106.98
19	D	401	HEC	C2C-C1C-NC	2.98	114.92	110.14
11	N	501	HEM	CAC-C3C-C4C	2.98	131.93	124.82
19	D	401	HEC	C2D-C1D-ND	2.97	114.91	110.14
19	O	401	HEC	CMD-C2D-C1D	2.97	129.94	125.42
14	L	3002	CDL	OA5-PA1-OA3	-2.94	97.26	108.94
19	D	401	HEC	CHB-C1B-C2B	-2.94	118.90	127.43
14	A	3002	CDL	OA6-CA4-CA6	2.93	118.84	108.34
19	O	401	HEC	CAA-C2A-C1A	2.91	130.77	124.85
11	N	502	HEM	CMA-C3A-C4A	2.91	129.85	125.42
11	N	502	HEM	CBD-CAD-C3D	-2.91	104.50	112.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
19	O	401	HEC	CBD-CAD-C3D	-2.89	104.53	112.53
11	N	501	HEM	CHD-C1D-C2D	-2.89	120.47	125.03
11	C	502	HEM	CBD-CAD-C3D	-2.87	104.61	112.53
19	O	401	HEC	CHB-C1B-C2B	-2.86	119.12	127.43
15	C	506	DCQ	C3-C4-C5	-2.85	115.46	120.69
11	C	502	HEM	O2A-CGA-CBA	2.84	122.99	114.00
12	N	503	PC1	O22-C21-C22	2.84	134.91	123.78
11	N	501	HEM	CHD-C4C-NC	2.84	127.55	124.45
15	N	507	DCQ	C1M-C1-C6	-2.83	119.80	124.45
19	O	401	HEC	CMA-C3A-C4A	2.82	129.70	124.73
16	J	101	LMT	C3-C2-C1	-2.81	101.25	113.47
18	L	3003	XP4	O5-C4-O6	-2.81	116.61	123.63
14	L	3002	CDL	OA6-CA4-CA6	2.78	118.32	108.34
11	C	502	HEM	C1A-CHA-C4D	-2.76	119.76	126.25
11	N	502	HEM	C4C-NC-C1C	2.75	110.30	105.82
19	O	401	HEC	C3A-C4A-NA	2.74	114.70	109.64
11	C	502	HEM	O2D-CGD-CBD	2.73	122.62	114.00
11	N	502	HEM	CHC-C4B-C3B	-2.71	119.66	125.07
11	N	501	HEM	C1A-CHA-C4D	-2.71	119.87	126.25
16	C	507	LMT	C1B-O5B-C5B	2.71	119.01	113.72
19	D	401	HEC	C3A-C4A-NA	2.71	114.64	109.64
11	C	501	HEM	CHA-C4D-C3D	-2.71	120.24	125.23
14	A	3002	CDL	OA2-PA1-OA3	2.70	119.64	108.94
16	C	507	LMT	C3B-C4B-C5B	2.70	115.12	110.23
19	O	401	HEC	C2D-C1D-ND	2.69	114.45	110.14
11	N	502	HEM	CHA-C4D-C3D	-2.68	120.28	125.23
11	C	502	HEM	O1D-CGD-CBD	-2.67	114.61	123.09
16	N	508	LMT	C3B-C4B-C5B	2.67	115.07	110.23
15	N	507	DCQ	C8-C7-C6	2.67	120.76	113.10
19	O	401	HEC	CBB-CAB-C3B	-2.64	122.15	127.43
14	H	701	CDL	OB6-CB4-CB6	2.59	117.63	108.34
11	C	501	HEM	CHC-C1C-NC	2.58	127.26	124.45
19	D	401	HEC	CHC-C4B-C3B	-2.58	120.86	125.21
11	C	501	HEM	CAD-CBD-CGD	-2.57	106.84	113.67
16	N	508	LMT	O1'-C1'-C2'	-2.57	104.37	108.27
11	N	501	HEM	C4C-CHD-C1D	-2.57	120.56	126.02
11	N	502	HEM	C1D-C2D-C3D	-2.56	104.28	106.98
19	O	401	HEC	C4A-C3A-C2A	-2.56	103.18	106.97
11	N	501	HEM	CAA-C2A-C1A	2.56	129.93	124.94
14	A	3001	CDL	OB8-CB6-CB4	2.55	117.86	105.85
16	N	508	LMT	O5B-C5B-C4B	2.55	114.29	109.70
18	L	3003	XP4	O5-C4-C5	2.55	119.60	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	C	502	HEM	C1C-CHC-C4B	-2.54	120.62	126.02
11	N	501	HEM	CAD-CBD-CGD	-2.52	106.99	113.67
19	D	401	HEC	CAD-C3D-C4D	2.51	129.85	124.94
16	C	507	LMT	O2'-C2'-C1'	-2.51	104.09	110.08
11	N	501	HEM	C3B-C2B-C1B	-2.51	104.53	106.41
11	C	502	HEM	CHD-C4C-C3C	-2.50	120.99	125.21
11	C	501	HEM	C4C-CHD-C1D	-2.49	120.72	126.02
11	N	502	HEM	CHA-C1A-NA	2.49	128.38	123.86
19	O	401	HEC	CMC-C2C-C1C	2.49	129.21	125.42
19	D	401	HEC	CHD-C4C-C3C	-2.49	121.01	125.21
14	S	101	CDL	OA6-CA4-CA6	2.48	117.26	108.34
16	C	507	LMT	C4B-C3B-C2B	2.48	115.19	110.83
11	C	502	HEM	C4A-C3A-C2A	-2.48	103.97	106.82
11	N	502	HEM	CHD-C4C-C3C	-2.48	121.03	125.21
19	D	401	HEC	CBB-CAB-C3B	-2.47	122.50	127.43
14	L	3002	CDL	OA2-PA1-OA3	2.46	118.69	108.94
12	N	503	PC1	O21-C21-O22	-2.45	117.97	123.70
14	H	702	CDL	OA5-PA1-OA3	-2.44	99.27	108.94
19	D	401	HEC	CMC-C2C-C1C	2.43	129.12	125.42
16	C	507	LMT	C1'-O5'-C5'	2.42	118.45	113.72
11	C	502	HEM	CBA-CAA-C2A	-2.41	105.86	112.53
18	A	3003	XP4	O5-C4-C5	2.41	119.19	111.83
11	N	501	HEM	CHA-C1A-NA	2.41	128.23	123.86
11	N	502	HEM	C3B-C4B-NB	-2.39	107.75	109.47
19	O	401	HEC	CHC-C1C-C2C	-2.39	120.48	127.43
19	D	401	HEC	C4A-C3A-C2A	-2.39	103.43	106.97
19	O	401	HEC	CHB-C4A-NA	-2.37	121.87	124.45
14	C	505	CDL	CB4-OB6-CB5	2.37	123.47	117.80
11	N	502	HEM	O2D-CGD-O1D	-2.36	117.26	123.33
15	N	507	DCQ	O5-C5-C4	-2.35	116.05	121.03
11	C	501	HEM	CHD-C4C-NC	2.35	127.01	124.45
11	C	501	HEM	CHB-C1B-NB	2.34	127.27	124.37
18	L	3003	XP4	O4-P1-O2	-2.33	100.13	106.44
14	C	505	CDL	OB6-CB4-CB3	2.33	116.72	108.34
11	C	502	HEM	C2A-C1A-NA	-2.32	107.58	110.15
11	N	502	HEM	CHC-C1C-NC	2.31	126.97	124.45
19	D	401	HEC	CAD-CBD-CGD	-2.30	107.57	113.67
14	A	3002	CDL	OA5-PA1-OA3	-2.29	99.87	108.94
19	D	401	HEC	CMA-C3A-C4A	2.28	128.74	124.73
13	N	505	PTY	O14-P1-O13	-2.28	99.92	108.94
11	C	501	HEM	CAC-C3C-C2C	-2.27	121.03	128.43
11	N	502	HEM	C4A-C3A-C2A	-2.26	104.22	106.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	C	502	HEM	C1A-C2A-C3A	2.25	110.35	106.87
11	C	502	HEM	CHA-C1A-NA	2.25	127.94	123.86
14	L	3001	CDL	OB2-PB2-OB3	2.24	117.80	108.94
11	N	501	HEM	CHB-C4A-NA	2.22	127.89	123.86
11	C	501	HEM	C3B-C4B-NB	-2.22	107.87	109.47
14	N	504	CDL	OA6-CA4-CA3	2.22	116.30	108.34
16	N	508	LMT	O4'-C4B-C3B	-2.21	105.16	110.38
19	D	401	HEC	CBC-CAC-C3C	-2.21	123.02	127.43
19	D	401	HEC	CHD-C1D-C2D	-2.20	121.05	127.43
11	N	502	HEM	CBA-CAA-C2A	-2.19	106.46	112.53
12	N	503	PC1	C23-C22-C21	2.18	121.69	113.69
19	O	401	HEC	CHD-C1D-C2D	-2.18	121.10	127.43
16	U	101	LMT	C4B-C3B-C2B	2.18	114.65	110.83
12	C	503	PC1	O21-C21-O22	-2.17	118.62	123.70
11	N	501	HEM	O2D-CGD-O1D	-2.16	117.77	123.33
14	N	506	CDL	CB4-OB6-CB5	2.16	122.96	117.80
18	L	3003	XP4	O5-C3-C2	-2.13	102.25	108.40
18	L	3003	XP4	O7-C2-C3	-2.12	100.72	108.34
11	N	501	HEM	CAC-C3C-C2C	-2.12	121.54	128.43
16	J	101	LMT	O1'-C1'-C2'	2.11	111.48	108.27
16	N	508	LMT	C4B-C3B-C2B	2.11	114.54	110.83
13	N	505	PTY	O7-C8-C11	2.11	116.04	111.48
14	N	506	CDL	OB6-CB4-CB3	2.10	115.88	108.34
19	O	401	HEC	CAD-CBD-CGD	-2.10	108.10	113.67
16	U	101	LMT	C1-O1'-C1'	-2.10	110.10	113.68
18	A	3003	XP4	O4-P1-O2	-2.10	100.78	106.44
11	C	501	HEM	O2D-CGD-O1D	-2.09	117.96	123.33
14	H	701	CDL	OB4-PB2-OB3	2.09	122.16	112.44
19	D	401	HEC	CHA-C1A-NA	-2.09	122.18	124.45
14	S	101	CDL	OA4-PA1-OA5	2.08	116.98	107.57
19	O	401	HEC	CMC-C2C-C3C	2.08	131.43	126.55
16	U	101	LMT	O6'-C6'-C5'	-2.08	104.27	111.33
12	I	201	PC1	O12-P-O14	2.07	122.08	112.44
12	C	503	PC1	C23-C22-C21	2.07	121.28	113.69
19	D	401	HEC	CHC-C1C-C2C	-2.06	121.45	127.43
16	U	101	LMT	C4-C3-C2	-2.06	103.96	114.37
12	N	503	PC1	O21-C21-C22	-2.05	107.04	111.48
11	N	502	HEM	C4C-C3C-C2C	2.05	108.59	106.81
14	S	101	CDL	CA4-OA6-CA5	2.05	122.69	117.80
14	S	101	CDL	OA2-PA1-OA3	2.05	117.04	108.94
13	P	302	PTY	O12-P1-O14	-2.03	98.36	107.57
19	O	401	HEC	CHD-C4C-C3C	-2.02	121.80	125.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
18	A	3003	XP4	O7-C2-C3	-2.01	101.11	108.34
14	S	101	CDL	OB6-CB5-C51	2.01	114.67	111.09

There are no chirality outliers.

All (504) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
12	I	201	PC1	C11-O13-P-O12
12	I	201	PC1	C11-O13-P-O11
12	T	201	PC1	C11-O13-P-O12
12	T	201	PC1	C11-O13-P-O14
12	T	201	PC1	C11-O13-P-O11
12	T	201	PC1	O21-C2-C3-O31
13	C	504	PTY	N1-C2-C3-O11
13	C	504	PTY	C11-C8-O7-C6
13	C	504	PTY	C3-O11-P1-O14
13	C	504	PTY	C5-O14-P1-O11
13	C	504	PTY	C5-O14-P1-O12
13	P	302	PTY	C5-O14-P1-O11
13	P	302	PTY	C5-O14-P1-O12
13	P	302	PTY	C5-O14-P1-O13
13	N	505	PTY	C11-C8-O7-C6
13	N	505	PTY	C3-O11-P1-O12
13	N	505	PTY	C3-O11-P1-O13
13	N	505	PTY	C5-O14-P1-O11
13	N	505	PTY	C5-O14-P1-O13
13	E	401	PTY	N1-C2-C3-O11
13	E	401	PTY	C3-O11-P1-O13
13	E	401	PTY	C5-O14-P1-O11
13	E	401	PTY	C5-O14-P1-O12
14	C	505	CDL	O1-C1-CA2-OA2
14	C	505	CDL	CA2-C1-CB2-OB2
14	C	505	CDL	CA3-OA5-PA1-OA3
14	C	505	CDL	C11-CA5-OA6-CA4
14	C	505	CDL	CB2-OB2-PB2-OB3
14	C	505	CDL	CB2-OB2-PB2-OB4
14	C	505	CDL	CB2-OB2-PB2-OB5
14	C	505	CDL	CB3-OB5-PB2-OB2
14	C	505	CDL	CB3-OB5-PB2-OB3
14	A	3001	CDL	OA5-CA3-CA4-OA6
14	A	3001	CDL	OB5-CB3-CB4-CB6
14	A	3001	CDL	CB3-CB4-CB6-OB8

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Mol	Chain	Res	Type	Atoms
14	A	3001	CDL	OB6-CB4-CB6-OB8
14	A	3002	CDL	CA2-C1-CB2-OB2
14	A	3002	CDL	CA2-OA2-PA1-OA3
14	A	3002	CDL	CA2-OA2-PA1-OA4
14	A	3002	CDL	CA2-OA2-PA1-OA5
14	A	3002	CDL	CB2-OB2-PB2-OB3
14	A	3002	CDL	CB3-OB5-PB2-OB3
14	A	3002	CDL	OB5-CB3-CB4-OB6
14	H	701	CDL	CB3-OB5-PB2-OB3
14	H	702	CDL	CA2-OA2-PA1-OA4
14	H	702	CDL	CA2-OA2-PA1-OA5
14	H	702	CDL	C31-CA7-OA8-CA6
14	H	702	CDL	CB2-OB2-PB2-OB3
14	N	504	CDL	CA2-OA2-PA1-OA3
14	N	504	CDL	CA2-OA2-PA1-OA4
14	N	504	CDL	CA2-OA2-PA1-OA5
14	N	506	CDL	O1-C1-CA2-OA2
14	N	506	CDL	CA3-OA5-PA1-OA3
14	N	506	CDL	C11-CA5-OA6-CA4
14	N	506	CDL	CB2-OB2-PB2-OB3
14	N	506	CDL	CB2-OB2-PB2-OB4
14	N	506	CDL	CB2-OB2-PB2-OB5
14	N	506	CDL	CB3-OB5-PB2-OB2
14	N	506	CDL	CB3-OB5-PB2-OB3
14	L	3001	CDL	CB2-OB2-PB2-OB4
14	L	3001	CDL	CB2-OB2-PB2-OB5
14	L	3001	CDL	CB3-CB4-CB6-OB8
14	L	3002	CDL	CA2-C1-CB2-OB2
14	L	3002	CDL	CA2-OA2-PA1-OA3
14	L	3002	CDL	CA2-OA2-PA1-OA4
14	L	3002	CDL	CA2-OA2-PA1-OA5
14	L	3002	CDL	CB3-OB5-PB2-OB2
14	L	3002	CDL	CB3-OB5-PB2-OB3
14	L	3002	CDL	OB5-CB3-CB4-OB6
14	S	101	CDL	CA2-OA2-PA1-OA4
14	S	101	CDL	CA2-OA2-PA1-OA5
14	S	101	CDL	C31-CA7-OA8-CA6
14	S	101	CDL	CB2-OB2-PB2-OB3
14	S	101	CDL	CB3-OB5-PB2-OB4
15	C	506	DCQ	C5-C6-C7-C8
15	C	506	DCQ	C1-C6-C7-C8
16	U	101	LMT	O5'-C1'-O1'-C1

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Mol	Chain	Res	Type	Atoms
14	H	702	CDL	C51-CB5-OB6-CB4
14	S	101	CDL	C51-CB5-OB6-CB4
14	H	702	CDL	OA9-CA7-OA8-CA6
14	H	702	CDL	OB9-CB7-OB8-CB6
14	S	101	CDL	OA9-CA7-OA8-CA6
14	S	101	CDL	OB9-CB7-OB8-CB6
14	A	3001	CDL	CB4-CB6-OB8-CB7
12	T	201	PC1	O32-C31-O31-C3
14	A	3001	CDL	OA9-CA7-OA8-CA6
14	H	701	CDL	OB9-CB7-OB8-CB6
14	N	504	CDL	OB9-CB7-OB8-CB6
14	H	702	CDL	OB7-CB5-OB6-CB4
14	S	101	CDL	OB7-CB5-OB6-CB4
13	C	504	PTY	O10-C8-O7-C6
13	N	505	PTY	O10-C8-O7-C6
14	C	505	CDL	OA7-CA5-OA6-CA4
14	L	3002	CDL	OA7-CA5-OA6-CA4
15	C	506	DCQ	C6-C7-C8-C9
12	T	201	PC1	C32-C31-O31-C3
14	H	701	CDL	C71-CB7-OB8-CB6
14	H	702	CDL	C71-CB7-OB8-CB6
14	N	504	CDL	C71-CB7-OB8-CB6
14	S	101	CDL	C71-CB7-OB8-CB6
14	A	3001	CDL	C31-CA7-OA8-CA6
14	H	701	CDL	OA9-CA7-OA8-CA6
14	H	701	CDL	C31-CA7-OA8-CA6
14	L	3001	CDL	OA9-CA7-OA8-CA6
14	N	506	CDL	OA7-CA5-OA6-CA4
14	N	504	CDL	OA9-CA7-OA8-CA6
14	A	3002	CDL	O1-C1-CB2-OB2
14	L	3001	CDL	OA5-CA3-CA4-OA6
14	L	3001	CDL	OB5-CB3-CB4-OB6
14	L	3002	CDL	O1-C1-CB2-OB2
14	S	101	CDL	O1-C1-CA2-OA2
14	L	3001	CDL	C31-CA7-OA8-CA6
16	C	507	LMT	O5'-C5'-C6'-O6'
14	A	3002	CDL	C11-CA5-OA6-CA4
14	N	504	CDL	C51-CB5-OB6-CB4
14	L	3002	CDL	C11-CA5-OA6-CA4
14	L	3001	CDL	OB9-CB7-OB8-CB6
16	N	508	LMT	O5'-C5'-C6'-O6'
16	U	101	LMT	C11-C10-C9-C8

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Mol	Chain	Res	Type	Atoms
14	L	3001	CDL	C71-CB7-OB8-CB6
13	N	505	PTY	C31-C32-C33-C34
16	N	508	LMT	O5B-C5B-C6B-O6B
12	I	201	PC1	C32-C33-C34-C35
13	P	302	PTY	C31-C30-O4-C1
16	C	507	LMT	O5B-C5B-C6B-O6B
13	P	302	PTY	C33-C34-C35-C36
13	N	505	PTY	C33-C34-C35-C36
14	N	504	CDL	OB7-CB5-OB6-CB4
14	C	505	CDL	CB2-C1-CA2-OA2
14	N	504	CDL	CA2-C1-CB2-OB2
14	N	506	CDL	CA2-C1-CB2-OB2
14	L	3001	CDL	CA2-C1-CB2-OB2
14	L	3001	CDL	OA5-CA3-CA4-CA6
14	L	3001	CDL	OB5-CB3-CB4-CB6
12	C	503	PC1	C32-C31-O31-C3
13	E	401	PTY	C13-C14-C15-C16
13	P	302	PTY	C37-C38-C39-C40
16	J	101	LMT	C4-C5-C6-C7
16	N	508	LMT	C4'-C5'-C6'-O6'
13	E	401	PTY	C30-C31-C32-C33
14	A	3001	CDL	CB7-C71-C72-C73
14	H	701	CDL	CA5-C11-C12-C13
14	H	702	CDL	O1-C1-CA2-OA2
14	N	506	CDL	O1-C1-CB2-OB2
16	J	101	LMT	O5B-C5B-C6B-O6B
13	C	504	PTY	C31-C32-C33-C34
13	N	505	PTY	C37-C38-C39-C40
16	J	101	LMT	C2-C3-C4-C5
16	C	507	LMT	C4'-C5'-C6'-O6'
13	N	505	PTY	C8-C11-C12-C13
14	N	504	CDL	C31-CA7-OA8-CA6
14	N	504	CDL	CA5-C11-C12-C13
13	C	504	PTY	C35-C36-C37-C38
12	N	503	PC1	C32-C31-O31-C3
14	A	3002	CDL	OA7-CA5-OA6-CA4
12	T	201	PC1	C21-C22-C23-C24
12	T	201	PC1	C31-C32-C33-C34
14	H	701	CDL	CB7-C71-C72-C73
14	N	506	CDL	CA7-C31-C32-C33
14	L	3002	CDL	CA7-C31-C32-C33
12	C	503	PC1	O32-C31-O31-C3

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Mol	Chain	Res	Type	Atoms
13	P	302	PTY	C30-C31-C32-C33
14	C	505	CDL	CA7-C31-C32-C33
13	E	401	PTY	C40-C41-C42-C43
13	E	401	PTY	C35-C36-C37-C38
14	C	505	CDL	O1-C1-CB2-OB2
14	H	702	CDL	O1-C1-CB2-OB2
14	N	504	CDL	O1-C1-CB2-OB2
14	L	3001	CDL	O1-C1-CB2-OB2
14	S	101	CDL	O1-C1-CB2-OB2
13	P	302	PTY	O30-C30-O4-C1
13	E	401	PTY	C8-C11-C12-C13
14	A	3002	CDL	CA7-C31-C32-C33
14	L	3001	CDL	OB6-CB4-CB6-OB8
14	N	504	CDL	CB7-C71-C72-C73
13	P	302	PTY	C35-C36-C37-C38
12	I	201	PC1	C21-C22-C23-C24
14	H	701	CDL	CB5-C51-C52-C53
14	A	3001	CDL	OA5-CA3-CA4-CA6
14	H	702	CDL	CB2-C1-CA2-OA2
14	H	702	CDL	CA2-C1-CB2-OB2
14	S	101	CDL	CB2-C1-CA2-OA2
14	S	101	CDL	CA2-C1-CB2-OB2
13	E	401	PTY	C31-C30-O4-C1
14	N	504	CDL	O1-C1-CA2-OA2
12	I	201	PC1	C32-C31-O31-C3
12	N	503	PC1	O32-C31-O31-C3
13	E	401	PTY	O14-C5-C6-O7
13	E	401	PTY	C33-C34-C35-C36
14	A	3002	CDL	C51-CB5-OB6-CB4
14	A	3002	CDL	C52-C53-C54-C55
16	C	507	LMT	C4-C5-C6-C7
14	L	3001	CDL	CB7-C71-C72-C73
12	I	201	PC1	O32-C31-O31-C3
16	N	508	LMT	C7-C8-C9-C10
13	E	401	PTY	C16-C17-C18-C19
14	H	701	CDL	C11-C12-C13-C14
12	I	201	PC1	C31-C32-C33-C34
13	C	504	PTY	C8-C11-C12-C13
16	U	101	LMT	C4-C5-C6-C7
13	N	505	PTY	C16-C17-C18-C19
14	H	701	CDL	C51-CB5-OB6-CB4
14	N	506	CDL	C51-CB5-OB6-CB4

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Mol	Chain	Res	Type	Atoms
13	E	401	PTY	C38-C39-C40-C41
14	L	3002	CDL	O1-C1-CA2-OA2
16	N	508	LMT	C4B-C5B-C6B-O6B
13	C	504	PTY	C37-C38-C39-C40
13	N	505	PTY	C32-C33-C34-C35
16	C	507	LMT	C6-C7-C8-C9
15	N	507	DCQ	C11-C10-C9-C8
16	C	507	LMT	C5-C6-C7-C8
14	L	3002	CDL	C11-C12-C13-C14
14	N	504	CDL	C74-C75-C76-C77
13	C	504	PTY	C14-C15-C16-C17
13	P	302	PTY	C16-C17-C18-C19
14	A	3001	CDL	C71-C72-C73-C74
15	N	507	DCQ	C9-C10-C11-C12
14	C	505	CDL	C51-CB5-OB6-CB4
16	N	508	LMT	C1-C2-C3-C4
13	E	401	PTY	O30-C30-O4-C1
13	C	504	PTY	C13-C14-C15-C16
13	P	302	PTY	C40-C41-C42-C43
14	L	3002	CDL	C12-C13-C14-C15
14	N	506	CDL	OB7-CB5-OB6-CB4
15	C	506	DCQ	C11-C10-C9-C8
16	C	507	LMT	C11-C10-C9-C8
15	N	507	DCQ	C6-C7-C8-C9
13	E	401	PTY	C11-C12-C13-C14
14	H	701	CDL	C71-C72-C73-C74
13	P	302	PTY	C34-C35-C36-C37
16	U	101	LMT	C3-C4-C5-C6
14	H	701	CDL	C73-C74-C75-C76
14	H	701	CDL	C74-C75-C76-C77
13	P	302	PTY	C41-C42-C43-C44
14	A	3002	CDL	C12-C13-C14-C15
15	C	506	DCQ	C10-C11-C12-C13
14	C	505	CDL	OB7-CB5-OB6-CB4
13	C	504	PTY	C11-C12-C13-C14
13	C	504	PTY	C38-C39-C40-C41
16	J	101	LMT	C6-C7-C8-C9
16	C	507	LMT	C1-C2-C3-C4
13	N	505	PTY	C14-C15-C16-C17
14	N	504	CDL	C71-C72-C73-C74
16	C	507	LMT	C7-C8-C9-C10
16	C	507	LMT	C3-C4-C5-C6

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Mol	Chain	Res	Type	Atoms
16	N	508	LMT	C6-C7-C8-C9
12	N	503	PC1	C28-C29-C2A-C2B
13	P	302	PTY	C13-C14-C15-C16
12	C	503	PC1	O21-C2-C3-O31
12	N	503	PC1	O21-C2-C3-O31
12	C	503	PC1	C28-C29-C2A-C2B
13	C	504	PTY	C12-C13-C14-C15
13	E	401	PTY	C14-C15-C16-C17
16	J	101	LMT	C7-C8-C9-C10
14	N	506	CDL	CB2-C1-CA2-OA2
14	A	3002	CDL	C33-C34-C35-C36
13	N	505	PTY	C30-C31-C32-C33
14	A	3002	CDL	OB7-CB5-OB6-CB4
13	N	505	PTY	C11-C12-C13-C14
16	N	508	LMT	C5-C6-C7-C8
13	N	505	PTY	O14-C5-C6-C1
14	H	701	CDL	OB7-CB5-OB6-CB4
16	J	101	LMT	C4B-C5B-C6B-O6B
13	C	504	PTY	C40-C41-C42-C43
14	L	3002	CDL	CA5-C11-C12-C13
16	C	507	LMT	C4B-C5B-C6B-O6B
12	T	201	PC1	C33-C34-C35-C36
14	L	3002	CDL	C51-CB5-OB6-CB4
13	C	504	PTY	C33-C34-C35-C36
13	P	302	PTY	C11-C12-C13-C14
13	C	504	PTY	O4-C1-C6-C5
13	N	505	PTY	O4-C1-C6-C5
14	H	701	CDL	CB3-CB4-CB6-OB8
14	N	504	CDL	C11-C12-C13-C14
15	N	507	DCQ	C10-C11-C12-C13
13	P	302	PTY	C8-C11-C12-C13
13	E	401	PTY	C12-C13-C14-C15
13	P	302	PTY	C38-C39-C40-C41
14	A	3002	CDL	CA6-CA4-OA6-CA5
14	L	3002	CDL	CA6-CA4-OA6-CA5
14	A	3001	CDL	C73-C74-C75-C76
14	N	504	CDL	OB5-CB3-CB4-OB6
16	N	508	LMT	C3-C4-C5-C6
13	N	505	PTY	C34-C35-C36-C37
13	N	505	PTY	C36-C37-C38-C39
14	L	3001	CDL	C71-C72-C73-C74
14	A	3001	CDL	OB5-CB3-CB4-OB6

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Mol	Chain	Res	Type	Atoms
13	C	504	PTY	C41-C42-C43-C44
16	J	101	LMT	C9-C10-C11-C12
12	N	503	PC1	C26-C27-C28-C29
13	E	401	PTY	C41-C42-C43-C44
16	N	508	LMT	C9-C10-C11-C12
13	C	504	PTY	C17-C18-C19-C20
13	N	505	PTY	C41-C42-C43-C44
14	H	701	CDL	C75-C76-C77-C78
14	N	504	CDL	C75-C76-C77-C78
12	N	503	PC1	C27-C28-C29-C2A
13	C	504	PTY	C15-C16-C17-C18
16	J	101	LMT	O5'-C5'-C6'-O6'
14	A	3002	CDL	C13-C14-C15-C16
12	C	503	PC1	C22-C23-C24-C25
14	A	3002	CDL	OA5-CA3-CA4-CA6
14	H	701	CDL	OB5-CB3-CB4-CB6
14	H	702	CDL	OB5-CB3-CB4-CB6
14	L	3002	CDL	OA5-CA3-CA4-CA6
14	S	101	CDL	OB5-CB3-CB4-CB6
16	C	507	LMT	C9-C10-C11-C12
12	C	503	PC1	C26-C27-C28-C29
14	N	504	CDL	CB2-C1-CA2-OA2
14	L	3002	CDL	CB2-C1-CA2-OA2
14	H	702	CDL	CB3-CB4-CB6-OB8
14	N	504	CDL	CA3-CA4-CA6-OA8
12	N	503	PC1	C22-C23-C24-C25
13	N	505	PTY	C38-C39-C40-C41
14	L	3002	CDL	OB7-CB5-OB6-CB4
14	N	506	CDL	C31-C32-C33-C34
14	L	3002	CDL	C31-C32-C33-C34
14	N	504	CDL	C73-C74-C75-C76
13	N	505	PTY	O4-C1-C6-O7
14	H	701	CDL	OA6-CA4-CA6-OA8
14	H	702	CDL	OB6-CB4-CB6-OB8
14	L	3001	CDL	C32-C33-C34-C35
13	C	504	PTY	C39-C40-C41-C42
12	C	503	PC1	C27-C28-C29-C2A
16	N	508	LMT	C11-C10-C9-C8
16	U	101	LMT	C5-C6-C7-C8
14	L	3002	CDL	C34-C35-C36-C37
12	I	201	PC1	C22-C23-C24-C25
13	N	505	PTY	C15-C16-C17-C18

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Mol	Chain	Res	Type	Atoms
13	E	401	PTY	O14-C5-C6-C1
14	H	702	CDL	OA5-CA3-CA4-CA6
14	N	504	CDL	OA5-CA3-CA4-CA6
16	C	507	LMT	O1'-C1-C2-C3
16	J	101	LMT	O1'-C1-C2-C3
13	P	302	PTY	C14-C15-C16-C17
13	E	401	PTY	C31-C32-C33-C34
14	A	3001	CDL	C72-C73-C74-C75
14	N	504	CDL	C51-C52-C53-C54
16	N	508	LMT	C2-C3-C4-C5
16	U	101	LMT	C7-C8-C9-C10
14	H	702	CDL	CB6-CB4-OB6-CB5
13	N	505	PTY	O14-C5-C6-O7
14	C	505	CDL	OA5-CA3-CA4-OA6
14	H	701	CDL	OA5-CA3-CA4-OA6
12	C	503	PC1	C1-C2-C3-O31
12	N	503	PC1	C1-C2-C3-O31
12	T	201	PC1	C1-C2-C3-O31
14	S	101	CDL	CB3-CB4-CB6-OB8
13	E	401	PTY	C36-C37-C38-C39
14	H	701	CDL	OB6-CB4-CB6-OB8
14	S	101	CDL	OB6-CB4-CB6-OB8
13	N	505	PTY	C12-C13-C14-C15
13	E	401	PTY	C34-C35-C36-C37
14	C	505	CDL	C31-C32-C33-C34
14	S	101	CDL	OA5-CA3-CA4-CA6
14	L	3001	CDL	O1-C1-CA2-OA2
14	N	506	CDL	C11-C12-C13-C14
13	E	401	PTY	C32-C33-C34-C35
13	P	302	PTY	C31-C32-C33-C34
14	H	701	CDL	C51-C52-C53-C54
13	P	302	PTY	O14-C5-C6-O7
14	H	701	CDL	OB5-CB3-CB4-OB6
14	H	702	CDL	OA5-CA3-CA4-OA6
14	N	506	CDL	OA5-CA3-CA4-OA6
14	S	101	CDL	OB5-CB3-CB4-OB6
14	A	3002	CDL	CA5-C11-C12-C13
13	C	504	PTY	O4-C1-C6-O7
14	N	504	CDL	OA6-CA4-CA6-OA8
14	L	3001	CDL	CB2-C1-CA2-OA2
14	L	3002	CDL	C33-C34-C35-C36
14	H	701	CDL	CA3-CA4-CA6-OA8

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Mol	Chain	Res	Type	Atoms
14	N	504	CDL	CB3-CB4-CB6-OB8
14	N	506	CDL	C13-C14-C15-C16
12	I	201	PC1	C1-O11-P-O14
13	C	504	PTY	C3-O11-P1-O13
13	C	504	PTY	C5-O14-P1-O13
13	N	505	PTY	C3-O11-P1-O14
13	N	505	PTY	C5-O14-P1-O12
14	C	505	CDL	CA3-OA5-PA1-OA2
14	C	505	CDL	CA3-OA5-PA1-OA4
14	C	505	CDL	CB3-OB5-PB2-OB4
14	A	3002	CDL	CB3-OB5-PB2-OB2
14	A	3002	CDL	CB3-OB5-PB2-OB4
14	H	701	CDL	CB3-OB5-PB2-OB2
14	H	702	CDL	CB3-OB5-PB2-OB2
14	H	702	CDL	CB3-OB5-PB2-OB3
14	H	702	CDL	CB3-OB5-PB2-OB4
14	N	506	CDL	CA3-OA5-PA1-OA2
14	N	506	CDL	CA3-OA5-PA1-OA4
14	N	506	CDL	CB3-OB5-PB2-OB4
14	L	3002	CDL	CB2-OB2-PB2-OB3
14	L	3002	CDL	CB3-OB5-PB2-OB4
14	S	101	CDL	CB3-OB5-PB2-OB2
14	S	101	CDL	CB3-OB5-PB2-OB3
14	L	3002	CDL	C52-C53-C54-C55
14	C	505	CDL	CA4-CA3-OA5-PA1
14	N	506	CDL	CA4-CA3-OA5-PA1
15	C	506	DCQ	C7-C8-C9-C10
14	A	3002	CDL	C31-C32-C33-C34
14	N	504	CDL	CA3-CA4-OA6-CA5
14	N	504	CDL	CA6-CA4-OA6-CA5
14	S	101	CDL	CB6-CB4-OB6-CB5
14	H	701	CDL	OA5-CA3-CA4-CA6
13	C	504	PTY	C31-C30-O4-C1
13	C	504	PTY	O30-C30-O4-C1
13	C	504	PTY	C32-C33-C34-C35
14	A	3002	CDL	OA6-CA4-CA6-OA8
12	I	201	PC1	C33-C34-C35-C36
14	N	504	CDL	CB5-C51-C52-C53
14	H	702	CDL	C32-C31-CA7-OA8
19	O	401	HEC	CAA-CBA-CGA-O2A
14	A	3002	CDL	C34-C35-C36-C37
14	N	504	CDL	OB5-CB3-CB4-CB6

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Mol	Chain	Res	Type	Atoms
16	U	101	LMT	O5'-C5'-C6'-O6'
19	D	401	HEC	CAA-CBA-CGA-O2A
11	C	501	HEM	C3D-CAD-CBD-CGD
11	N	501	HEM	C3D-CAD-CBD-CGD
14	A	3002	CDL	CB2-C1-CA2-OA2
14	A	3002	CDL	O1-C1-CA2-OA2
14	A	3002	CDL	C51-C52-C53-C54
14	H	701	CDL	CA3-CA4-OA6-CA5
11	C	502	HEM	CAA-CBA-CGA-O1A
11	N	502	HEM	CAD-CBD-CGD-O2D
15	N	507	DCQ	C12-C13-C14-C15
11	N	502	HEM	CAA-CBA-CGA-O1A
19	O	401	HEC	CAA-CBA-CGA-O1A
14	N	504	CDL	OA5-CA3-CA4-OA6
11	C	502	HEM	CAD-CBD-CGD-O1D
11	N	501	HEM	CAA-CBA-CGA-O2A
13	P	302	PTY	O14-C5-C6-C1
11	C	501	HEM	CAA-CBA-CGA-O2A
19	D	401	HEC	CAA-CBA-CGA-O1A
11	N	502	HEM	CAD-CBD-CGD-O1D
11	C	501	HEM	CAA-CBA-CGA-O1A
11	N	501	HEM	CAA-CBA-CGA-O1A
11	C	502	HEM	CAD-CBD-CGD-O2D
16	U	101	LMT	O1'-C1-C2-C3
15	N	507	DCQ	C7-C8-C9-C10
11	N	501	HEM	CAD-CBD-CGD-O2D
14	C	505	CDL	C13-C14-C15-C16
11	N	502	HEM	CAA-CBA-CGA-O2A
14	L	3002	CDL	CB5-C51-C52-C53
14	L	3002	CDL	OA5-CA3-CA4-OA6
14	L	3002	CDL	C51-C52-C53-C54
12	I	201	PC1	C22-C21-O21-C2
11	C	502	HEM	CAA-CBA-CGA-O2A
13	C	504	PTY	O14-C5-C6-C1
14	N	506	CDL	OA5-CA3-CA4-CA6
14	S	101	CDL	C72-C71-CB7-OB8
11	C	501	HEM	CAD-CBD-CGD-O2D
18	L	3003	XP4	C5-C6-C7-C8
14	L	3001	CDL	CA4-CA3-OA5-PA1
15	C	506	DCQ	C9-C10-C11-C12
14	H	702	CDL	C32-C31-CA7-OA9
13	C	504	PTY	C6-C5-O14-P1

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Mol	Chain	Res	Type	Atoms
14	S	101	CDL	OA5-CA3-CA4-OA6
11	N	501	HEM	CAD-CBD-CGD-O1D
12	N	503	PC1	C32-C33-C34-C35
11	C	501	HEM	CAD-CBD-CGD-O1D
14	N	506	CDL	C32-C33-C34-C35
14	C	505	CDL	C11-C12-C13-C14
14	C	505	CDL	OA5-CA3-CA4-CA6
14	S	101	CDL	C32-C31-CA7-OA8
15	C	506	DCQ	C2-C3-O3-C3M
14	S	101	CDL	CB4-CB3-OB5-PB2
14	H	701	CDL	C72-C71-CB7-OB8
14	A	3001	CDL	C72-C71-CB7-OB8
12	I	201	PC1	O21-C21-C22-C23
14	H	702	CDL	OB5-CB3-CB4-OB6
16	C	507	LMT	C2-C3-C4-C5
15	N	507	DCQ	C13-C14-C15-C16
14	N	504	CDL	C72-C71-CB7-OB8
16	J	101	LMT	C2-C1-O1'-C1'
13	P	302	PTY	O4-C30-C31-C32
19	D	401	HEC	CAD-CBD-CGD-O1D
19	D	401	HEC	CAD-CBD-CGD-O2D
12	C	503	PC1	O31-C31-C32-C33
13	N	505	PTY	C31-C30-O4-C1
14	C	505	CDL	CB6-CB4-OB6-CB5
14	N	506	CDL	CB6-CB4-OB6-CB5
12	I	201	PC1	O22-C21-O21-C2
19	O	401	HEC	CAD-CBD-CGD-O1D
13	N	505	PTY	C13-C14-C15-C16
14	A	3002	CDL	CB5-C51-C52-C53
12	N	503	PC1	O31-C31-C32-C33
15	N	507	DCQ	C5-C4-O4-C4M
12	I	201	PC1	O22-C21-C22-C23
14	H	701	CDL	C72-C71-CB7-OB9
14	C	505	CDL	C32-C31-CA7-OA8
14	S	101	CDL	C32-C31-CA7-OA9
12	N	503	PC1	C21-C22-C23-C24
14	C	505	CDL	OB6-CB4-CB6-OB8
14	H	701	CDL	CA2-C1-CB2-OB2
16	U	101	LMT	C2-C3-C4-C5
12	T	201	PC1	C32-C33-C34-C35
14	N	506	CDL	C32-C31-CA7-OA8
12	C	503	PC1	O32-C31-C32-C33

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Mol	Chain	Res	Type	Atoms
16	U	101	LMT	O5B-C5B-C6B-O6B
14	L	3001	CDL	C72-C71-CB7-OB8
12	C	503	PC1	C23-C24-C25-C26
12	N	503	PC1	O32-C31-C32-C33
13	P	302	PTY	O30-C30-C31-C32
14	A	3002	CDL	C31-CA7-OA8-CA6
14	N	504	CDL	C72-C71-CB7-OB9

There are no ring outliers.

31 monomers are involved in 98 short contacts:

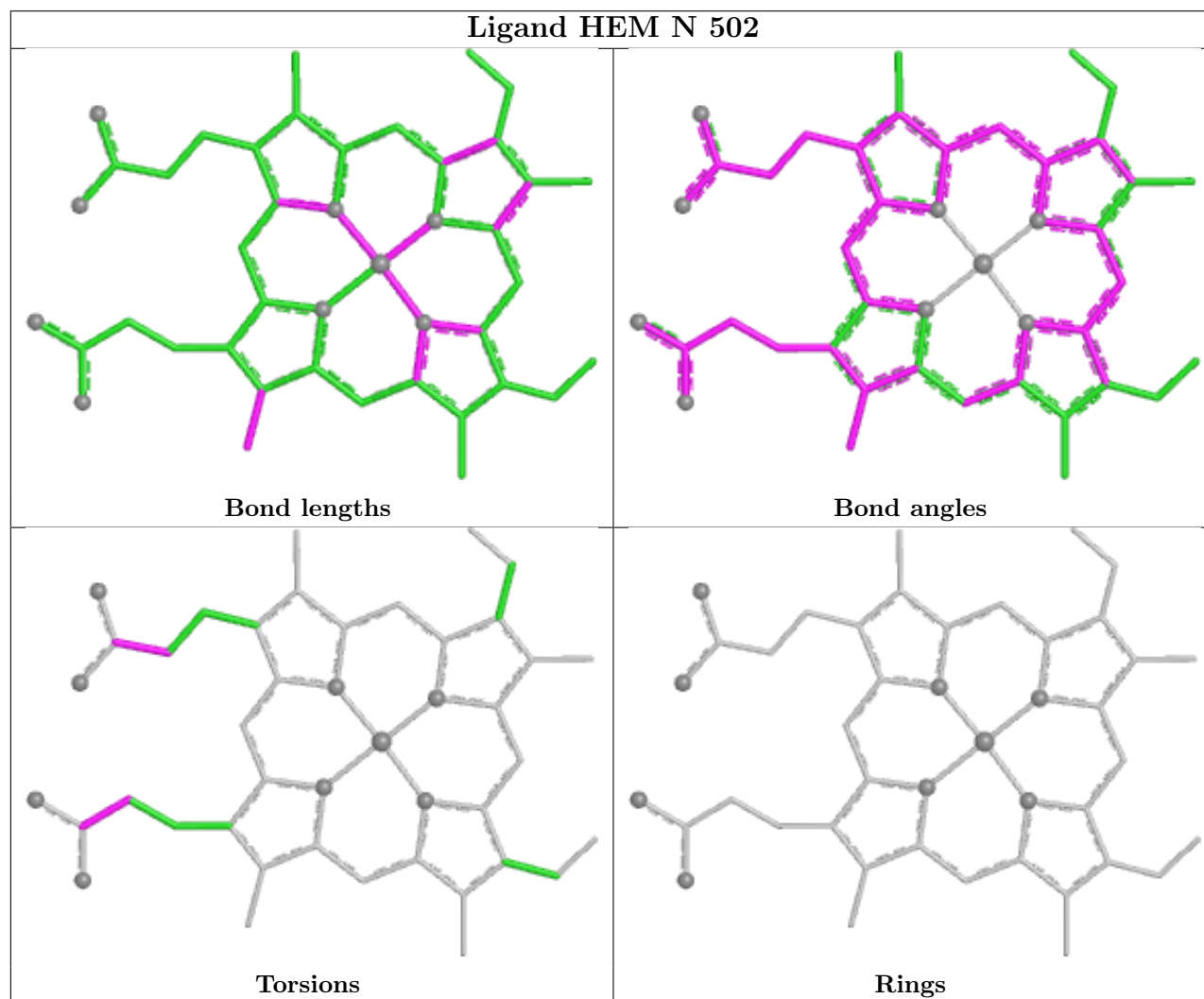
Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	N	502	HEM	3	0
19	D	401	HEC	11	0
18	L	3003	XP4	1	0
15	C	506	DCQ	4	0
16	C	507	LMT	1	0
16	U	101	LMT	1	0
14	N	506	CDL	2	0
16	J	101	LMT	1	0
14	H	701	CDL	4	0
13	N	505	PTY	9	0
14	S	101	CDL	4	0
14	H	702	CDL	4	0
11	N	501	HEM	4	0
12	N	503	PC1	2	0
14	A	3002	CDL	2	0
14	N	504	CDL	3	0
13	C	504	PTY	3	0
14	L	3001	CDL	1	0
14	L	3002	CDL	1	0
11	C	501	HEM	4	0
15	N	507	DCQ	3	0
11	C	502	HEM	4	0
12	I	201	PC1	1	0
12	T	201	PC1	4	0
14	A	3001	CDL	2	0
13	E	401	PTY	2	0
12	C	503	PC1	2	0
14	C	505	CDL	1	0
16	N	508	LMT	1	0
19	O	401	HEC	11	0

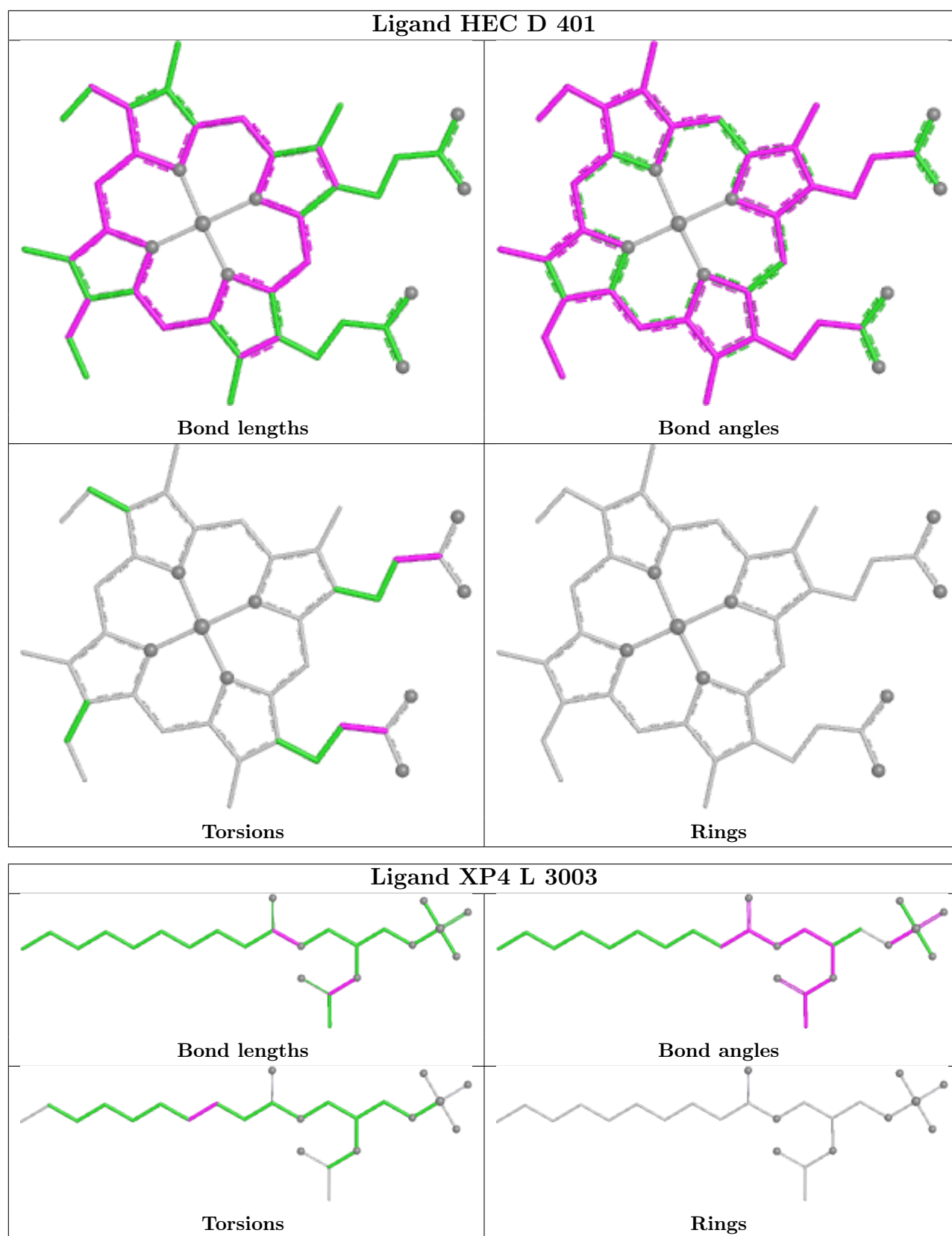
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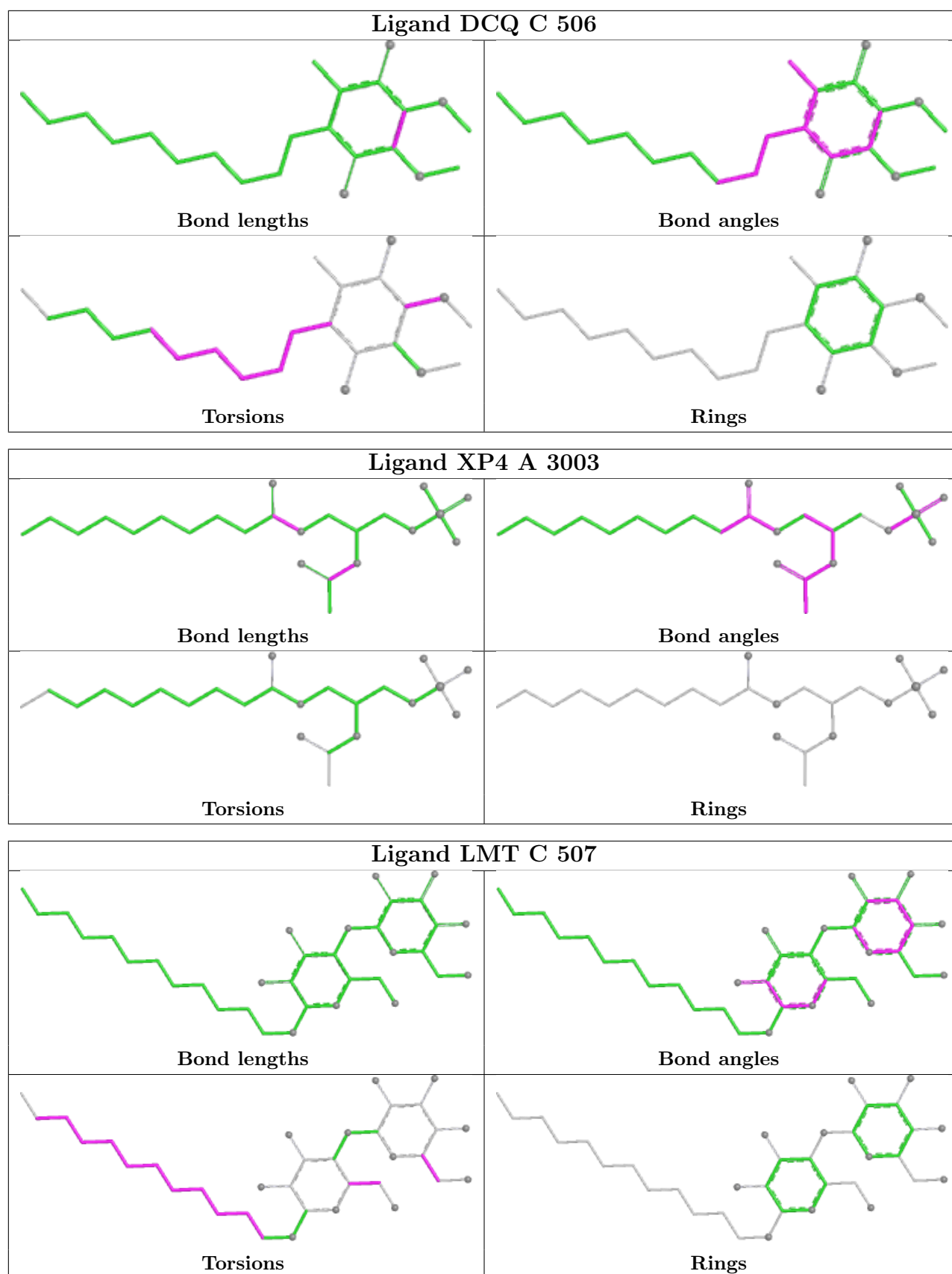
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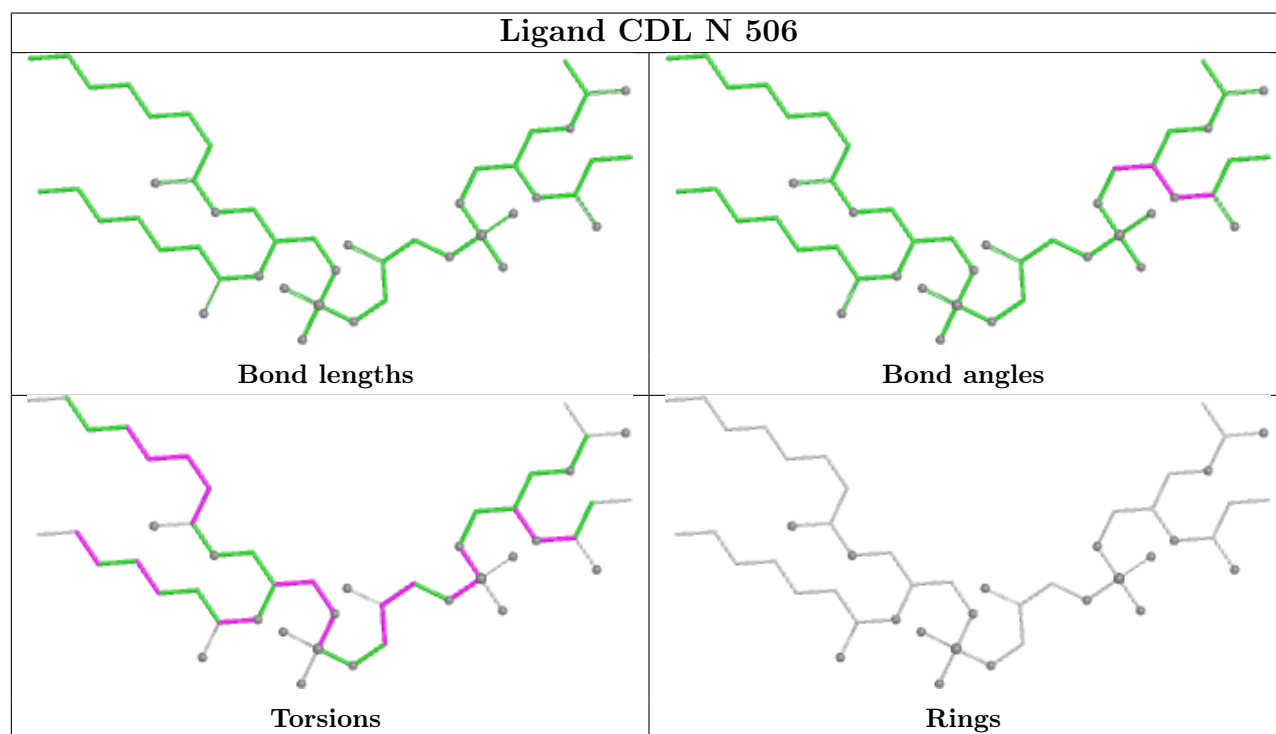
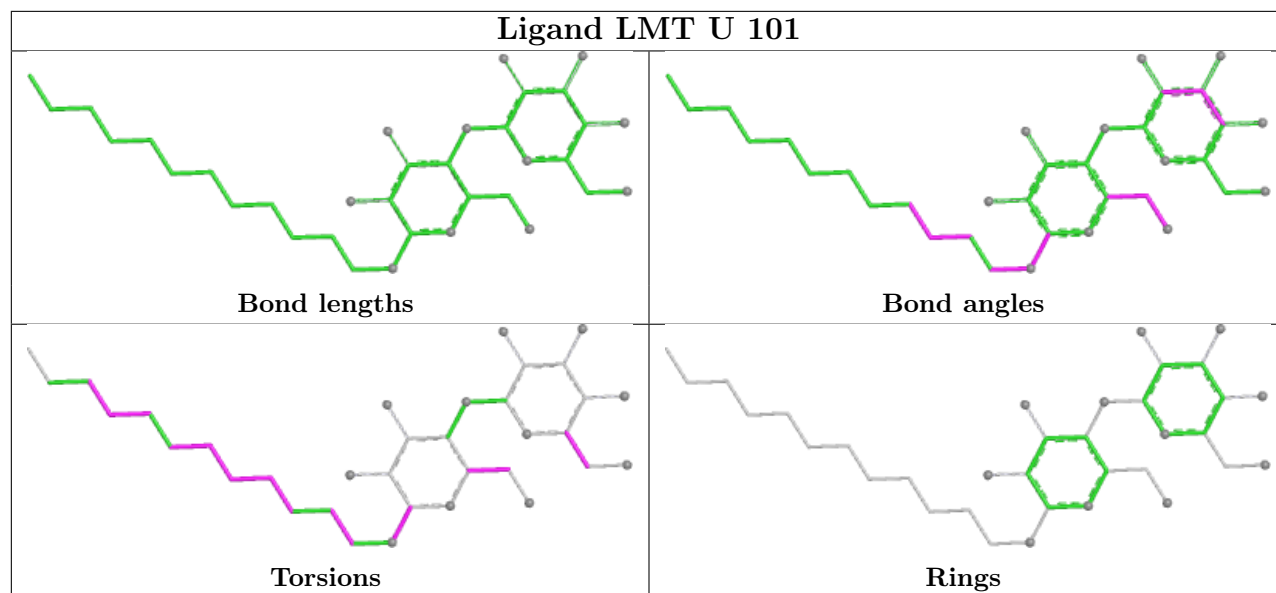
Mol	Chain	Res	Type	Clashes	Symm-Clashes
13	P	302	PTY	12	0

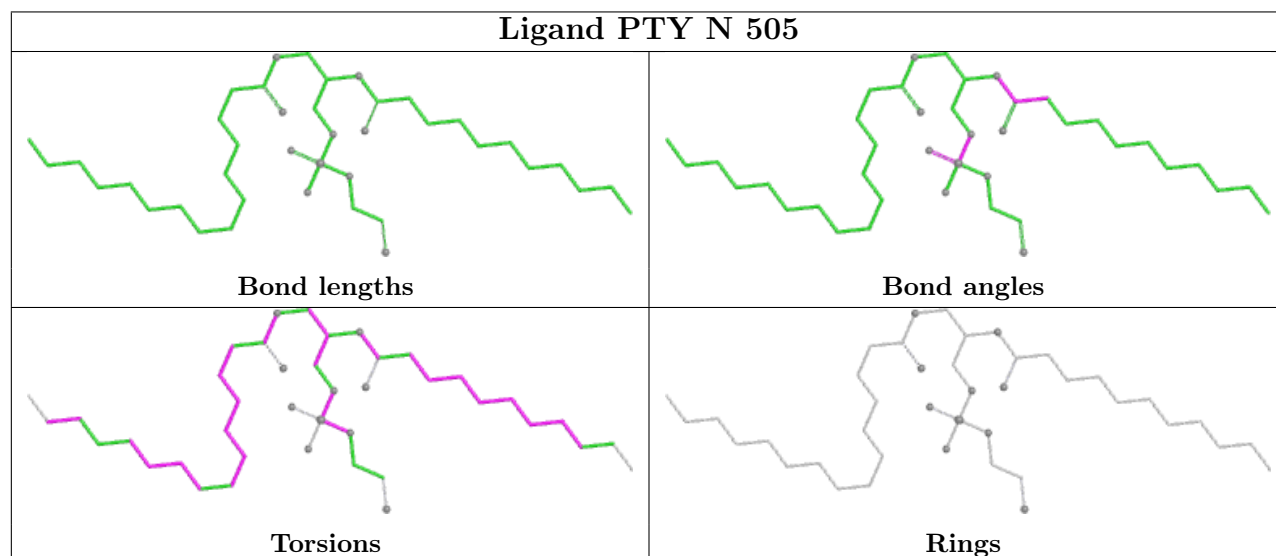
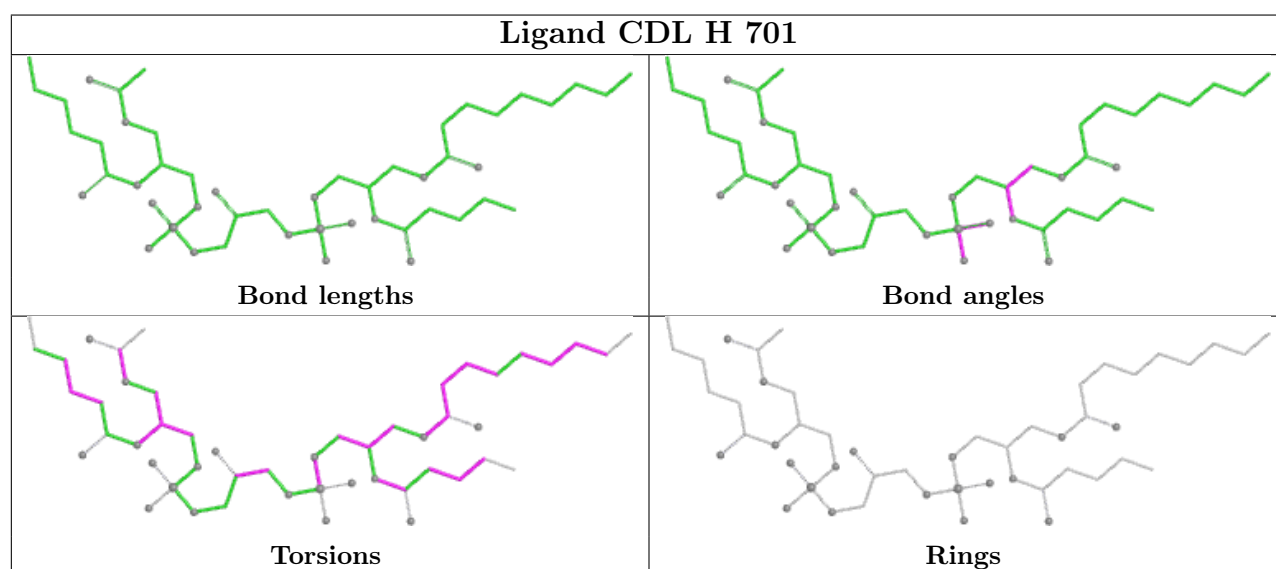
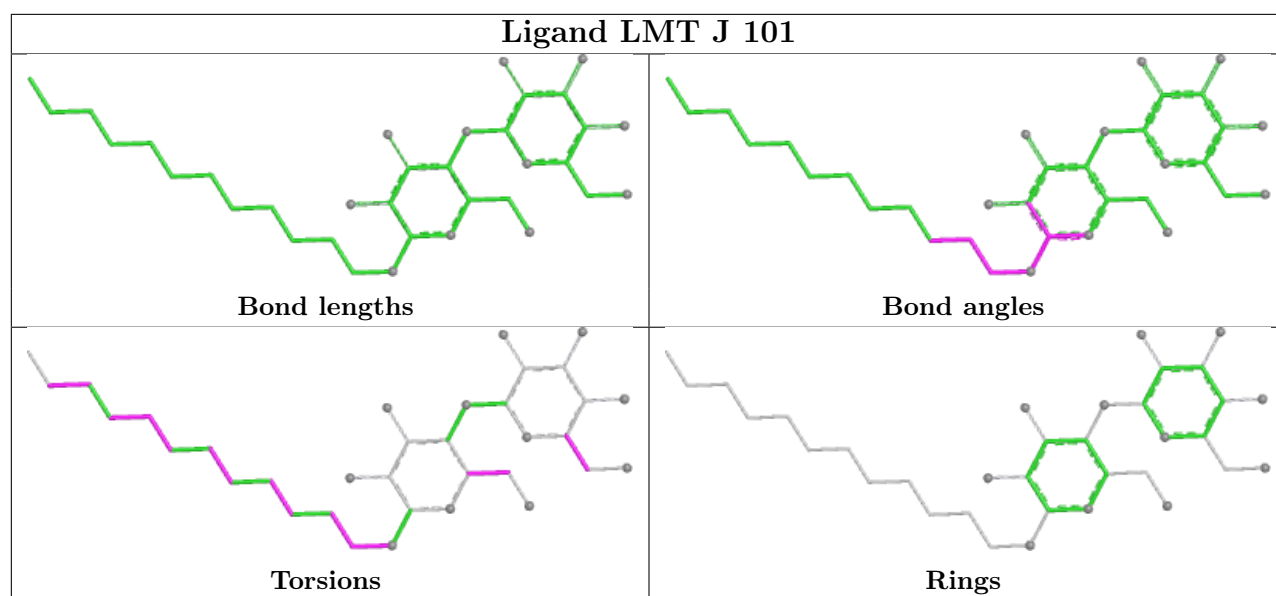
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

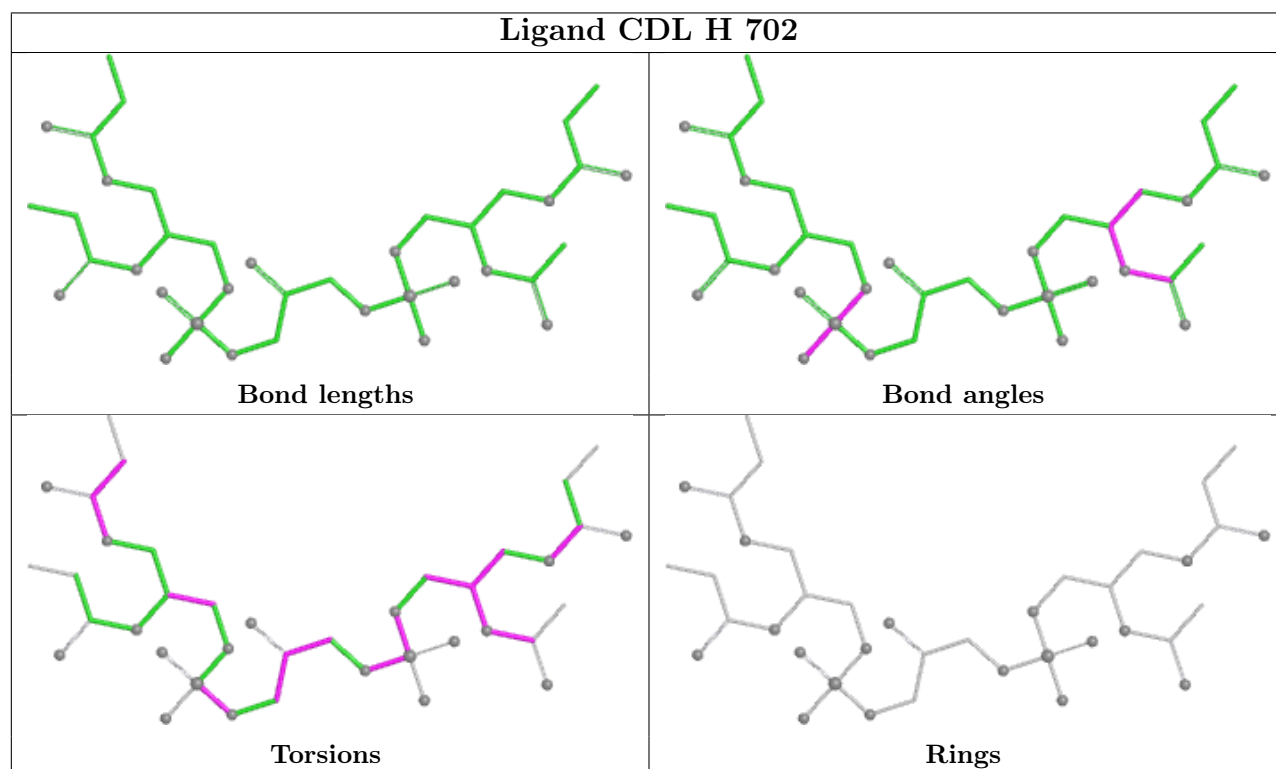
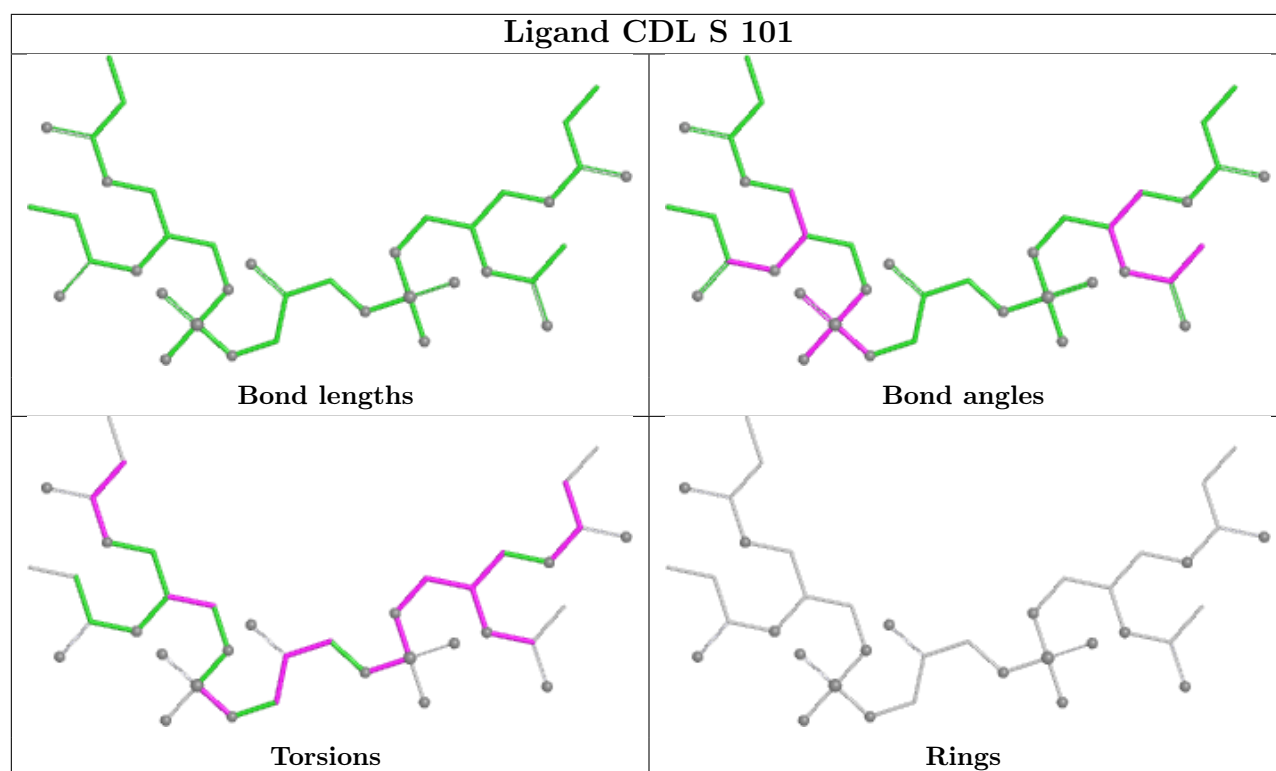


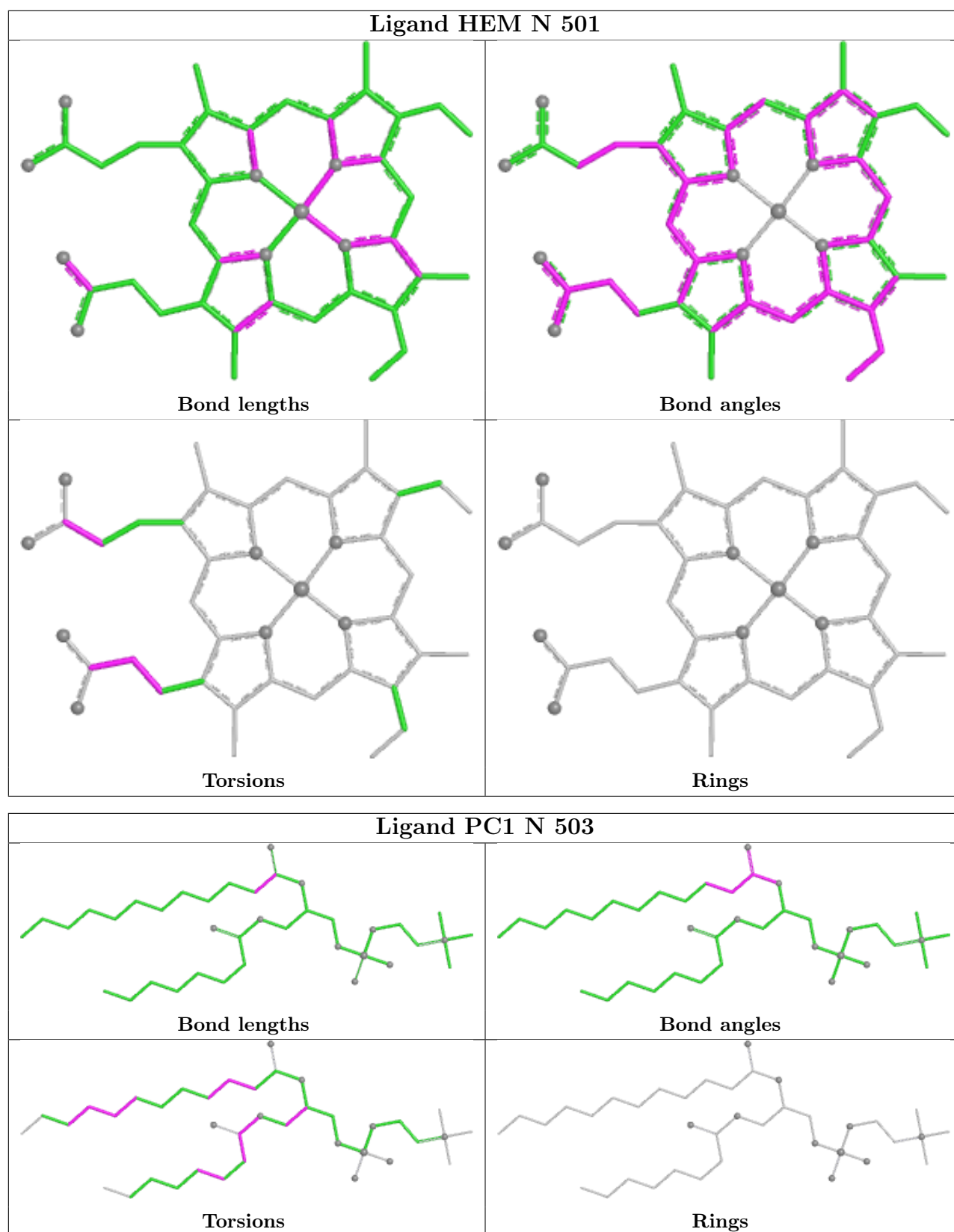


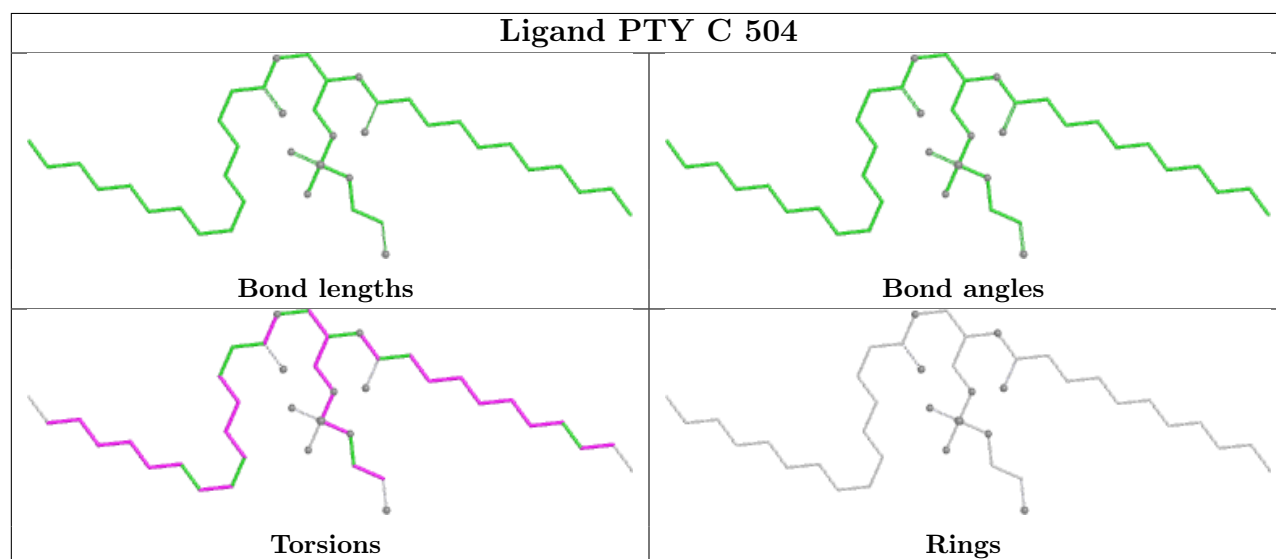
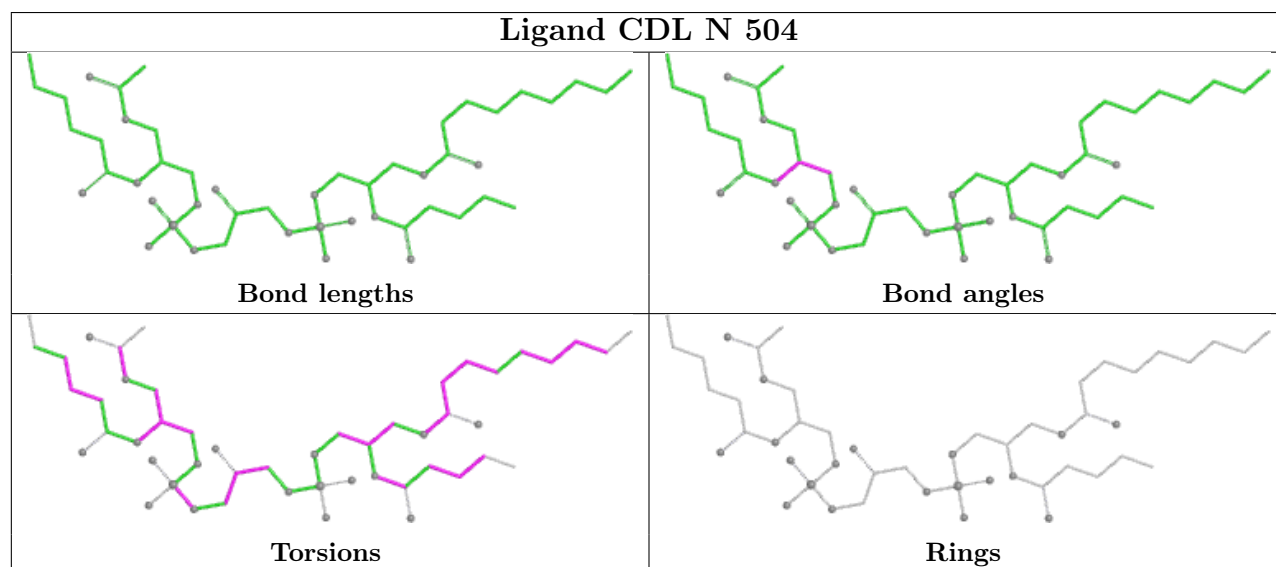
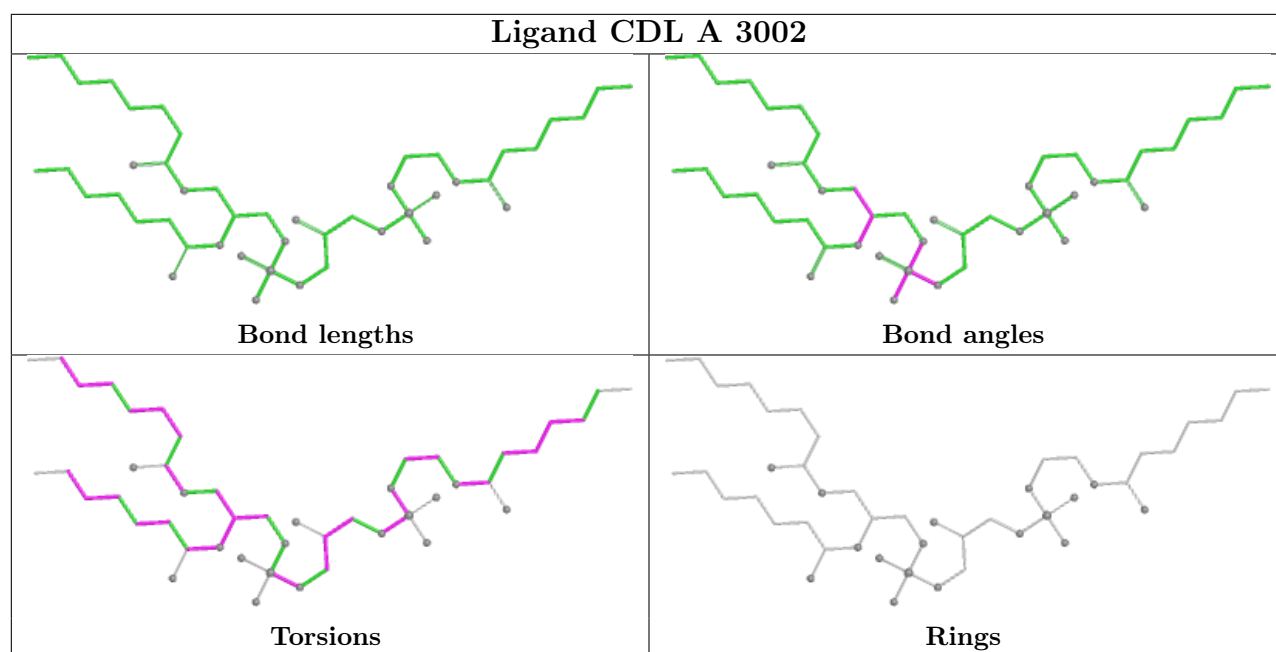


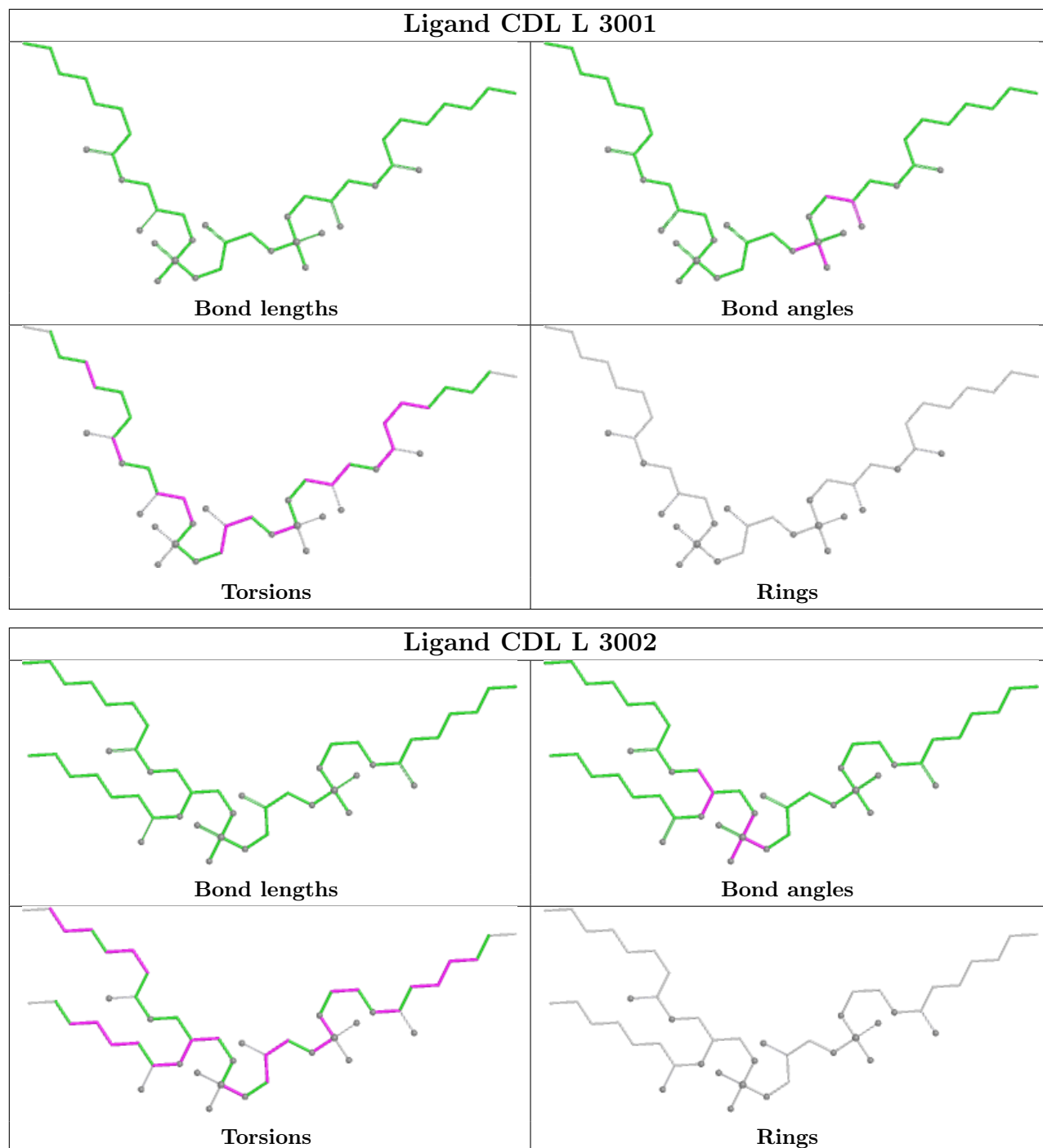


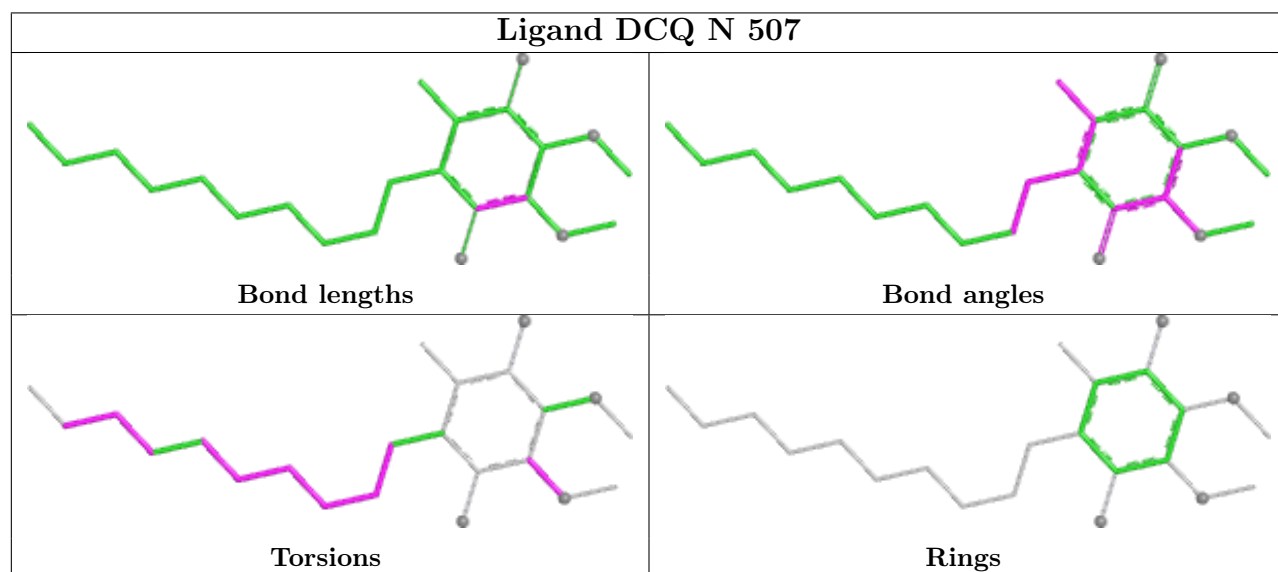
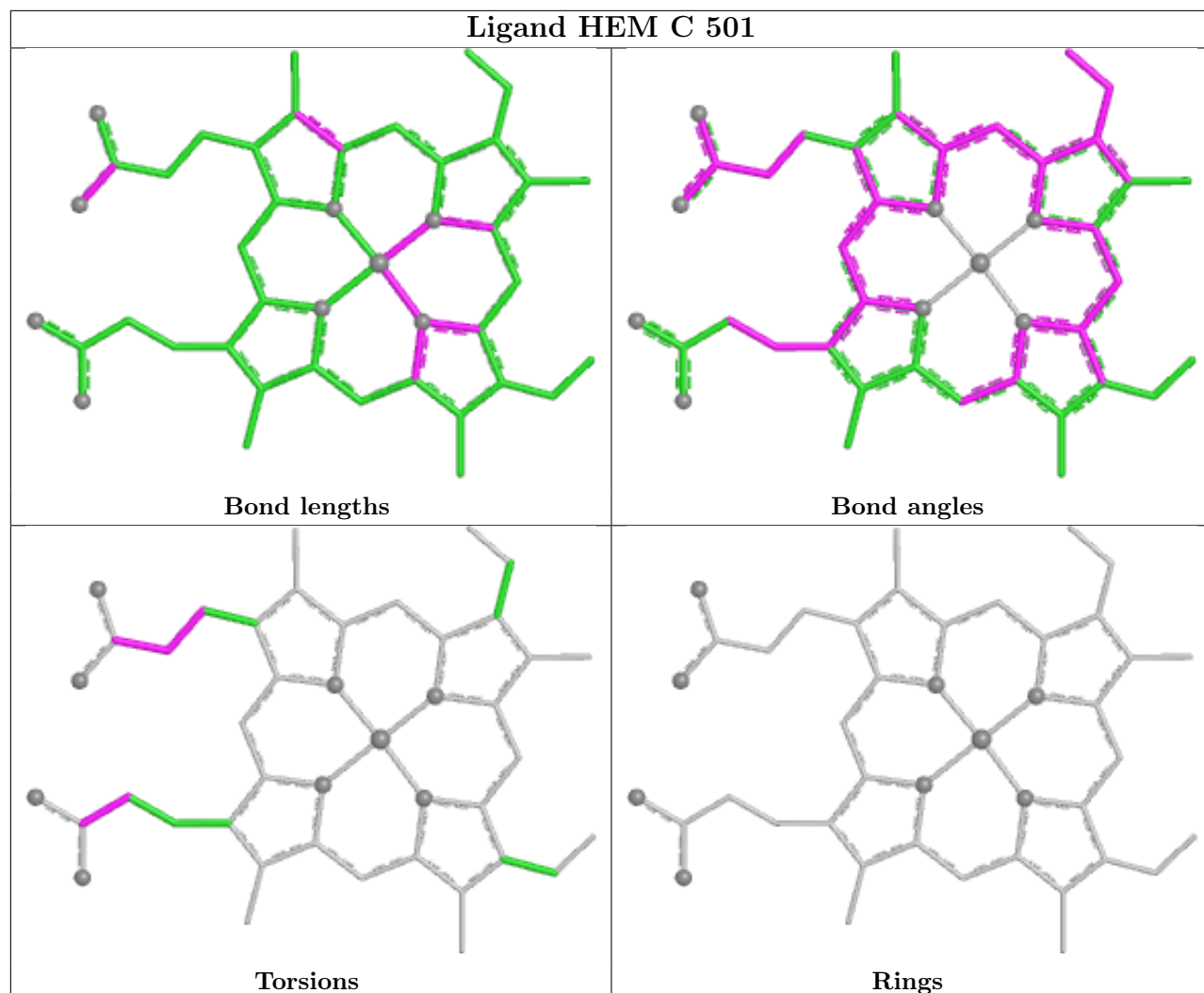


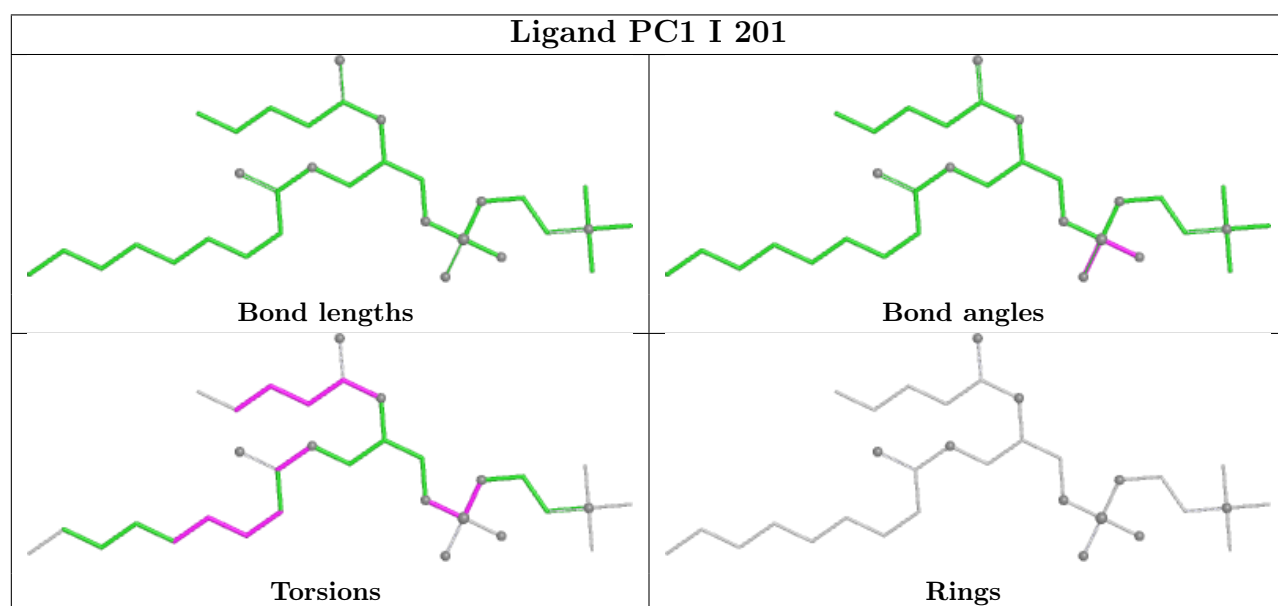
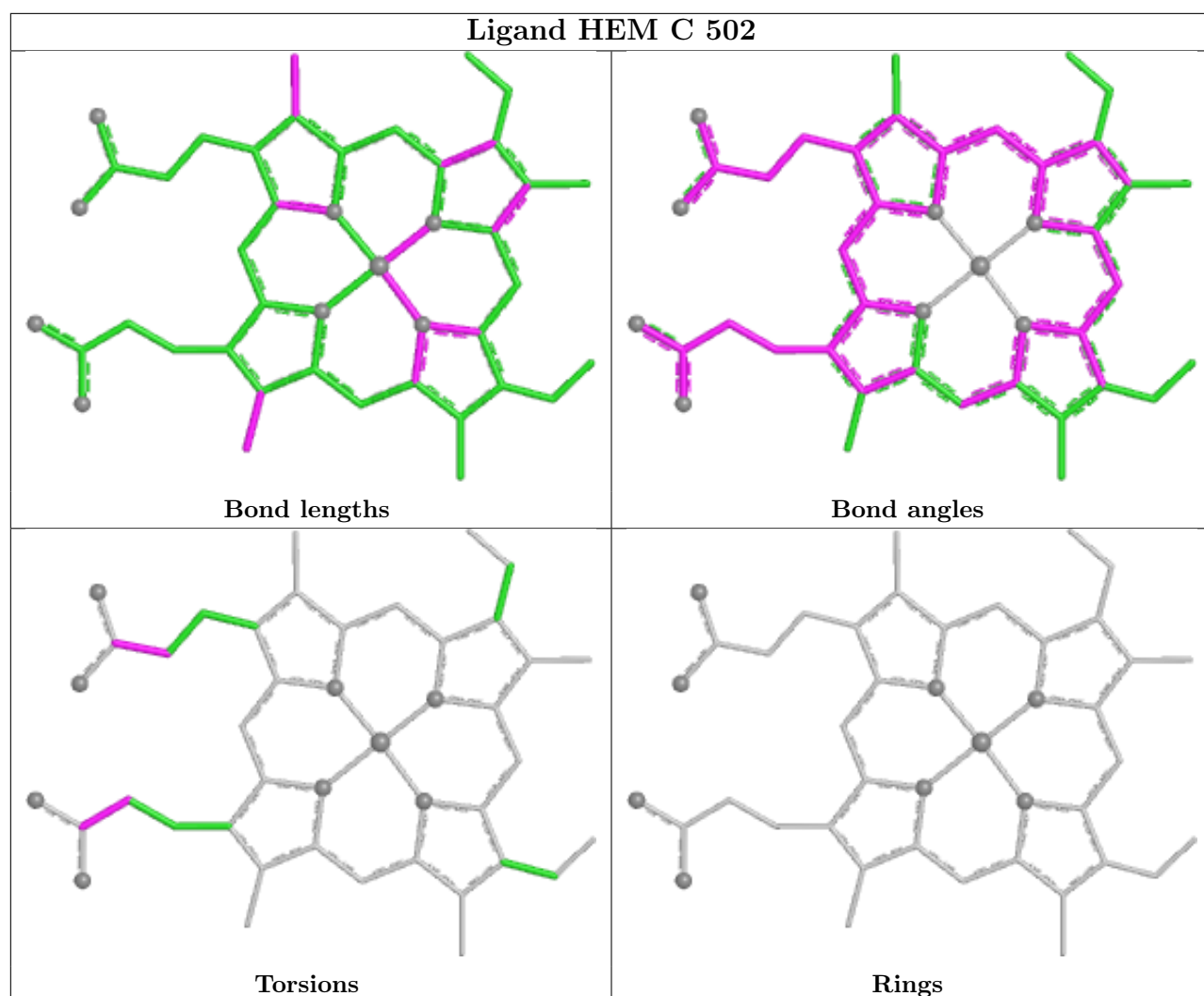


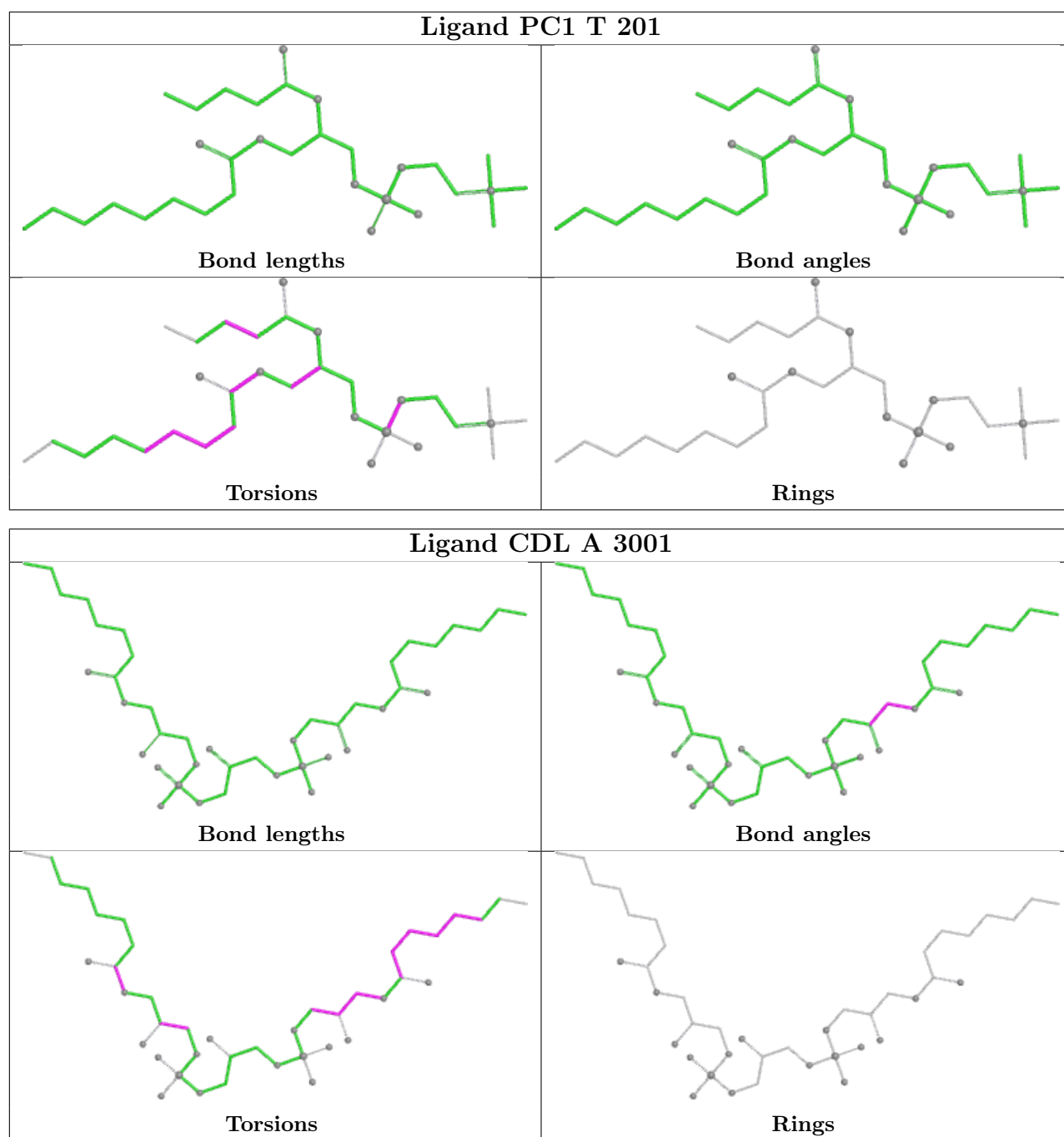


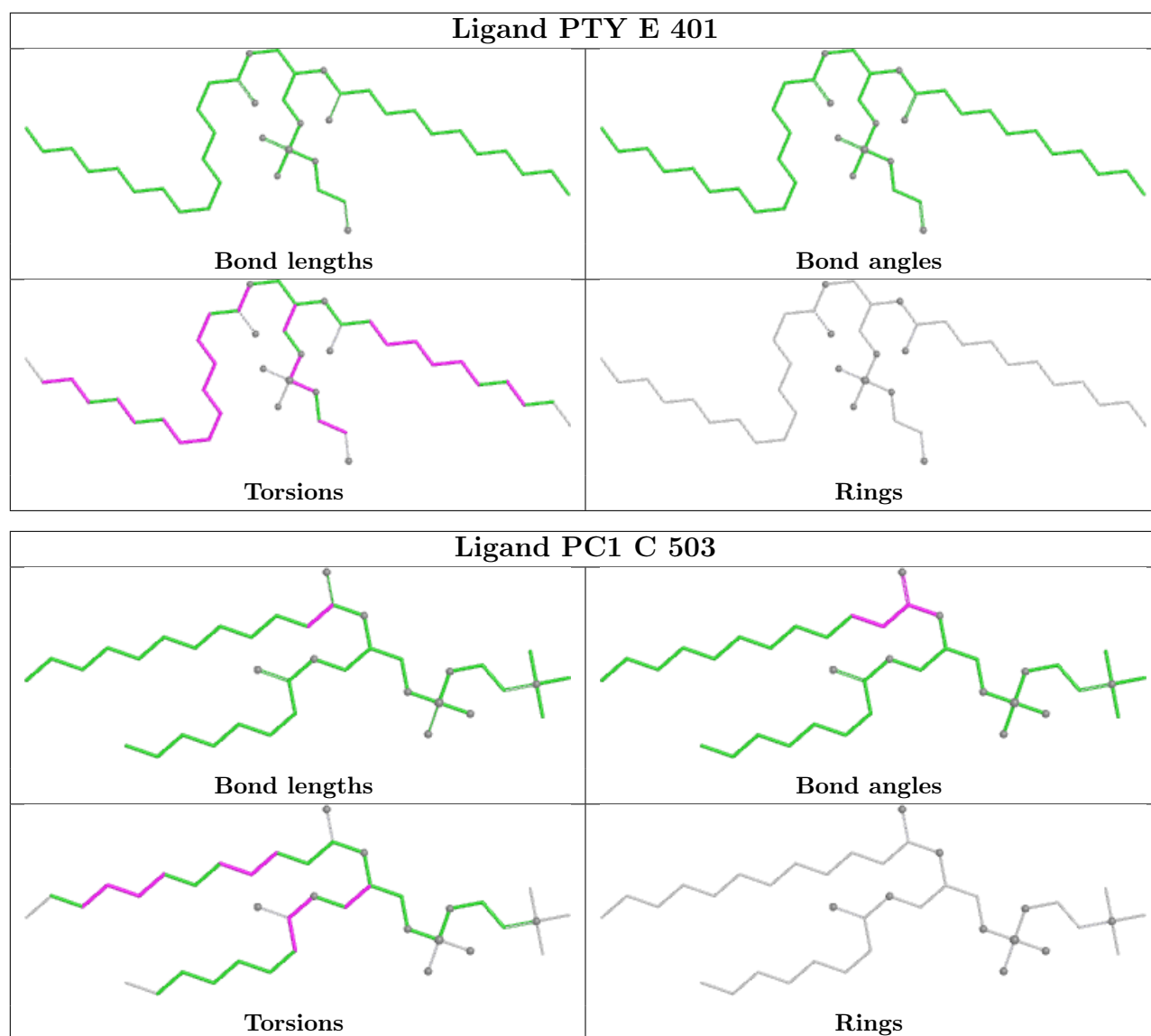


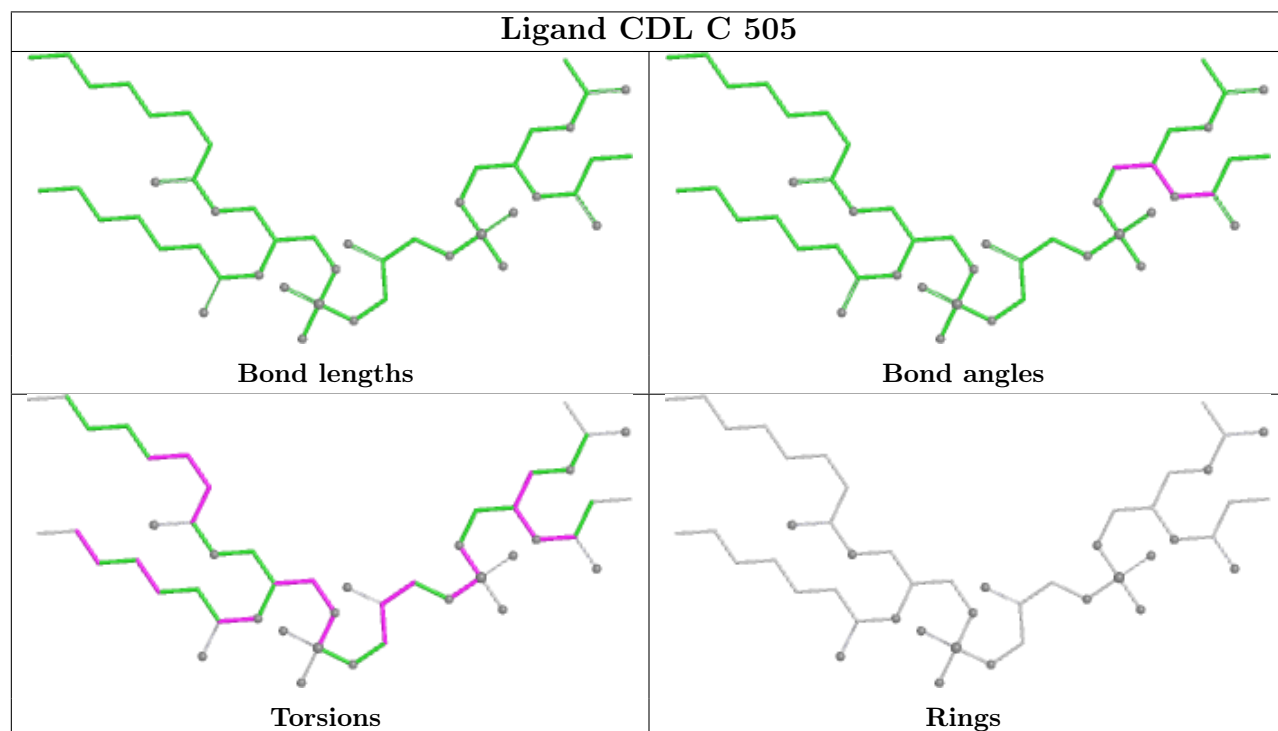
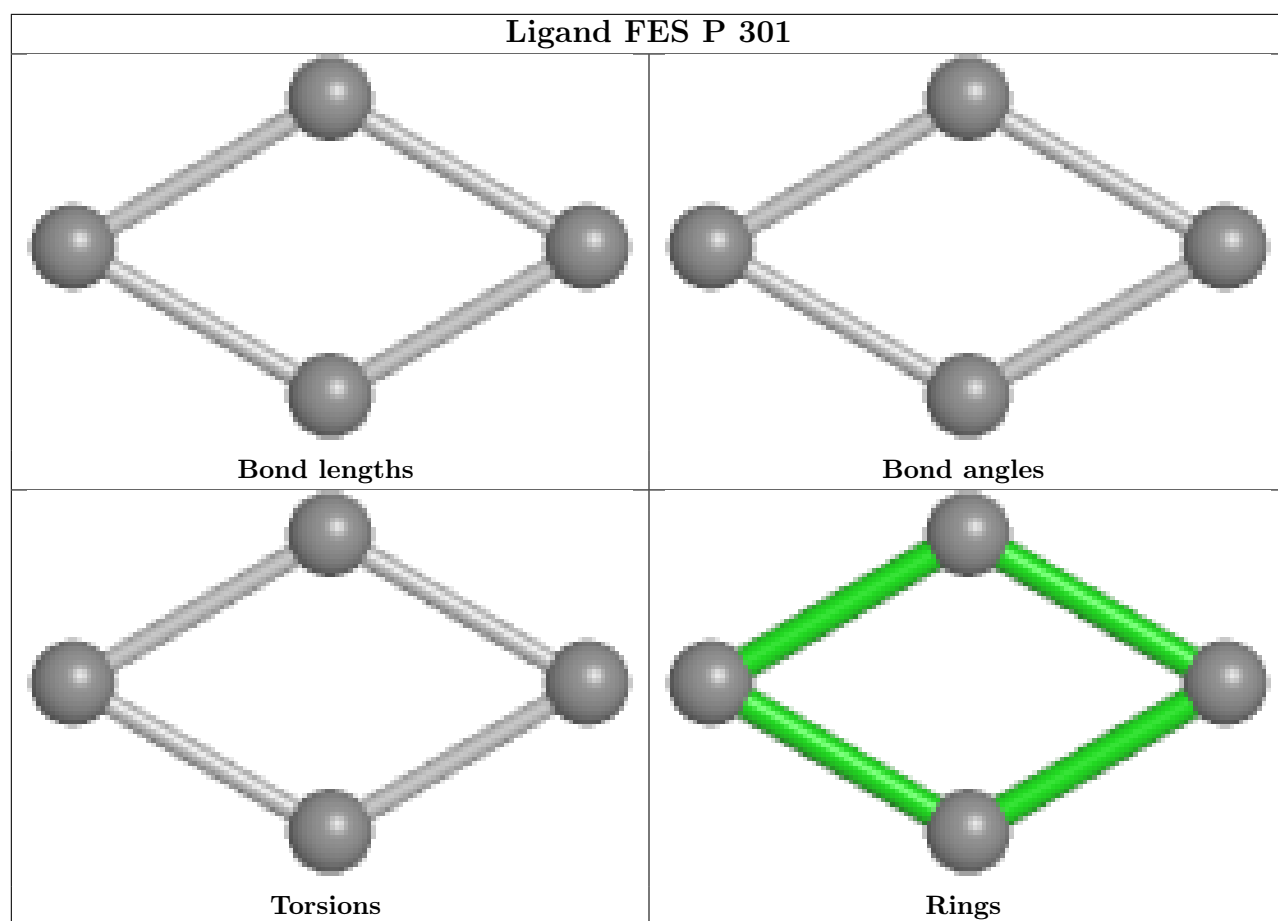


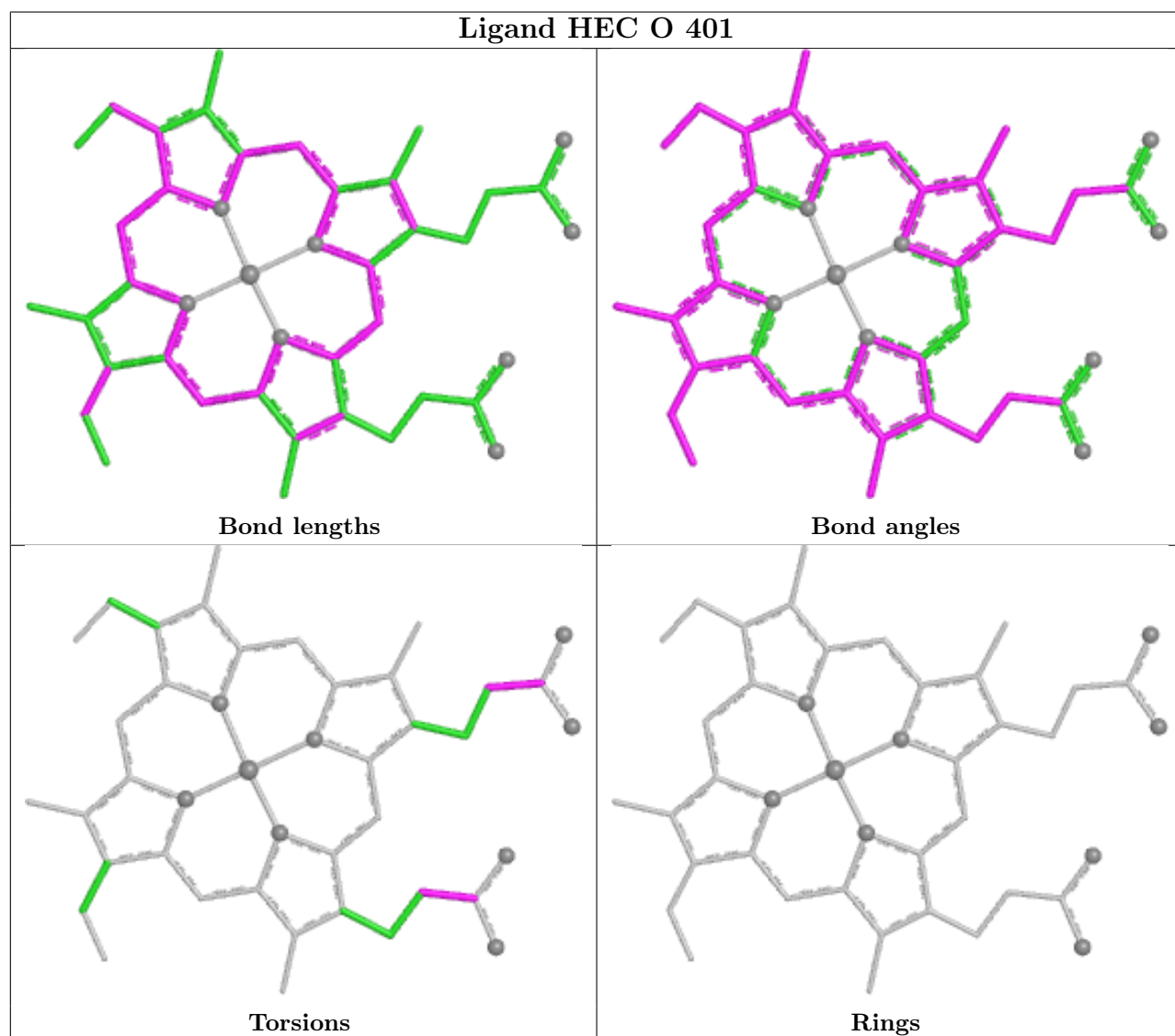
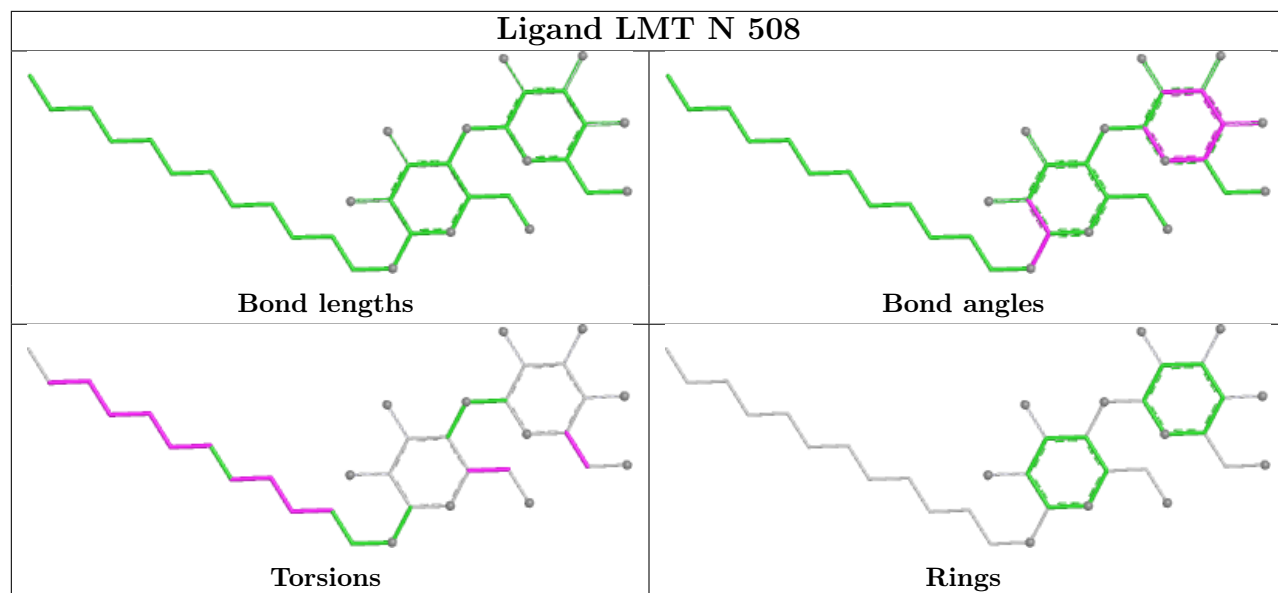


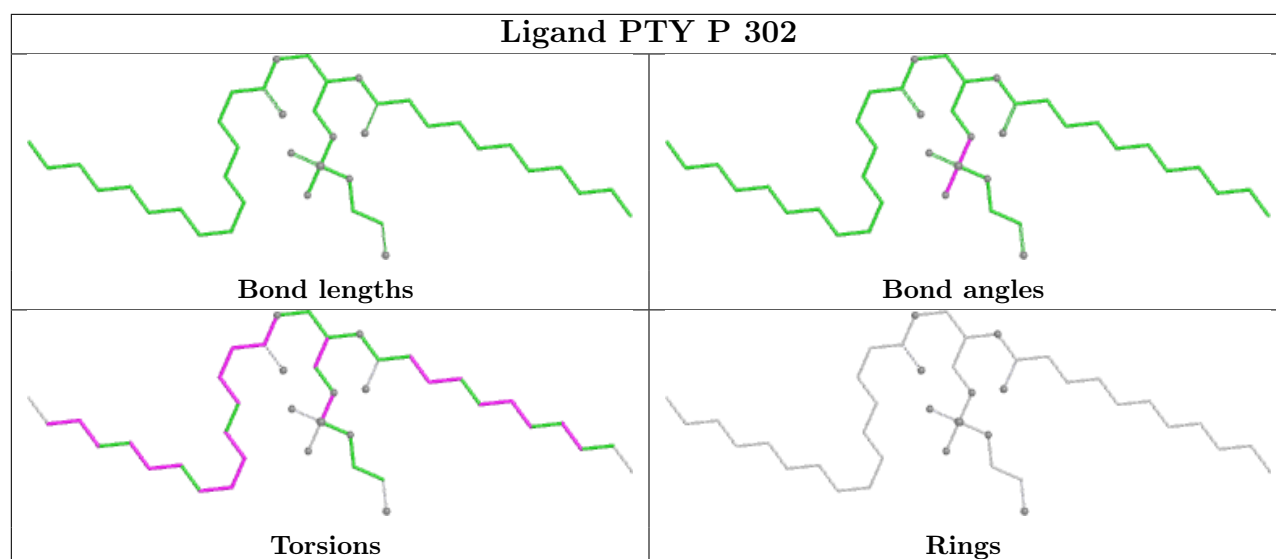












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

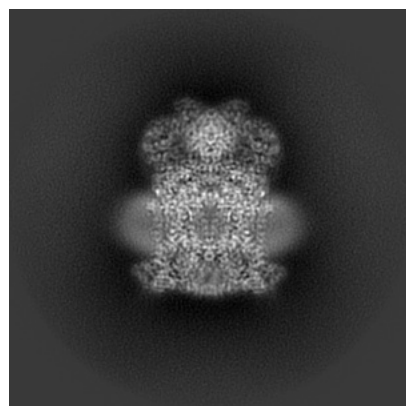
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-15324. These allow visual inspection of the internal detail of the map and identification of artifacts.

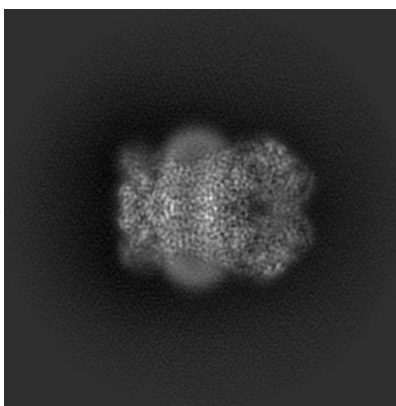
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

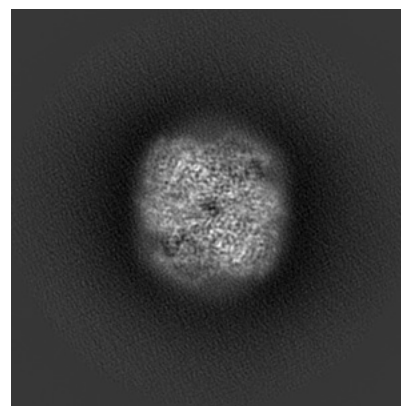
6.1.1 Primary map



X

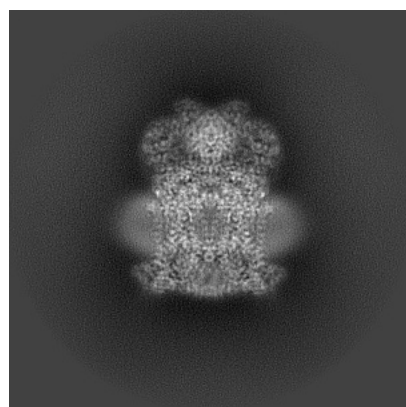


Y

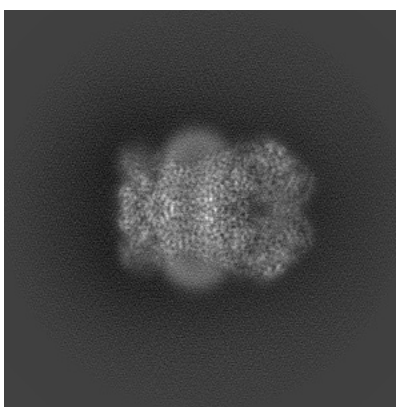


Z

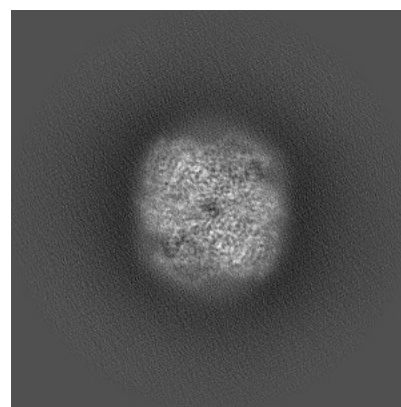
6.1.2 Raw map



X



Y

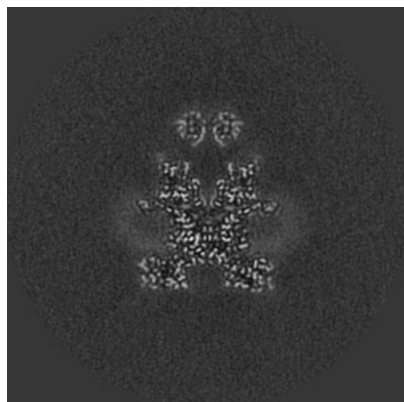


Z

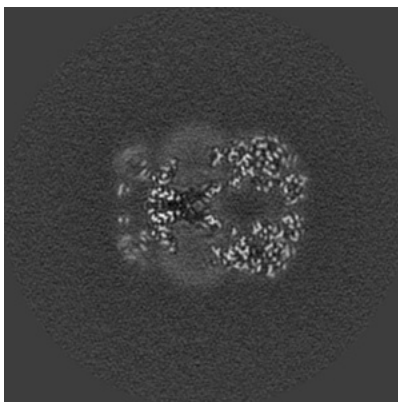
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

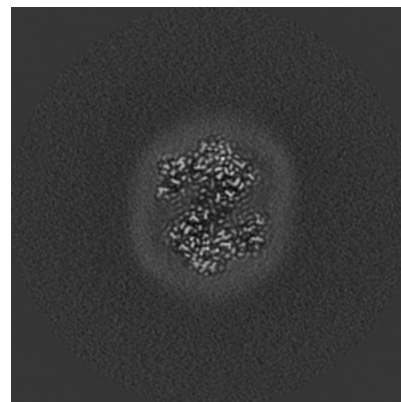
6.2.1 Primary map



X Index: 180

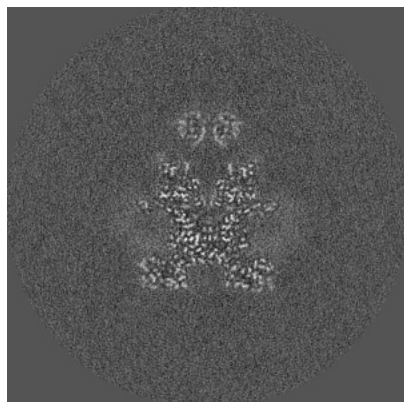


Y Index: 180

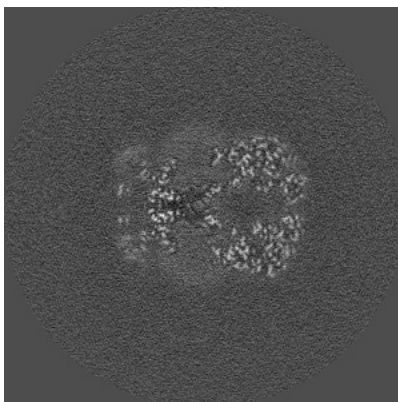


Z Index: 180

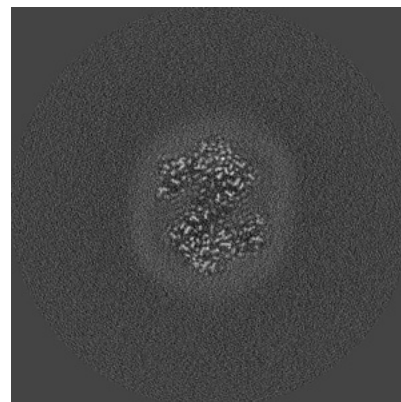
6.2.2 Raw map



X Index: 180



Y Index: 180

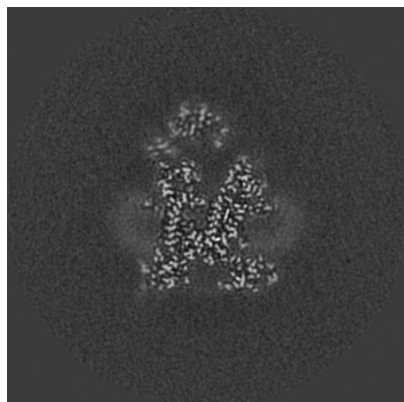


Z Index: 180

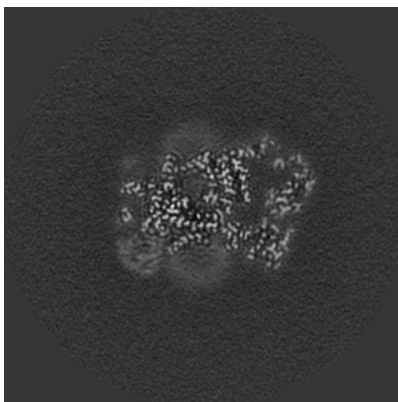
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

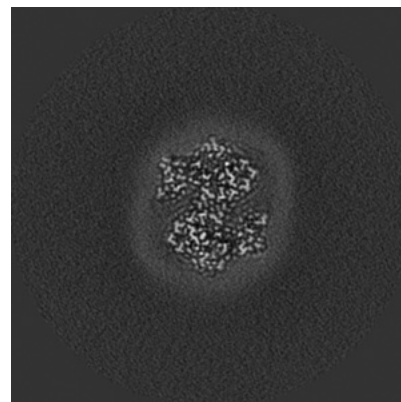
6.3.1 Primary map



X Index: 187

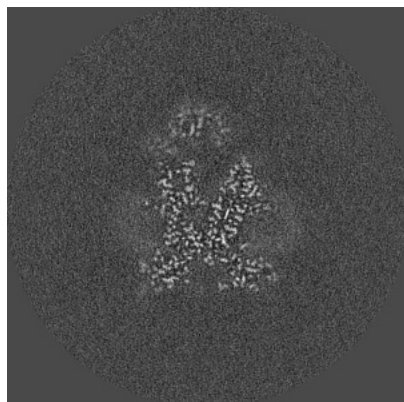


Y Index: 166

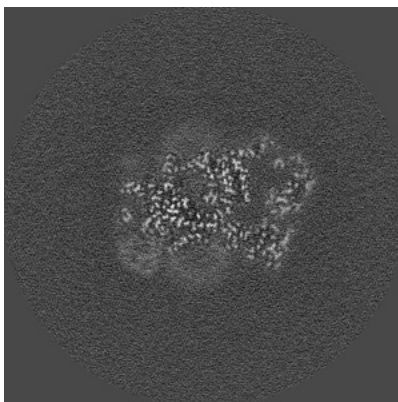


Z Index: 182

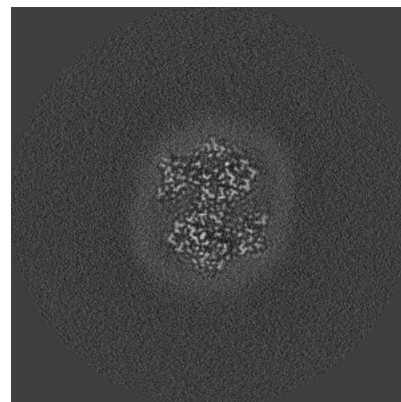
6.3.2 Raw map



X Index: 186



Y Index: 166

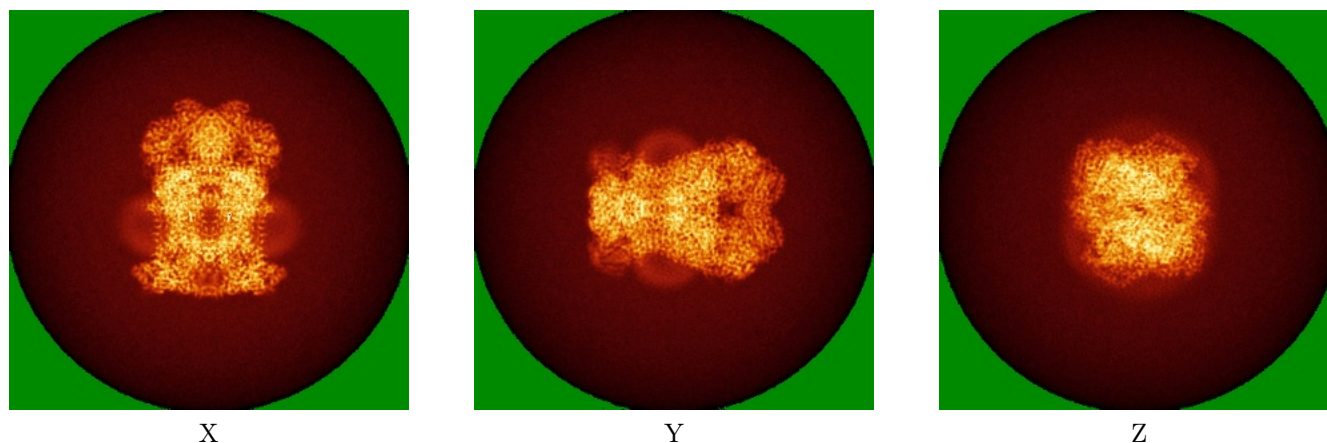


Z Index: 182

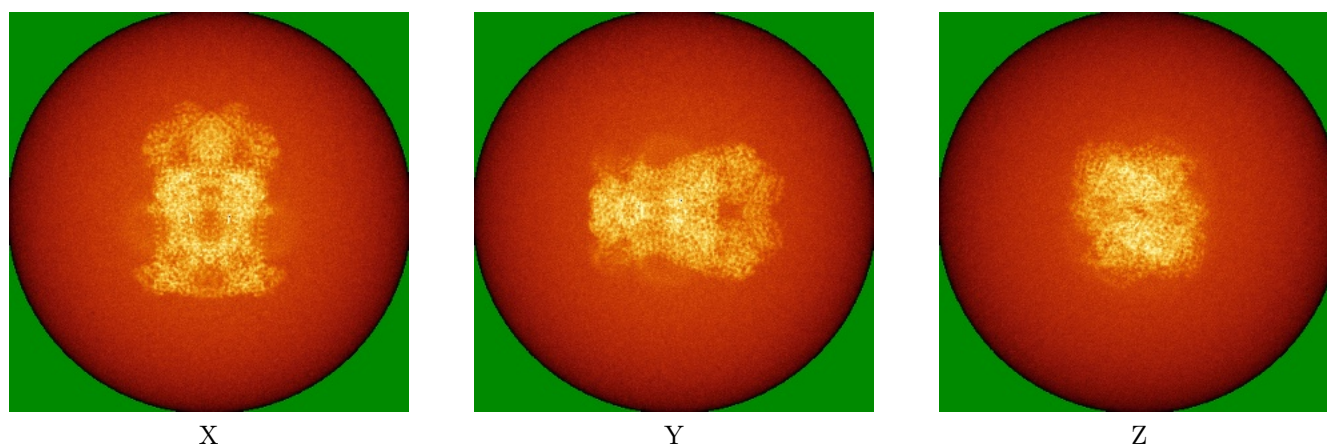
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



6.4.2 Raw map



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](#)

This section was not generated.

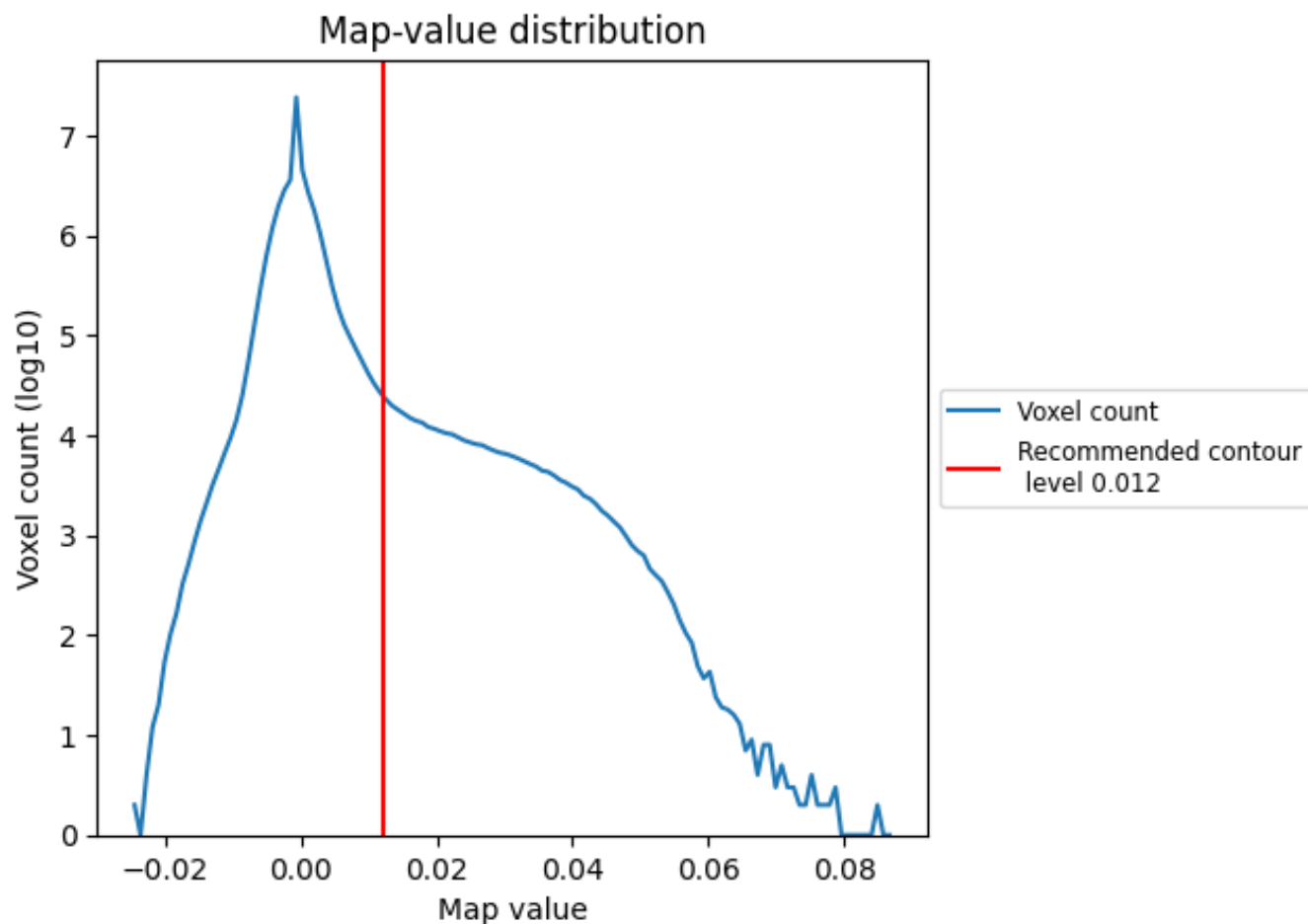
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

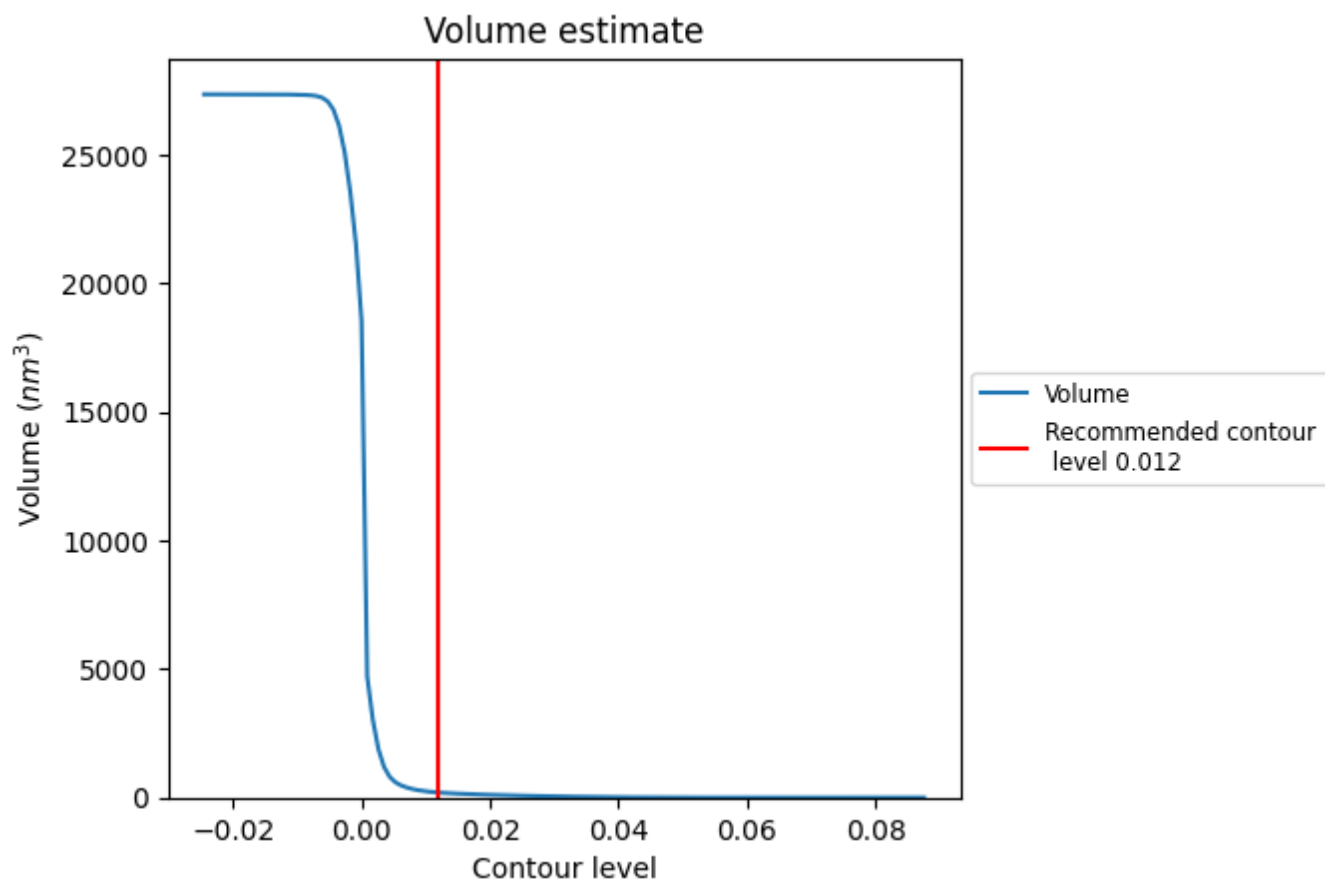
This section contains the results of statistical analysis of the map.

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.

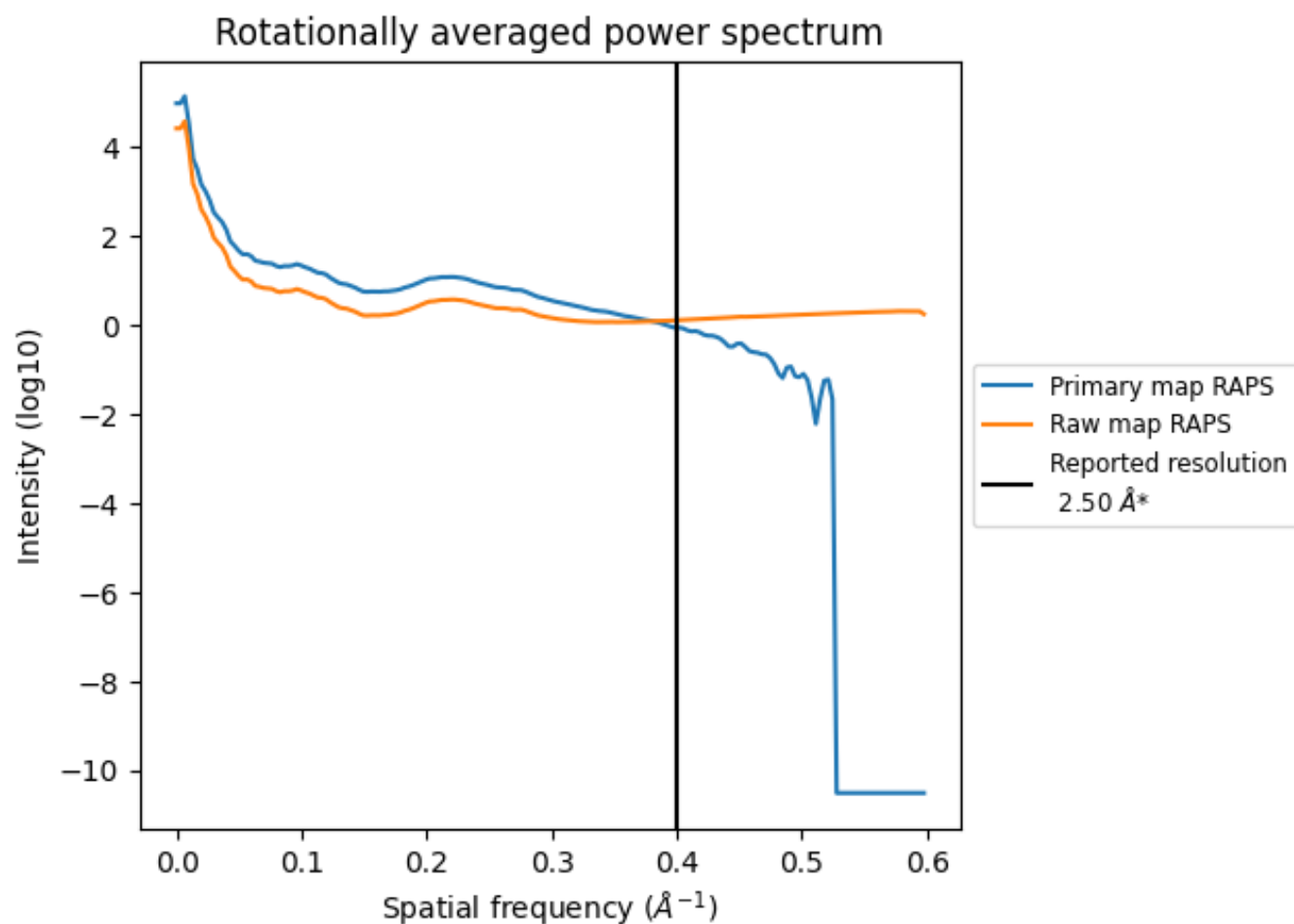
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 195 nm³; this corresponds to an approximate mass of 176 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 ⓘ

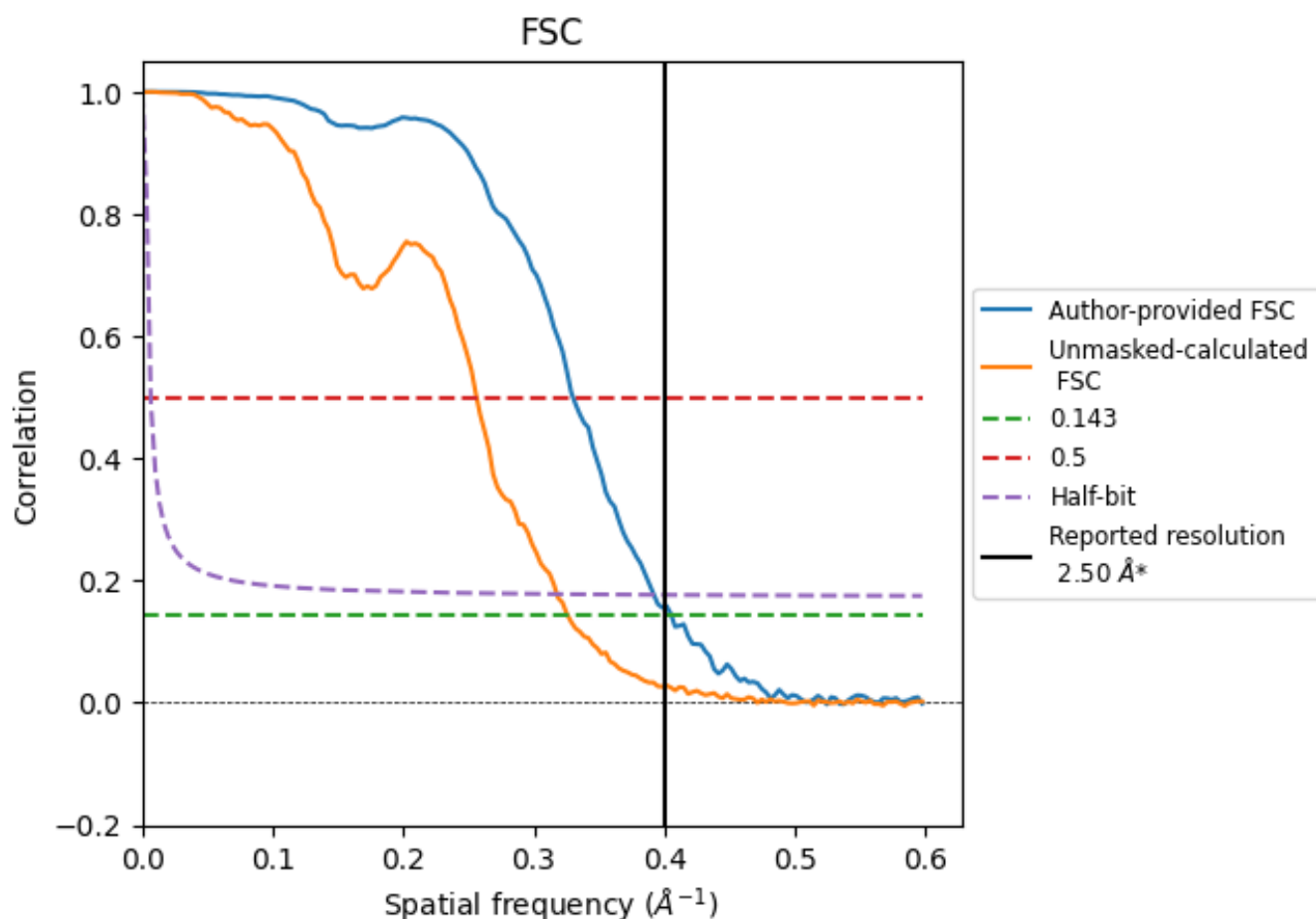


*Reported resolution corresponds to spatial frequency of 0.400 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.400 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

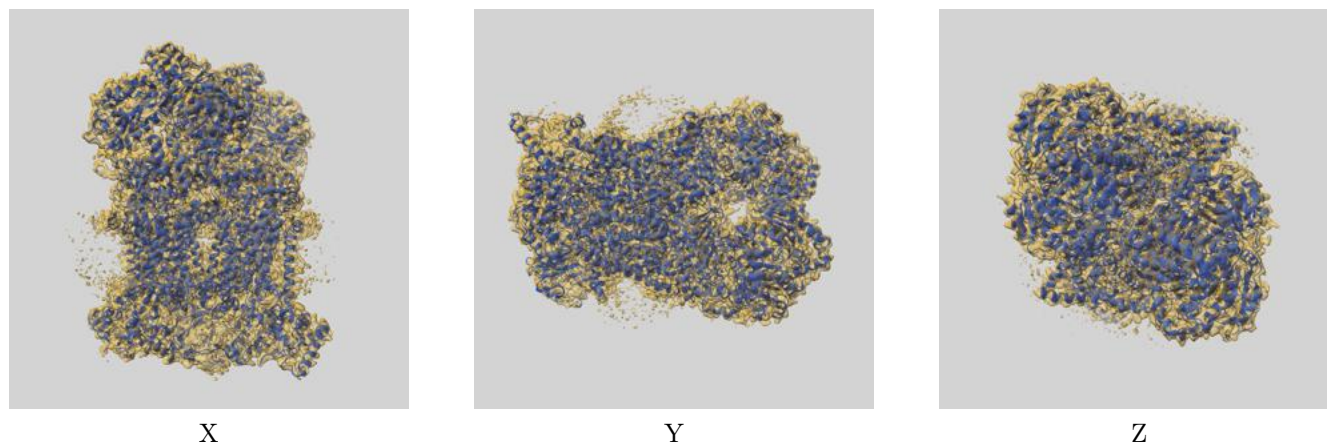
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.50	-	-
Author-provided FSC curve	2.47	3.03	2.55
Unmasked-calculated*	3.07	3.90	3.15

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.07 differs from the reported value 2.5 by more than 10 %

9 Map-model fit [i](#)

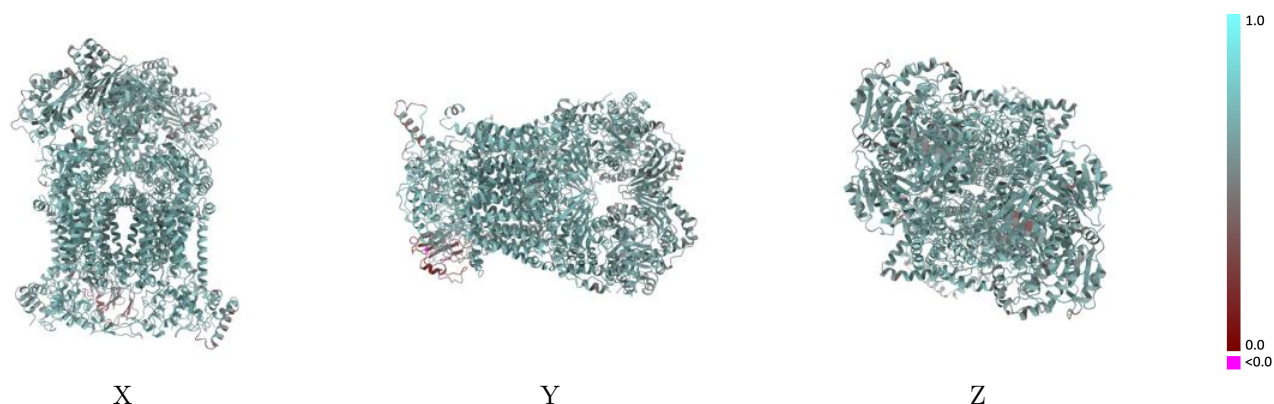
This section contains information regarding the fit between EMDB map EMD-15324 and PDB model 8ABK. Per-residue inclusion information can be found in section [3](#) on page [13](#).

9.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 0.012 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](#)

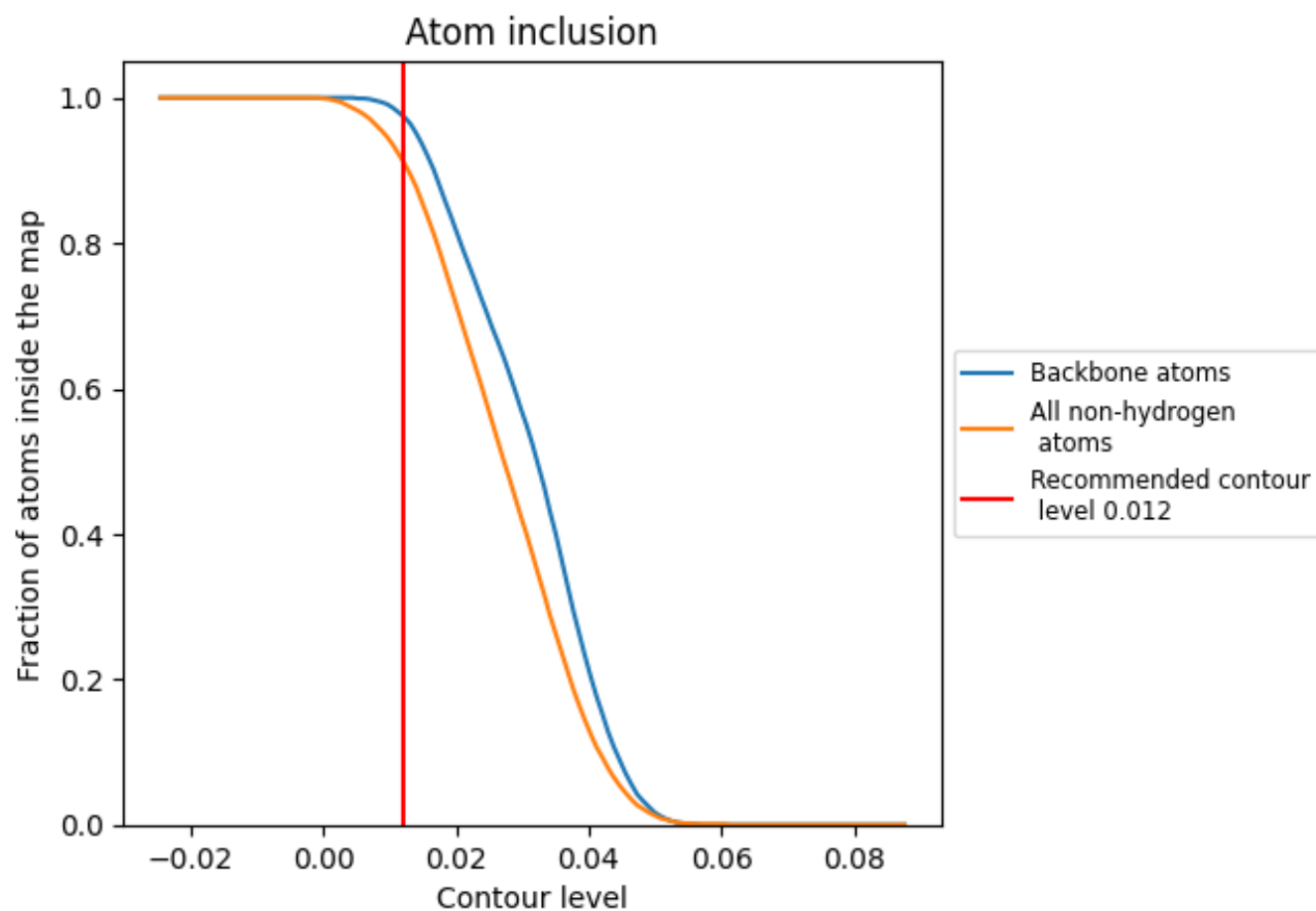


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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](#)

This section was not generated.

9.4 Atom inclusion ⓘ



At the recommended contour level, 98% of all backbone atoms, 91% 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 (0.012) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9130	 0.6330
A	 0.9080	 0.6330
B	 0.8990	 0.6130
C	 0.9700	 0.6770
D	 0.9460	 0.6580
E	 0.9440	 0.6600
F	 0.8040	 0.5540
G	 0.9230	 0.6580
H	 0.9530	 0.6520
I	 0.9600	 0.6660
J	 0.8990	 0.6250
L	 0.9060	 0.6330
M	 0.8980	 0.6160
N	 0.9680	 0.6780
O	 0.9510	 0.6580
P	 0.6640	 0.4220
Q	 0.7780	 0.5480
R	 0.9280	 0.6560
S	 0.9580	 0.6540
T	 0.9560	 0.6560
U	 0.8970	 0.6330

