



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 9, 2021 – 10:39 PM GMT

PDB ID : 6T0F
Title : Crystal structure of CYP124 in complex with cholest-4-en-3-one
Authors : Bukhdruker, S.; Marin, E.; Varaksa, T.; Gilep, A.; Strushkevich, N.; Borshchevskiy, V.
Deposited on : 2019-10-03
Resolution : 1.65 Å(reported)

This is a Full wwPDB X-ray Structure 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/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.16
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.16

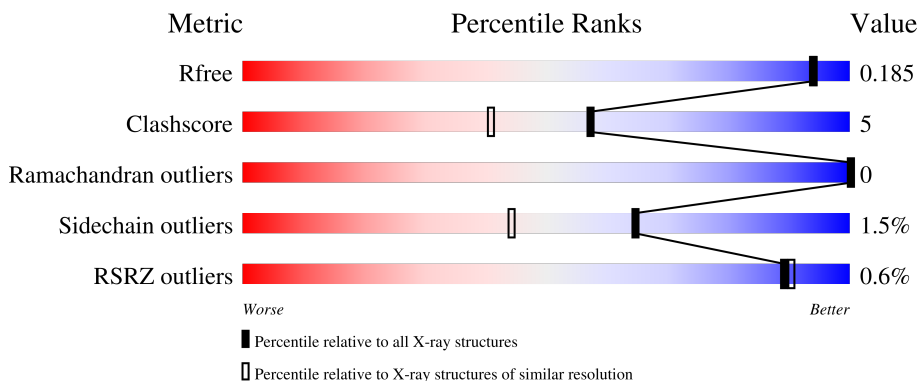
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION




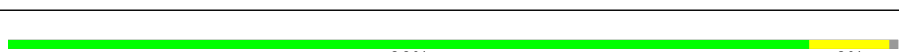
The reported resolution of this entry is 1.65 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1827 (1.66-1.66)
Clashscore	141614	1931 (1.66-1.66)
Ramachandran outliers	138981	1891 (1.66-1.66)
Sidechain outliers	138945	1891 (1.66-1.66)
RSRZ outliers	127900	1791 (1.66-1.66)

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

Mol	Chain	Length	Quality of chain
1	A	435	
1	B	435	
1	C	435	
1	D	435	

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
6	PGE	A	508[B]	-	-	X	-
6	PGE	C	506[A]	-	-	X	-

2 Entry composition i

There are 7 unique types of molecules in this entry. The entry contains 18282 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a protein called Methyl-branched lipid omega-hydroxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	431	Total 3792	C 2377	N 688	O 715	S 12	0	42	0
1	B	430	Total 3758	C 2351	N 680	O 713	S 14	0	44	0
1	C	430	Total 3699	C 2321	N 659	O 705	S 14	0	39	0
1	D	431	Total 3761	C 2357	N 683	O 708	S 13	0	44	0

There are 32 discrepancies between the modelled and reference sequences:

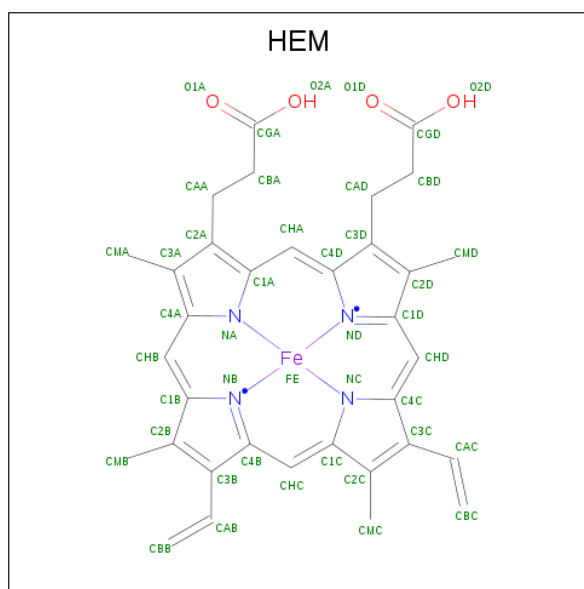
Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	MET	-	initiating methionine	UNP P9WPP3
A	-5	HIS	-	expression tag	UNP P9WPP3
A	-4	HIS	-	expression tag	UNP P9WPP3
A	-3	HIS	-	expression tag	UNP P9WPP3
A	-2	HIS	-	expression tag	UNP P9WPP3
A	-1	HIS	-	expression tag	UNP P9WPP3
A	0	HIS	-	expression tag	UNP P9WPP3
A	65	THR	ALA	engineered mutation	UNP P9WPP3
B	-6	MET	-	initiating methionine	UNP P9WPP3
B	-5	HIS	-	expression tag	UNP P9WPP3
B	-4	HIS	-	expression tag	UNP P9WPP3
B	-3	HIS	-	expression tag	UNP P9WPP3
B	-2	HIS	-	expression tag	UNP P9WPP3
B	-1	HIS	-	expression tag	UNP P9WPP3
B	0	HIS	-	expression tag	UNP P9WPP3
B	65	THR	ALA	engineered mutation	UNP P9WPP3
C	-6	MET	-	initiating methionine	UNP P9WPP3
C	-5	HIS	-	expression tag	UNP P9WPP3
C	-4	HIS	-	expression tag	UNP P9WPP3
C	-3	HIS	-	expression tag	UNP P9WPP3
C	-2	HIS	-	expression tag	UNP P9WPP3

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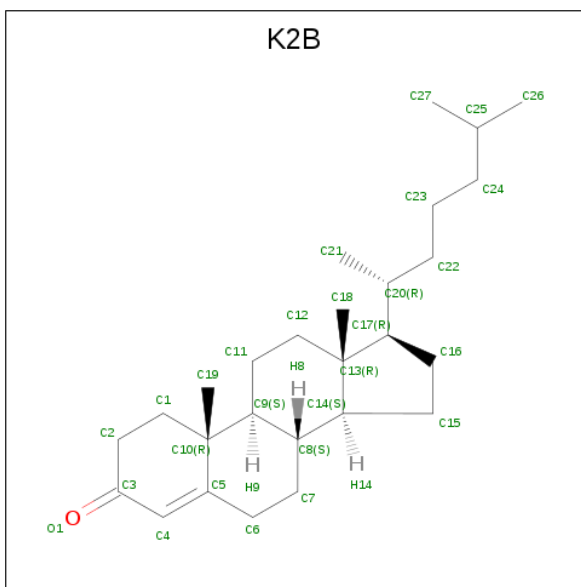
Chain	Residue	Modelled	Actual	Comment	Reference
C	-1	HIS	-	expression tag	UNP P9WPP3
C	0	HIS	-	expression tag	UNP P9WPP3
C	65	THR	ALA	engineered mutation	UNP P9WPP3
D	-6	MET	-	initiating methionine	UNP P9WPP3
D	-5	HIS	-	expression tag	UNP P9WPP3
D	-4	HIS	-	expression tag	UNP P9WPP3
D	-3	HIS	-	expression tag	UNP P9WPP3
D	-2	HIS	-	expression tag	UNP P9WPP3
D	-1	HIS	-	expression tag	UNP P9WPP3
D	0	HIS	-	expression tag	UNP P9WPP3
D	65	THR	ALA	engineered mutation	UNP P9WPP3

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



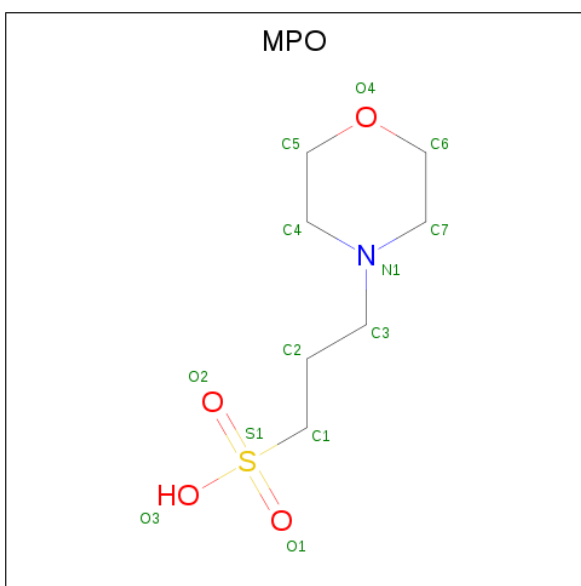
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
2	A	1	Total	C	Fe	N	O	0	1
			86	68	2	8	8		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	D	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is (8ALPHA,9BETA)-CHOLEST-4-EN-3-ONE (three-letter code: K2B) (formula: $C_{27}H_{44}O$) (labeled as "Ligand of Interest" by depositor).



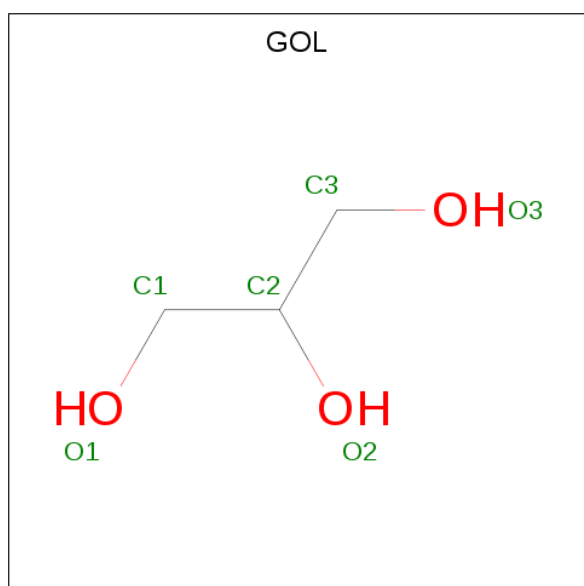
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	C O	0	0
			28	27 1		
3	B	1	Total	C O	0	0
			28	27 1		
3	C	1	Total	C O	0	0
			28	27 1		
3	D	1	Total	C O	0	0
			28	27 1		

- Molecule 4 is 3[N-MORPHOLINO]PROPANE SULFONIC ACID (three-letter code: MPO) (formula: C₇H₁₅NO₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	1
			26	14	2	8	2		
4	B	1	Total	C	N	O	S	0	1
			26	14	2	8	2		
4	C	1	Total	C	N	O	S	0	1
			13	7	1	4	1		
4	D	1	Total	C	N	O	S	0	1
			26	14	2	8	2		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



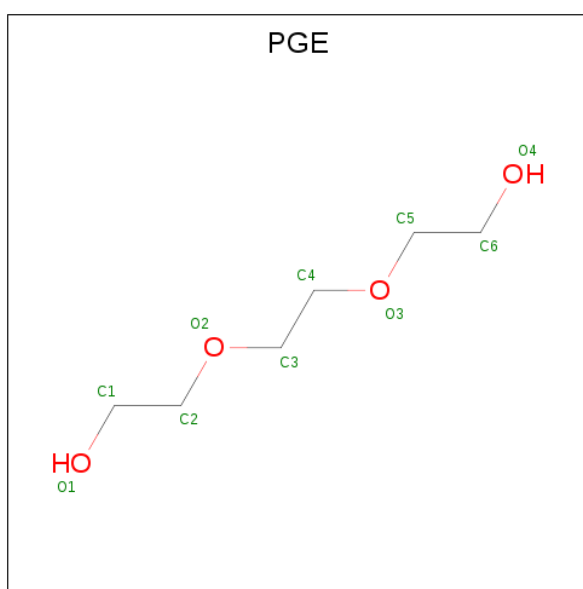
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	C	O	0	0
			6	3	3		
5	C	1	Total	C	O	0	0
			6	3	3		
5	D	1	Total	C	O	0	0
			6	3	3		
5	D	1	Total	C	O	0	1
			6	3	3		

- Molecule 6 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C₆H₁₄O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	1
			7	4	3		
6	B	1	Total	C	O	0	1
			7	4	3		
6	B	1	Total	C	O	0	1
			7	4	3		
6	C	1	Total	C	O	0	1
			10	6	4		
6	C	1	Total	C	O	0	1
			7	4	3		
6	D	1	Total	C	O	0	0
			10	6	4		

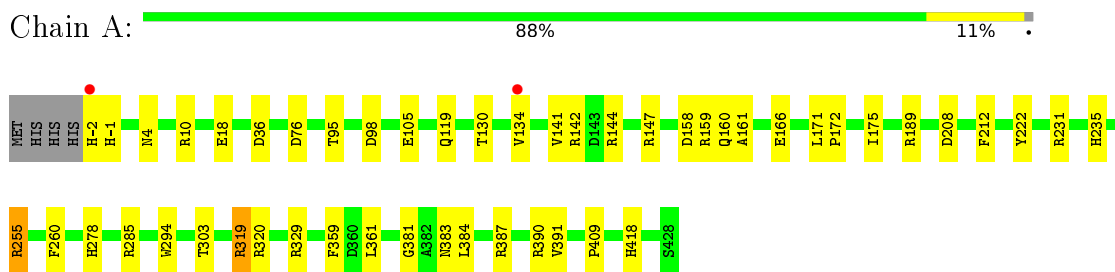
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	640	Total 671	O 671	0	64
7	B	604	Total 644	O 644	0	60
7	C	638	Total 672	O 672	0	56
7	D	709	Total 747	O 747	0	78

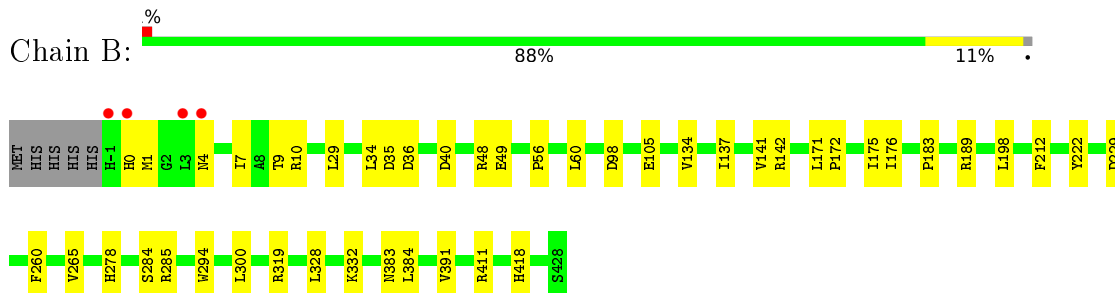
3 Residue-property plots [i](#)

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

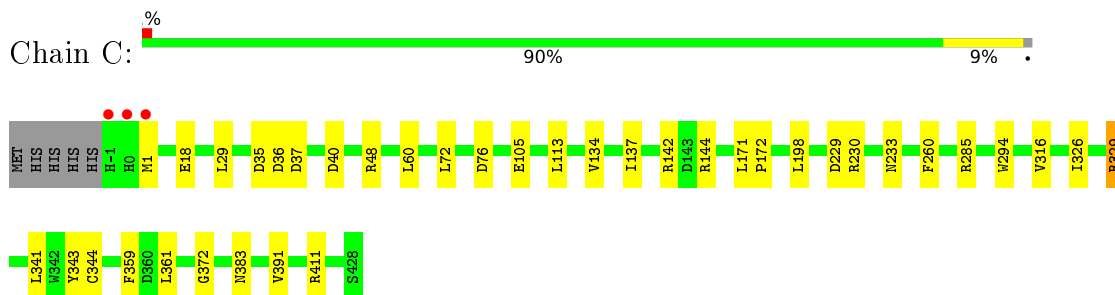
- Molecule 1: Methyl-branched lipid omega-hydroxylase



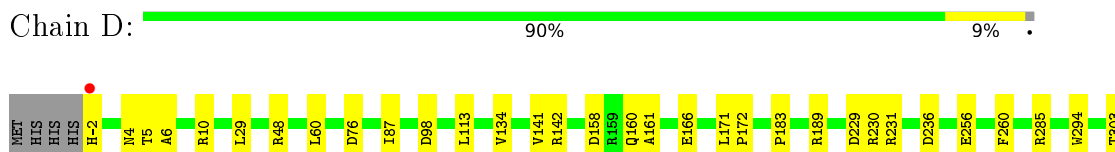
- Molecule 1: Methyl-branched lipid omega-hydroxylase



- Molecule 1: Methyl-branched lipid omega-hydroxylase



- Molecule 1: Methyl-branched lipid omega-hydroxylase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	93.67Å 81.27Å 155.85Å 90.00° 107.29° 90.00°	Depositor
Resolution (Å)	29.75 – 1.65 29.75 – 1.65	Depositor EDS
% Data completeness (in resolution range)	99.5 (29.75-1.65) 99.4 (29.75-1.65)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.30 (at 1.65Å)	Xtrriage
Refinement program	PHENIX 1.16_3549	Depositor
R, R_{free}	0.152 , 0.185 0.152 , 0.185	Depositor DCC
R_{free} test set	13343 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	21.9	Xtrriage
Anisotropy	0.342	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 44.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	0.196 for h,-k,-h-l	Xtrriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	18282	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 64.95 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 7.6461e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: K2B, GOL, PGE, HEM, MPO

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.55	0/3887	0.69	1/5289 (0.0%)
1	B	0.56	0/3848	0.67	0/5236
1	C	0.61	0/3793	0.69	0/5162
1	D	0.60	0/3865	0.70	0/5259
All	All	0.58	0/15393	0.69	1/20946 (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	A	208	ASP	CB-CG-OD1	5.21	122.99	118.30

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	A	3792	0	3638	44	0
1	B	3758	0	3606	41	0
1	C	3699	0	3537	33	0
1	D	3761	0	3605	29	0
2	A	86	0	60	2	0
2	B	43	0	30	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	43	0	30	0	0
2	D	43	0	30	0	0
3	A	28	0	44	1	0
3	B	28	0	44	1	0
3	C	28	0	44	0	0
3	D	28	0	44	0	0
4	A	26	0	28	1	0
4	B	26	0	28	5	0
4	C	13	0	15	0	0
4	D	26	0	30	1	0
5	A	30	0	39	2	0
5	B	18	0	23	3	0
5	C	12	0	16	0	0
5	D	12	0	14	1	0
6	A	7	0	9	10	0
6	B	14	0	18	8	0
6	C	17	0	23	12	0
6	D	10	0	14	2	0
7	A	671	0	0	14	0
7	B	644	0	0	13	0
7	C	672	0	0	7	0
7	D	747	0	0	9	0
All	All	18282	0	14969	152	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 5.

All (152) 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)
1:A:-1:HIS:HB2	5:A:505:GOL:H2	1.52	0.89
1:B:141:VAL:HG12	1:B:391[B]:VAL:HG21	1.56	0.88
6:C:506[A]:PGE:H42	7:C:1005:HOH:O	1.77	0.84
1:A:141:VAL:HG12	1:A:391[B]:VAL:HG11	1.61	0.82
1:D:141:VAL:HG12	1:D:391[B]:VAL:HG21	1.58	0.81
1:D:183:PRO:HG3	4:D:503[A]:MPO:H22	1.66	0.77
1:A:98:ASP:OD1	1:A:319[B]:ARG:NH2	2.17	0.76
2:A:501[B]:HEM:HHD	2:A:501[B]:HEM:HBC2	1.66	0.76
1:C:411[B]:ARG:NH2	7:C:601:HOH:O	2.21	0.74
5:B:506:GOL:H31	7:B:1070:HOH:O	1.89	0.73
1:C:37[A]:ASP:OD1	6:C:506[A]:PGE:H3	1.90	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1[B]:MET:HE1	1:D:303:THR:HA	1.75	0.69
1:A:278:HIS:CE1	6:A:508[B]:PGE:H1	2.28	0.68
1:D:10[B]:ARG:HD2	1:D:285:ARG:HD3	1.76	0.68
6:B:507[A]:PGE:H32	7:B:605:HOH:O	1.95	0.66
1:B:98:ASP:OD1	1:B:319[A]:ARG:NH2	2.24	0.66
1:A:142:ARG:HA	1:A:391[B]:VAL:HG21	1.79	0.65
1:C:316:VAL:HG13	1:C:344[B]:CYS:SG	2.37	0.65
1:D:256:GLU:HG3	5:D:506[A]:GOL:H32	1.79	0.64
1:A:18[B]:GLU:OE2	7:A:601[B]:HOH:O	2.15	0.64
1:B:10[B]:ARG:NH1	7:B:1113[B]:HOH:O	2.31	0.64
1:D:98:ASP:OD1	1:D:319[B]:ARG:NH2	2.27	0.64
1:B:134:VAL:O	1:B:137:ILE:HG22	1.97	0.64
1:A:409:PRO:HG2	6:A:508[B]:PGE:H42	1.80	0.63
1:A:418:HIS:CD2	6:A:508[B]:PGE:H12	2.33	0.63
1:C:134:VAL:O	1:C:137:ILE:HG22	1.99	0.63
1:D:387[B]:ARG:HD2	1:D:387[B]:ARG:N	2.12	0.63
1:B:383[A]:ASN:ND2	7:B:604:HOH:O	2.32	0.63
1:B:10[B]:ARG:HD2	1:B:285[B]:ARG:HD3	1.79	0.62
1:C:326[B]:ILE:HG22	7:C:605:HOH:O	1.99	0.62
1:D:76[B]:ASP:OD1	1:D:329:ARG:NH2	2.33	0.61
1:C:113:LEU:HD11	6:C:507[A]:PGE:H5	1.81	0.61
1:C:285:ARG:HH12	6:C:506[A]:PGE:H52	1.66	0.61
1:C:229[B]:ASP:OD2	1:C:230:ARG:NH1	2.34	0.60
1:C:316:VAL:HA	1:C:344[B]:CYS:SG	2.41	0.59
1:A:144:ARG:NH2	7:A:604:HOH:O	2.27	0.59
1:D:383[B]:ASN:ND2	7:D:912[B]:HOH:O	2.25	0.58
1:B:222:TYR:CE2	4:B:503[B]:MPO:H52	2.38	0.58
6:A:508[B]:PGE:H32	7:A:621:HOH:O	2.05	0.57
1:C:1[B]:MET:HG3	1:D:364[B]:ASN:OD1	2.05	0.56
1:A:383[A]:ASN:ND2	7:A:611[A]:HOH:O	2.38	0.56
1:C:36[A]:ASP:HB2	6:C:506[A]:PGE:H5	1.87	0.56
1:D:4[B]:ASN:HB3	1:D:160[B]:GLN:OE1	2.06	0.56
1:D:142:ARG:HA	1:D:391[A]:VAL:HG11	1.88	0.56
1:A:278:HIS:CE1	6:A:508[B]:PGE:C1	2.89	0.56
1:C:142:ARG:HA	1:C:391:VAL:HG11	1.88	0.55
1:B:4[A]:ASN:O	7:B:601[A]:HOH:O	2.18	0.55
1:D:4[B]:ASN:OD1	1:D:5[B]:THR:N	2.32	0.55
1:A:158:ASP:HB2	7:A:1054[A]:HOH:O	2.08	0.54
1:A:159:ARG:NH1	7:A:605:HOH:O	2.31	0.54
1:A:189[B]:ARG:NH2	7:A:614[B]:HOH:O	2.39	0.54
1:C:285:ARG:HH12	6:C:506[A]:PGE:H6	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:332:LYS:NZ	7:B:612:HOH:O	2.40	0.54
1:B:229:ASP:OD2	4:B:503[B]:MPO:H32	2.08	0.54
1:B:189[B]:ARG:NH2	7:B:1029[B]:HOH:O	2.40	0.53
1:C:285:ARG:HH12	6:C:506[A]:PGE:C5	2.22	0.53
1:C:72:LEU:HD12	1:C:341[B]:LEU:CD2	2.38	0.53
1:C:29:LEU:HD22	1:C:60[A]:LEU:HD21	1.90	0.52
1:A:134[B]:VAL:HG12	1:A:387[B]:ARG:HD2	1.93	0.51
1:B:48[B]:ARG:NH1	1:B:49:GLU:OE2	2.43	0.51
1:D:161:ALA:HB1	1:D:166[B]:GLU:HG3	1.93	0.51
1:D:171:LEU:HB3	1:D:172:PRO:HD3	1.91	0.51
1:C:29:LEU:HD22	1:C:60[B]:LEU:HD21	1.93	0.50
1:B:176[B]:ILE:HD12	1:B:265:VAL:HA	1.94	0.50
6:D:504:PGE:H32	7:D:667:HOH:O	2.11	0.50
1:B:35:ASP:HA	1:B:411[B]:ARG:CZ	2.41	0.50
1:C:1[A]:MET:HE2	7:D:1168:HOH:O	2.10	0.50
1:B:171:LEU:HB3	1:B:172:PRO:HD3	1.94	0.49
1:C:76[B]:ASP:OD1	1:C:329:ARG:NH2	2.43	0.49
1:D:134:VAL:HG12	1:D:387[A]:ARG:HD2	1.93	0.49
1:B:142:ARG:HA	1:B:391[A]:VAL:HG11	1.94	0.49
1:B:29:LEU:HD22	1:B:60[A]:LEU:HD21	1.94	0.49
1:A:4[B]:ASN:N	1:A:4[B]:ASN:OD1	2.44	0.49
1:A:142:ARG:HA	1:A:391[A]:VAL:HG11	1.94	0.49
1:C:383[A]:ASN:HB3	7:C:857[A]:HOH:O	2.12	0.49
1:A:130:THR:O	1:A:134[B]:VAL:HG23	2.13	0.49
1:A:212:PHE:CZ	3:A:502:K2B:H12C	2.48	0.49
1:A:390[B]:ARG:NH1	7:A:623[B]:HOH:O	2.46	0.48
1:A:76[B]:ASP:OD1	1:A:329:ARG:NH2	2.45	0.48
1:A:171:LEU:HB3	1:A:172:PRO:HD3	1.95	0.48
1:B:411[A]:ARG:HD2	6:B:507[A]:PGE:H12	1.94	0.48
1:B:40:ASP:OD2	6:B:508[A]:PGE:H1	2.13	0.48
1:B:278:HIS:HD2	7:B:1082[B]:HOH:O	1.95	0.48
6:A:508[B]:PGE:H4	7:A:633:HOH:O	2.12	0.48
1:A:36[B]:ASP:OD2	6:A:508[B]:PGE:H2	2.14	0.48
2:B:501:HEM:HBC2	2:B:501:HEM:HHD	1.96	0.47
1:C:285:ARG:NH1	6:C:506[A]:PGE:H6	2.28	0.47
1:A:278:HIS:HE1	6:A:508[B]:PGE:C1	2.27	0.47
1:C:359:PHE:CZ	1:C:361:LEU:HD21	2.49	0.47
1:A:303:THR:HG22	1:B:1[B]:MET:HE1	1.97	0.47
1:D:4[B]:ASN:HD22	1:D:160[B]:GLN:CD	2.16	0.47
1:D:189[B]:ARG:NH1	7:D:990[B]:HOH:O	2.36	0.47
1:A:10[A]:ARG:HG3	1:A:285:ARG:HD3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:171:LEU:HB3	1:C:172:PRO:HD3	1.97	0.47
1:D:29:LEU:HD22	1:D:60:LEU:HD21	1.96	0.47
1:D:-2:HIS:HB2	7:D:1049[A]:HOH:O	2.15	0.47
1:B:183:PRO:HG3	4:B:503[B]:MPO:H21	1.97	0.46
1:B:9[B]:THR:HG22	7:B:731:HOH:O	2.13	0.46
1:A:95:THR:O	1:A:319[A]:ARG:HD2	2.15	0.46
1:D:113:LEU:HD11	6:D:504:PGE:H2	1.98	0.46
1:B:418:HIS:NE2	6:B:507[A]:PGE:H1	2.29	0.46
1:A:4[B]:ASN:O	1:A:160[B]:GLN:NE2	2.49	0.46
1:C:198:LEU:HD23	1:C:198:LEU:HA	1.83	0.46
1:B:56:PRO:CD	5:B:504:GOL:H31	2.46	0.46
1:C:285:ARG:HH12	6:C:506[A]:PGE:C6	2.29	0.46
1:C:37[A]:ASP:OD1	6:C:506[A]:PGE:H1	2.15	0.46
1:B:222:TYR:CZ	4:B:503[B]:MPO:H52	2.50	0.46
1:C:37[A]:ASP:HA	6:C:506[A]:PGE:H3	1.97	0.45
1:D:87:ILE:HG21	1:D:326[A]:ILE:HG21	1.97	0.45
1:B:36[A]:ASP:HB2	6:B:508[A]:PGE:H3	1.98	0.45
1:A:359:PHE:CZ	1:A:361:LEU:HD21	2.52	0.45
1:B:411[A]:ARG:NH1	7:B:634:HOH:O	2.50	0.45
1:A:319[A]:ARG:HG2	1:A:320:ARG:N	2.31	0.45
1:C:48:ARG:NH1	7:C:629:HOH:O	2.49	0.45
1:B:36[A]:ASP:HB2	6:B:508[A]:PGE:H42	1.97	0.45
1:D:236[A]:ASP:HB3	7:D:992[A]:HOH:O	2.17	0.45
1:B:35:ASP:HA	1:B:411[B]:ARG:NH1	2.32	0.45
6:B:508[A]:PGE:H12	7:B:1046:HOH:O	2.15	0.44
1:C:40:ASP:OD2	6:C:506[A]:PGE:H32	2.17	0.44
1:A:222:TYR:CZ	4:A:503[A]:MPO:H62	2.53	0.44
1:A:409:PRO:HG2	7:A:968[A]:HOH:O	2.16	0.44
1:A:119[B]:GLN:HG3	7:A:685[B]:HOH:O	2.17	0.43
1:B:198:LEU:HD23	1:B:198:LEU:HA	1.90	0.43
1:A:175[B]:ILE:HG21	1:A:384:LEU:HD11	2.00	0.43
1:B:137:ILE:HD13	1:B:137:ILE:HG21	1.67	0.43
1:A:161:ALA:HB1	1:A:166[B]:GLU:HG3	2.01	0.43
1:B:48[B]:ARG:NH1	7:B:618:HOH:O	2.43	0.43
1:A:-2:HIS:HD2	7:A:831[A]:HOH:O	2.02	0.43
1:B:56:PRO:HD2	5:B:504:GOL:H31	2.00	0.43
1:A:147:ARG:CZ	1:D:231[B]:ARG:HD2	2.48	0.43
1:A:255[B]:ARG:NE	7:A:629[B]:HOH:O	2.52	0.43
1:B:411[A]:ARG:HD3	6:B:507[A]:PGE:H4	2.01	0.42
1:B:278:HIS:HD2	7:B:1082[A]:HOH:O	2.02	0.42
1:A:141:VAL:HG22	1:A:175[B]:ILE:HD13	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:7[B]:ILE:HD11	1:B:284:SER:HB3	2.01	0.42
1:C:233:ASN:HB2	7:C:682[A]:HOH:O	2.20	0.42
1:B:212:PHE:CZ	3:B:502:K2B:H12C	2.55	0.42
1:D:48[B]:ARG:NH1	7:D:605[B]:HOH:O	2.31	0.42
1:C:144:ARG:NH2	7:C:609:HOH:O	2.31	0.42
1:B:34:LEU:O	1:B:411[B]:ARG:NH1	2.50	0.42
1:A:166[A]:GLU:OE1	5:A:509:GOL:H2	2.19	0.41
1:D:158[B]:ASP:OD1	1:D:158[B]:ASP:N	2.53	0.41
1:D:387[B]:ARG:HD3	7:D:1025[B]:HOH:O	2.19	0.41
1:A:381:GLY:HA3	2:A:501[B]:HEM:C3C	2.56	0.41
1:D:6[B]:ALA:HB1	7:D:850:HOH:O	2.20	0.41
1:B:175:ILE:HG21	1:B:384:LEU:HD11	2.02	0.41
1:D:229[B]:ASP:OD1	1:D:230:ARG:NH1	2.53	0.41
1:B:183:PRO:HG3	4:B:503[A]:MPO:H12	2.03	0.41
1:A:409:PRO:CG	6:A:508[B]:PGE:H42	2.50	0.40
1:C:343:TYR:CE1	1:C:372:GLY:HA2	2.57	0.40
1:A:418:HIS:NE2	6:A:508[B]:PGE:H12	2.36	0.40
1:A:4[C]:ASN:ND2	7:A:641:HOH:O	2.55	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	475/435 (109%)	465 (98%)	10 (2%)	0	100	100
1	B	474/435 (109%)	461 (97%)	13 (3%)	0	100	100
1	C	468/435 (108%)	457 (98%)	11 (2%)	0	100	100
1	D	473/435 (109%)	458 (97%)	15 (3%)	0	100	100
All	All	1890/1740 (109%)	1841 (97%)	49 (3%)	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 X-ray entries followed by that with respect to entries of similar resolution.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	406/365 (111%)	395 (97%)	11 (3%)	44	19
1	B	402/365 (110%)	394 (98%)	8 (2%)	55	32
1	C	395/365 (108%)	387 (98%)	8 (2%)	55	32
1	D	403/365 (110%)	399 (99%)	4 (1%)	76	62
All	All	1606/1460 (110%)	1575 (98%)	31 (2%)	65	34

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	105[A]	GLU
1	A	105[B]	GLU
1	A	231[A]	ARG
1	A	231[B]	ARG
1	A	235	HIS
1	A	255[A]	ARG
1	A	255[B]	ARG
1	A	260	PHE
1	A	294	TRP
1	A	319[A]	ARG
1	A	319[B]	ARG
1	B	0	HIS
1	B	105[A]	GLU
1	B	105[B]	GLU
1	B	260	PHE
1	B	294	TRP
1	B	300[A]	LEU
1	B	300[B]	LEU
1	B	328	LEU
1	C	18	GLU
1	C	35[A]	ASP
1	C	35[B]	ASP
1	C	105[A]	GLU
1	C	105[B]	GLU

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Mol	Chain	Res	Type
1	C	260	PHE
1	C	294	TRP
1	C	329	ARG
1	D	260	PHE
1	D	294	TRP
1	D	319[A]	ARG
1	D	319[B]	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

34 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
5	GOL	B	504	-	5,5,5	1.46	1 (20%)	5,5,5	1.07	1 (20%)
6	PGE	C	507[A]	-	6,6,9	0.46	0	5,5,8	0.75	0
5	GOL	B	506	-	5,5,5	1.05	0	5,5,5	0.92	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	MPO	A	503[B]	-	13,13,13	2.73	2 (15%)	17,17,17	1.84	6 (35%)
6	PGE	C	506[A]	-	9,9,9	0.38	0	8,8,8	0.75	0
5	GOL	A	506	-	5,5,5	1.09	0	5,5,5	0.88	0
6	PGE	A	508[B]	-	6,6,9	0.38	0	5,5,8	0.63	0
5	GOL	C	505	-	5,5,5	1.12	0	5,5,5	0.81	0
5	GOL	D	506[A]	-	5,5,5	1.03	0	5,5,5	1.03	0
5	GOL	B	505	-	5,5,5	1.01	0	5,5,5	0.84	0
6	PGE	D	504	-	9,9,9	0.37	0	8,8,8	0.20	0
3	K2B	A	502	-	31,31,31	0.90	0	48,48,48	1.12	3 (6%)
5	GOL	A	504	-	5,5,5	1.58	1 (20%)	5,5,5	0.74	0
2	HEM	A	501[A]	1	27,50,50	1.84	4 (14%)	17,82,82	1.84	6 (35%)
2	HEM	A	501[B]	1	27,50,50	1.82	4 (14%)	17,82,82	1.90	7 (41%)
3	K2B	B	502	-	31,31,31	0.77	1 (3%)	48,48,48	1.23	4 (8%)
3	K2B	D	502	-	31,31,31	0.88	0	48,48,48	1.22	6 (12%)
2	HEM	D	501	1	27,50,50	2.00	6 (22%)	17,82,82	2.24	9 (52%)
5	GOL	A	505	-	5,5,5	0.93	0	5,5,5	0.96	0
6	PGE	B	508[A]	-	6,6,9	0.25	0	5,5,8	0.52	0
3	K2B	C	502	-	31,31,31	0.82	0	48,48,48	1.24	5 (10%)
4	MPO	C	503[A]	-	13,13,13	2.30	2 (15%)	17,17,17	1.32	2 (11%)
5	GOL	A	507	-	5,5,5	0.95	0	5,5,5	0.96	0
4	MPO	A	503[A]	-	13,13,13	2.70	2 (15%)	17,17,17	1.28	3 (17%)
2	HEM	C	501	1	27,50,50	1.83	7 (25%)	17,82,82	2.09	6 (35%)
4	MPO	D	503[B]	-	13,13,13	2.42	2 (15%)	17,17,17	1.87	5 (29%)
4	MPO	D	503[A]	-	13,13,13	2.38	2 (15%)	17,17,17	1.48	1 (5%)
4	MPO	B	503[A]	-	13,13,13	2.73	2 (15%)	17,17,17	1.63	6 (35%)
6	PGE	B	507[A]	-	6,6,9	0.37	0	5,5,8	0.41	0
5	GOL	D	505	-	5,5,5	1.56	1 (20%)	5,5,5	0.90	0
4	MPO	B	503[B]	-	13,13,13	2.68	2 (15%)	17,17,17	1.92	5 (29%)
5	GOL	A	509	-	5,5,5	0.95	0	5,5,5	1.06	0
2	HEM	B	501	1	27,50,50	1.91	8 (29%)	17,82,82	2.26	9 (52%)
5	GOL	C	504	-	5,5,5	1.06	0	5,5,5	0.89	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	GOL	B	504	-	-	0/4/4/4	-
6	PGE	C	507[A]	-	-	2/4/4/7	-
5	GOL	B	506	-	-	4/4/4/4	-
4	MPO	A	503[B]	-	-	0/7/15/15	0/1/1/1
6	PGE	C	506[A]	-	-	4/7/7/7	-
5	GOL	A	506	-	-	2/4/4/4	-
6	PGE	A	508[B]	-	-	3/4/4/7	-
5	GOL	C	505	-	-	3/4/4/4	-
5	GOL	D	506[A]	-	-	2/4/4/4	-
5	GOL	B	505	-	-	0/4/4/4	-
6	PGE	D	504	-	-	3/7/7/7	-
3	K2B	A	502	-	-	4/10/68/68	0/4/4/4
5	GOL	A	504	-	-	0/4/4/4	-
2	HEM	A	501[A]	1	-	0/6/54/54	-
2	HEM	A	501[B]	1	-	0/6/54/54	-
3	K2B	B	502	-	-	4/10/68/68	0/4/4/4
3	K2B	D	502	-	-	4/10/68/68	0/4/4/4
2	HEM	D	501	1	-	0/6/54/54	-
5	GOL	A	505	-	-	2/4/4/4	-
6	PGE	B	508[A]	-	-	2/4/4/7	-
3	K2B	C	502	-	-	4/10/68/68	0/4/4/4
4	MPO	C	503[A]	-	-	0/7/15/15	0/1/1/1
5	GOL	A	507	-	-	2/4/4/4	-
4	MPO	A	503[A]	-	-	0/7/15/15	0/1/1/1
2	HEM	C	501	1	-	0/6/54/54	-
4	MPO	D	503[B]	-	-	1/7/15/15	0/1/1/1
4	MPO	D	503[A]	-	-	6/7/15/15	0/1/1/1
4	MPO	B	503[A]	-	-	2/7/15/15	0/1/1/1
6	PGE	B	507[A]	-	-	3/4/4/7	-
5	GOL	D	505	-	-	4/4/4/4	-
4	MPO	B	503[B]	-	-	4/7/15/15	0/1/1/1
5	GOL	A	509	-	-	2/4/4/4	-
2	HEM	B	501	1	-	0/6/54/54	-
5	GOL	C	504	-	-	2/4/4/4	-

All (47) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	503[B]	MPO	C1-S1	-7.29	1.67	1.77
4	D	503[B]	MPO	C1-S1	-7.25	1.67	1.77
4	B	503[A]	MPO	C1-S1	-7.24	1.67	1.77
4	A	503[A]	MPO	C1-S1	-7.12	1.67	1.77
4	D	503[A]	MPO	C1-S1	-7.09	1.67	1.77
4	B	503[B]	MPO	C1-S1	-7.04	1.67	1.77
4	C	503[A]	MPO	C1-S1	-6.76	1.67	1.77
4	A	503[B]	MPO	O1-S1	6.45	1.64	1.45
4	A	503[A]	MPO	O1-S1	6.44	1.64	1.45
4	B	503[B]	MPO	O1-S1	6.44	1.64	1.45
4	B	503[A]	MPO	O1-S1	6.43	1.64	1.45
2	D	501	HEM	C3C-C2C	-5.29	1.33	1.40
4	D	503[A]	MPO	O3-S1	4.69	1.64	1.47
2	A	501[B]	HEM	C3C-C2C	-4.69	1.33	1.40
4	D	503[B]	MPO	O3-S1	4.68	1.64	1.47
2	A	501[A]	HEM	C3C-C2C	-4.67	1.33	1.40
4	C	503[A]	MPO	O3-S1	4.66	1.64	1.47
2	D	501	HEM	C3B-CAB	4.34	1.56	1.47
2	D	501	HEM	C3B-C2B	-3.93	1.34	1.40
2	A	501[B]	HEM	C3B-C2B	-3.90	1.35	1.40
2	A	501[A]	HEM	C3B-C2B	-3.89	1.35	1.40
2	B	501	HEM	C3B-CAB	3.83	1.55	1.47
2	C	501	HEM	C3C-C2C	-3.76	1.35	1.40
2	A	501[B]	HEM	C3B-CAB	3.59	1.55	1.47
2	A	501[A]	HEM	C3B-CAB	3.57	1.55	1.47
2	B	501	HEM	C3C-C2C	-3.56	1.35	1.40
2	C	501	HEM	C3B-CAB	3.54	1.55	1.47
2	B	501	HEM	CAA-C2A	3.33	1.56	1.52
2	C	501	HEM	C3B-C2B	-3.31	1.35	1.40
2	C	501	HEM	CAA-C2A	3.30	1.56	1.52
5	D	505	GOL	O2-C2	-3.27	1.33	1.43
2	B	501	HEM	C3B-C2B	-3.24	1.35	1.40
5	A	504	GOL	O2-C2	-3.18	1.33	1.43
2	A	501[A]	HEM	C3C-CAC	3.17	1.54	1.47
2	D	501	HEM	C3C-CAC	3.09	1.54	1.47
2	A	501[B]	HEM	C3C-CAC	3.03	1.54	1.47
5	B	504	GOL	O2-C2	-2.89	1.34	1.43
2	C	501	HEM	CMB-C2B	2.54	1.57	1.51
2	B	501	HEM	C3C-CAC	2.54	1.53	1.47
2	C	501	HEM	C3C-CAC	2.46	1.52	1.47
2	C	501	HEM	C4D-C3D	2.43	1.48	1.42
2	B	501	HEM	CAD-C3D	2.40	1.56	1.52
2	B	501	HEM	C1A-NA	2.31	1.40	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	501	HEM	C1A-NA	2.21	1.40	1.36
2	B	501	HEM	CMB-C2B	2.15	1.56	1.51
3	B	502	K2B	C4-C3	2.03	1.50	1.45
2	D	501	HEM	CMD-C2D	2.03	1.55	1.51

All (84) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	503[B]	MPO	C5-C4-N1	4.18	116.44	110.10
4	B	503[B]	MPO	C5-C4-N1	4.09	116.31	110.10
2	B	501	HEM	C1D-C2D-C3D	3.94	109.74	107.00
2	D	501	HEM	C4A-C3A-C2A	3.93	109.73	107.00
2	B	501	HEM	C4A-C3A-C2A	3.86	109.68	107.00
4	D	503[A]	MPO	O3-S1-C1	3.86	112.00	105.77
2	A	501[A]	HEM	C1D-C2D-C3D	3.78	109.62	107.00
2	C	501	HEM	CBD-CAD-C3D	-3.73	105.60	112.48
2	A	501[B]	HEM	CBD-CAD-C3D	-3.67	105.72	112.48
2	C	501	HEM	CMD-C2D-C1D	-3.60	122.92	128.46
3	C	502	K2B	C22-C20-C17	3.60	117.72	110.28
4	A	503[B]	MPO	C6-O4-C5	3.59	121.89	109.89
2	B	501	HEM	CMD-C2D-C1D	-3.56	122.99	128.46
3	A	502	K2B	C16-C17-C20	3.42	117.44	112.15
2	D	501	HEM	CBD-CAD-C3D	-3.41	106.19	112.48
2	C	501	HEM	CBA-CAA-C2A	-3.40	106.21	112.49
2	A	501[B]	HEM	C1D-C2D-C3D	3.39	109.35	107.00
2	D	501	HEM	CMD-C2D-C1D	-3.23	123.50	128.46
2	D	501	HEM	C1D-C2D-C3D	3.22	109.23	107.00
2	C	501	HEM	C1D-C2D-C3D	3.21	109.23	107.00
2	B	501	HEM	CMB-C2B-C3B	3.17	130.61	124.68
2	B	501	HEM	CBA-CAA-C2A	-3.14	106.70	112.49
2	D	501	HEM	C4C-C3C-C2C	2.93	108.94	106.90
4	B	503[B]	MPO	O4-C5-C4	2.92	118.22	111.80
3	C	502	K2B	C7-C8-C9	-2.88	106.92	110.49
4	B	503[B]	MPO	O1-S1-C1	2.83	110.33	106.92
2	A	501[A]	HEM	CBD-CAD-C3D	-2.80	107.33	112.48
3	D	502	K2B	C16-C17-C20	2.79	116.47	112.15
4	D	503[B]	MPO	O4-C6-C7	-2.76	105.72	111.80
4	D	503[B]	MPO	O3-S1-C1	2.76	110.22	105.77
4	B	503[A]	MPO	O2-S1-C1	2.72	110.19	106.92
2	D	501	HEM	CMB-C2B-C3B	2.71	129.74	124.68
4	A	503[B]	MPO	C5-C4-N1	-2.69	106.02	110.10
4	B	503[A]	MPO	C7-N1-C4	2.68	114.86	108.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501[B]	HEM	C4C-C3C-C2C	2.64	108.74	106.90
4	D	503[B]	MPO	O1-S1-C1	2.64	110.09	106.92
4	A	503[A]	MPO	O2-S1-C1	2.60	110.05	106.92
3	B	502	K2B	C2-C3-C4	2.60	120.75	116.74
4	A	503[B]	MPO	C2-C1-S1	-2.59	109.27	113.25
4	B	503[B]	MPO	O2-S1-C1	2.58	110.02	106.92
3	D	502	K2B	C22-C20-C17	2.58	115.61	110.28
2	B	501	HEM	CMA-C3A-C4A	-2.56	124.52	128.46
2	C	501	HEM	C4C-C3C-C2C	2.56	108.69	106.90
2	D	501	HEM	C3C-C4C-NC	-2.49	106.25	110.94
2	D	501	HEM	CBA-CAA-C2A	-2.47	107.93	112.49
3	B	502	K2B	C1-C2-C3	-2.47	106.34	111.62
2	A	501[A]	HEM	C4C-C3C-C2C	2.46	108.62	106.90
4	A	503[B]	MPO	O2-S1-C1	2.46	109.88	106.92
3	A	502	K2B	C22-C20-C17	2.45	115.34	110.28
3	D	502	K2B	C1-C2-C3	-2.42	106.44	111.62
2	A	501[A]	HEM	CMD-C2D-C1D	-2.41	124.76	128.46
2	C	501	HEM	CMB-C2B-C3B	2.41	129.19	124.68
3	D	502	K2B	C11-C9-C8	2.41	115.22	111.75
4	B	503[A]	MPO	O1-S1-C1	2.37	109.77	106.92
4	C	503[A]	MPO	O2-S1-C1	2.34	109.73	106.92
3	B	502	K2B	C22-C20-C17	2.34	115.12	110.28
2	A	501[B]	HEM	CMD-C2D-C1D	-2.30	124.92	128.46
3	C	502	K2B	C12-C13-C14	2.30	110.83	107.27
2	A	501[B]	HEM	CMB-C2B-C3B	2.28	128.94	124.68
4	A	503[B]	MPO	C2-C3-N1	-2.28	108.10	113.84
4	A	503[B]	MPO	O3-S1-C1	2.25	109.40	105.77
4	C	503[A]	MPO	O1-S1-C1	2.24	109.62	106.92
2	D	501	HEM	CMA-C3A-C4A	-2.23	125.04	128.46
5	B	504	GOL	C3-C2-C1	-2.18	103.21	111.70
2	A	501[B]	HEM	CBA-CAA-C2A	-2.17	108.48	112.49
4	B	503[A]	MPO	C5-C4-N1	2.17	113.39	110.10
2	A	501[A]	HEM	CMB-C2B-C3B	2.15	128.71	124.68
3	D	502	K2B	C7-C6-C5	-2.15	107.89	111.93
2	B	501	HEM	C4C-C3C-C2C	2.14	108.39	106.90
4	B	503[A]	MPO	C6-C7-N1	2.14	113.35	110.10
2	A	501[A]	HEM	CMA-C3A-C4A	-2.14	125.18	128.46
4	A	503[A]	MPO	O3-S1-C1	2.14	109.22	105.77
4	B	503[B]	MPO	C6-C7-N1	-2.13	106.88	110.10
3	C	502	K2B	C21-C20-C17	-2.12	109.67	112.92
4	D	503[B]	MPO	C7-N1-C4	2.10	113.56	108.83
2	B	501	HEM	CBD-CAD-C3D	-2.10	108.62	112.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501[B]	HEM	CMA-C3A-C4A	-2.07	125.28	128.46
4	A	503[A]	MPO	O1-S1-C1	2.06	109.39	106.92
3	C	502	K2B	C16-C17-C20	2.05	115.32	112.15
3	B	502	K2B	C16-C17-C20	2.05	115.32	112.15
4	B	503[A]	MPO	O4-C6-C7	-2.03	107.33	111.80
3	D	502	K2B	C6-C5-C4	-2.03	117.48	120.87
3	A	502	K2B	C6-C7-C8	-2.03	108.02	111.69
2	B	501	HEM	C3C-C4C-NC	-2.02	107.12	110.94

There are no chirality outliers.

All (69) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	506	GOL	O1-C1-C2-O2
5	B	506	GOL	O1-C1-C2-C3
5	A	506	GOL	O1-C1-C2-C3
4	D	503[A]	MPO	C2-C1-S1-O1
4	D	503[A]	MPO	C2-C1-S1-O3
5	D	505	GOL	O1-C1-C2-C3
5	D	505	GOL	C1-C2-C3-O3
5	A	509	GOL	O1-C1-C2-C3
5	C	504	GOL	O1-C1-C2-C3
4	B	503[A]	MPO	C2-C3-N1-C4
4	B	503[B]	MPO	C2-C3-N1-C4
6	C	506[A]	PGE	C6-C5-O3-C4
3	A	502	K2B	C13-C17-C20-C22
3	D	502	K2B	C13-C17-C20-C22
3	C	502	K2B	C13-C17-C20-C22
3	B	502	K2B	C13-C17-C20-C22
6	D	504	PGE	O2-C3-C4-O3
3	A	502	K2B	C16-C17-C20-C21
3	D	502	K2B	C16-C17-C20-C21
3	C	502	K2B	C16-C17-C20-C21
5	A	505	GOL	O2-C2-C3-O3
5	A	507	GOL	O2-C2-C3-O3
4	B	503[A]	MPO	C2-C3-N1-C7
3	B	502	K2B	C16-C17-C20-C21
3	A	502	K2B	C13-C17-C20-C21
6	C	506[A]	PGE	O2-C3-C4-O3
3	B	502	K2B	C13-C17-C20-C21
3	D	502	K2B	C13-C17-C20-C21
3	C	502	K2B	C13-C17-C20-C21

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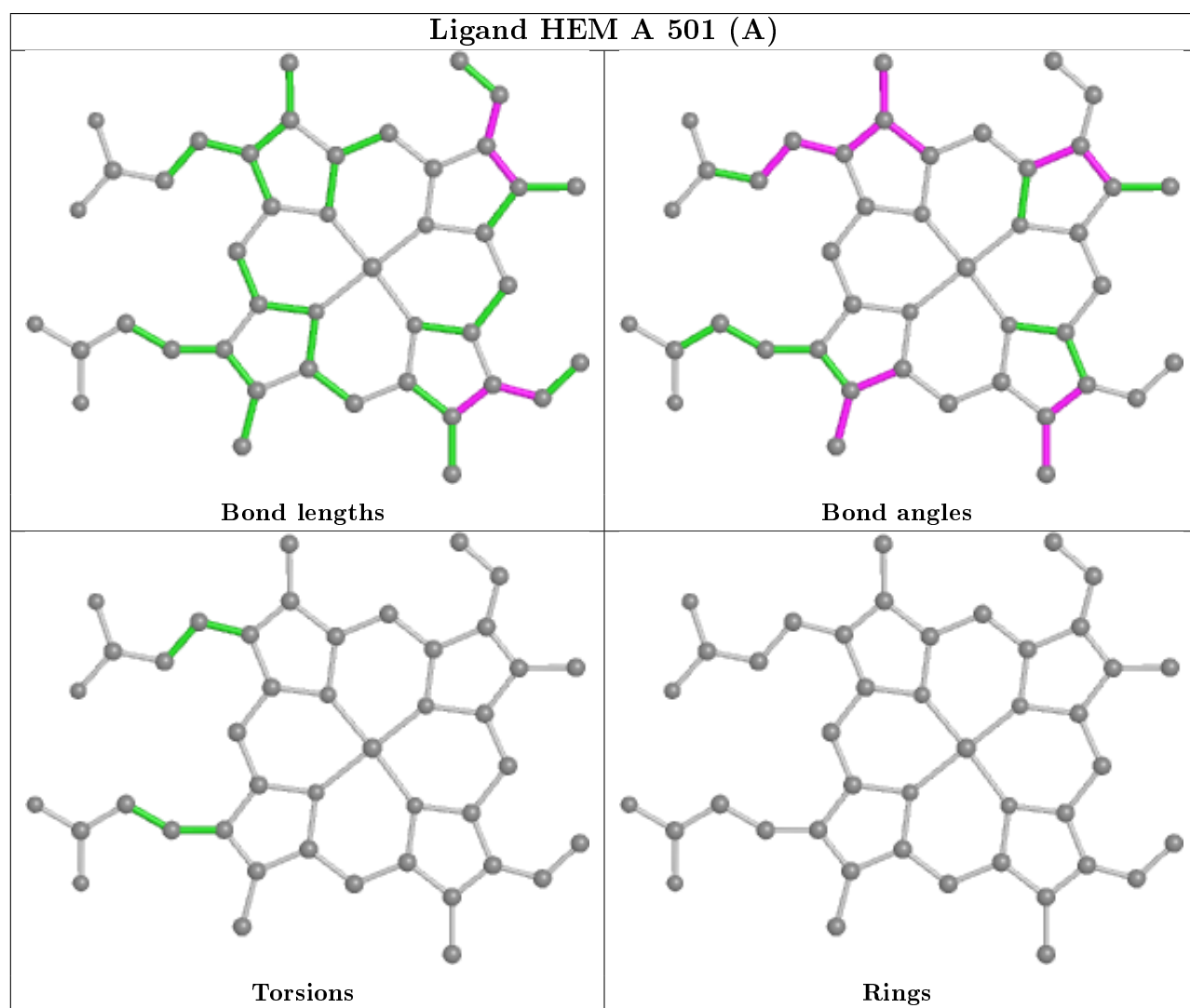
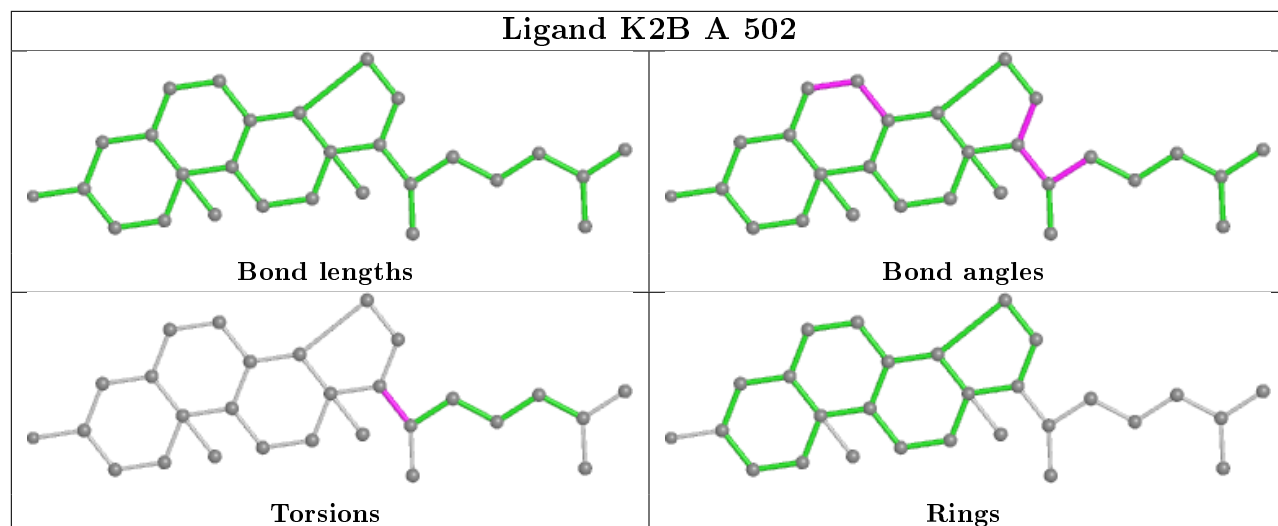
Mol	Chain	Res	Type	Atoms
6	D	504	PGE	O1-C1-C2-O2
6	B	508[A]	PGE	O1-C1-C2-O2
6	B	507[A]	PGE	O1-C1-C2-O2
5	B	506	GOL	C1-C2-C3-O3
5	C	505	GOL	C1-C2-C3-O3
5	D	506[A]	GOL	O1-C1-C2-C3
5	A	505	GOL	C1-C2-C3-O3
5	A	507	GOL	C1-C2-C3-O3
6	C	507[A]	PGE	O3-C5-C6-O4
4	B	503[B]	MPO	C1-C2-C3-N1
5	A	506	GOL	O1-C1-C2-O2
5	D	505	GOL	O1-C1-C2-O2
5	A	509	GOL	O1-C1-C2-O2
6	C	506[A]	PGE	O3-C5-C6-O4
6	B	508[A]	PGE	O2-C3-C4-O3
5	B	506	GOL	O2-C2-C3-O3
5	C	504	GOL	O1-C1-C2-O2
6	C	507[A]	PGE	O2-C3-C4-O3
4	D	503[A]	MPO	C2-C1-S1-O2
4	B	503[B]	MPO	C2-C1-S1-O2
3	B	502	K2B	C16-C17-C20-C22
3	C	502	K2B	C16-C17-C20-C22
3	A	502	K2B	C16-C17-C20-C22
6	C	506[A]	PGE	O1-C1-C2-O2
3	D	502	K2B	C16-C17-C20-C22
6	B	507[A]	PGE	C1-C2-O2-C3
6	D	504	PGE	C1-C2-O2-C3
5	D	505	GOL	O2-C2-C3-O3
6	A	508[B]	PGE	C4-C3-O2-C2
4	D	503[A]	MPO	C2-C3-N1-C4
4	D	503[B]	MPO	C1-C2-C3-N1
4	D	503[A]	MPO	C1-C2-C3-N1
6	A	508[B]	PGE	O2-C3-C4-O3
6	A	508[B]	PGE	C1-C2-O2-C3
6	B	507[A]	PGE	C4-C3-O2-C2
4	D	503[A]	MPO	C2-C3-N1-C7
5	C	505	GOL	O2-C2-C3-O3
4	B	503[B]	MPO	C2-C1-S1-O3
5	D	506[A]	GOL	O1-C1-C2-O2
5	C	505	GOL	O1-C1-C2-C3

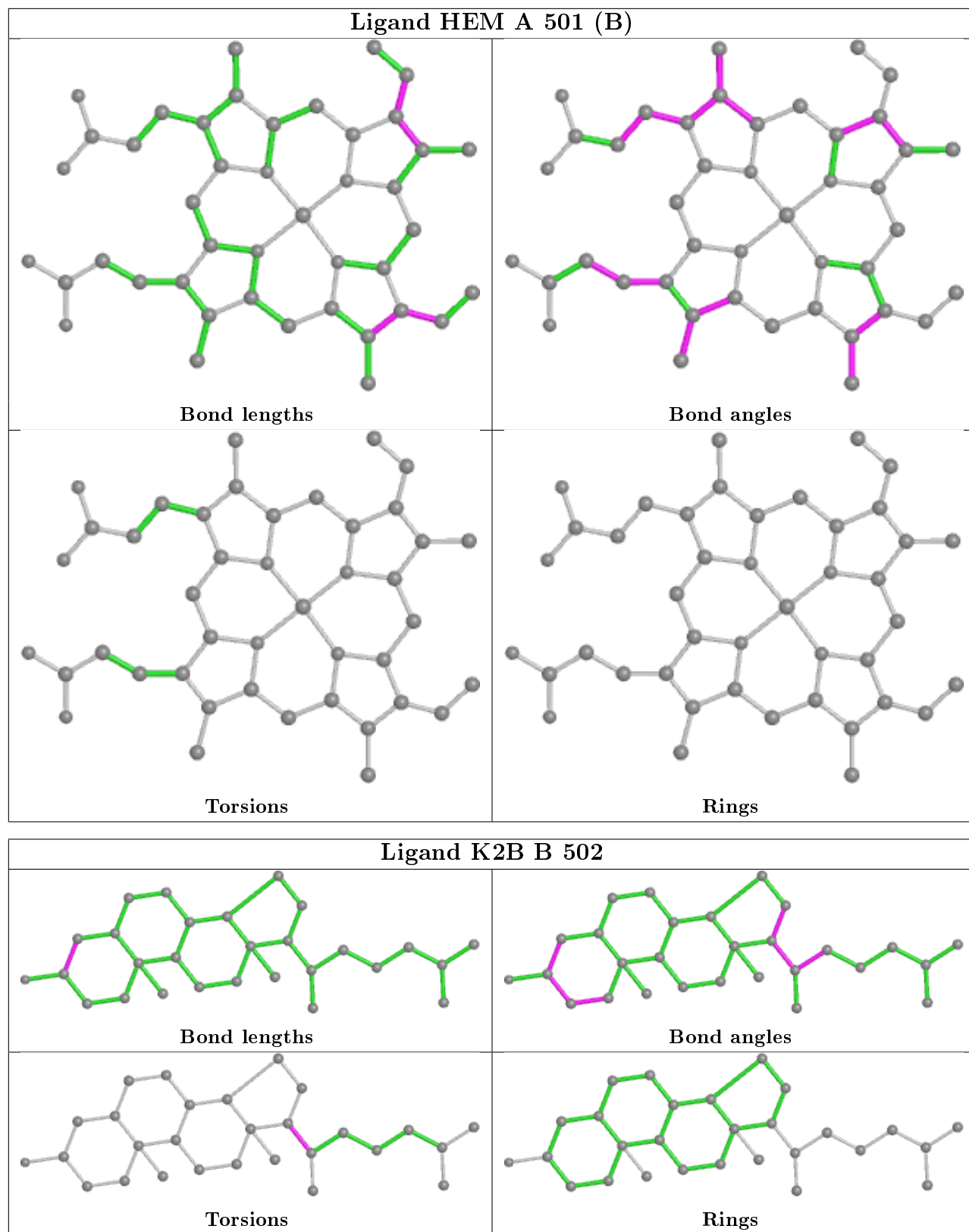
There are no ring outliers.

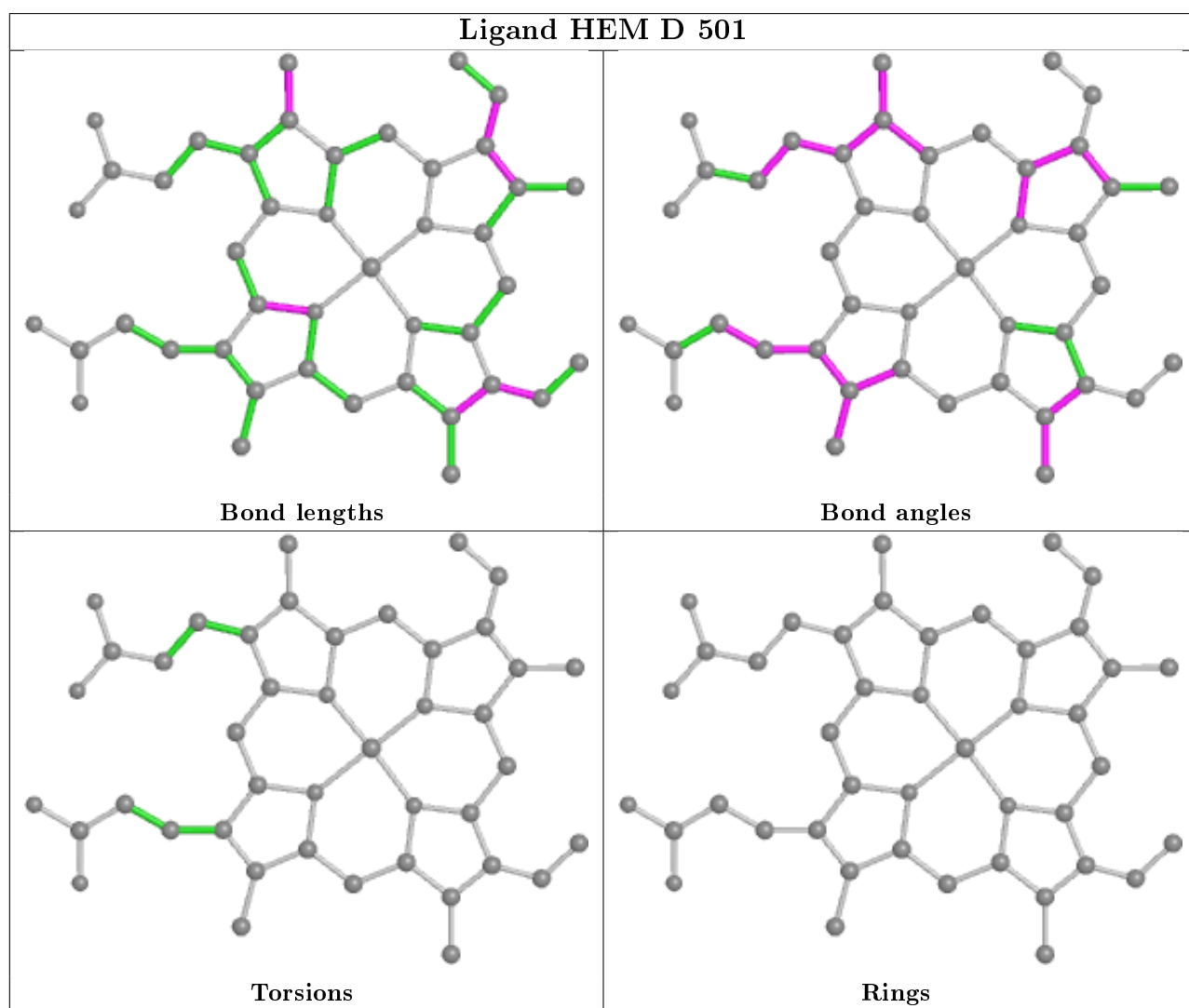
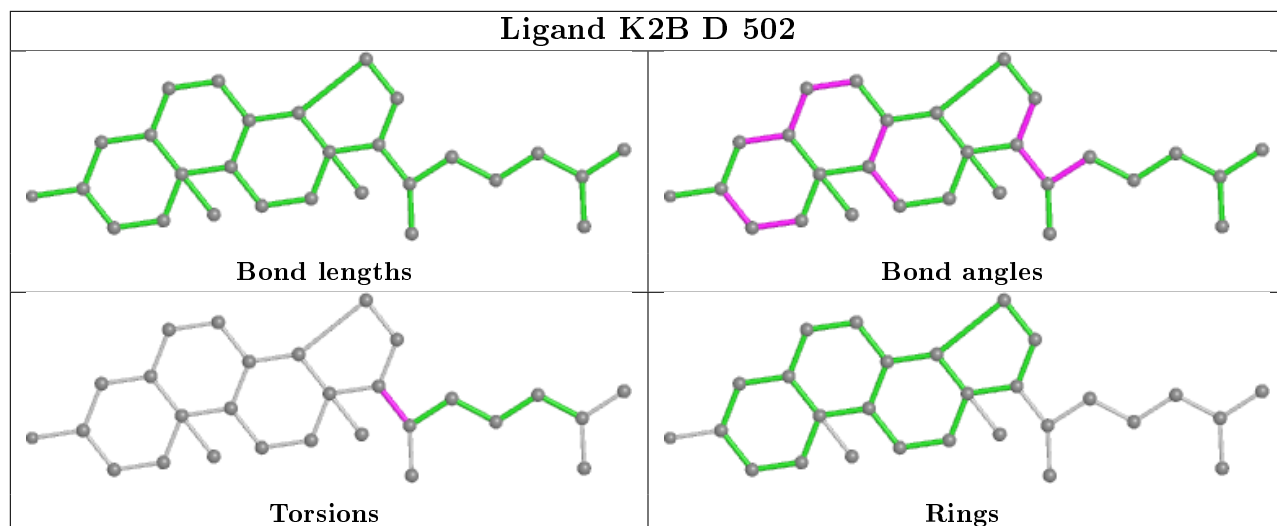
19 monomers are involved in 50 short contacts:

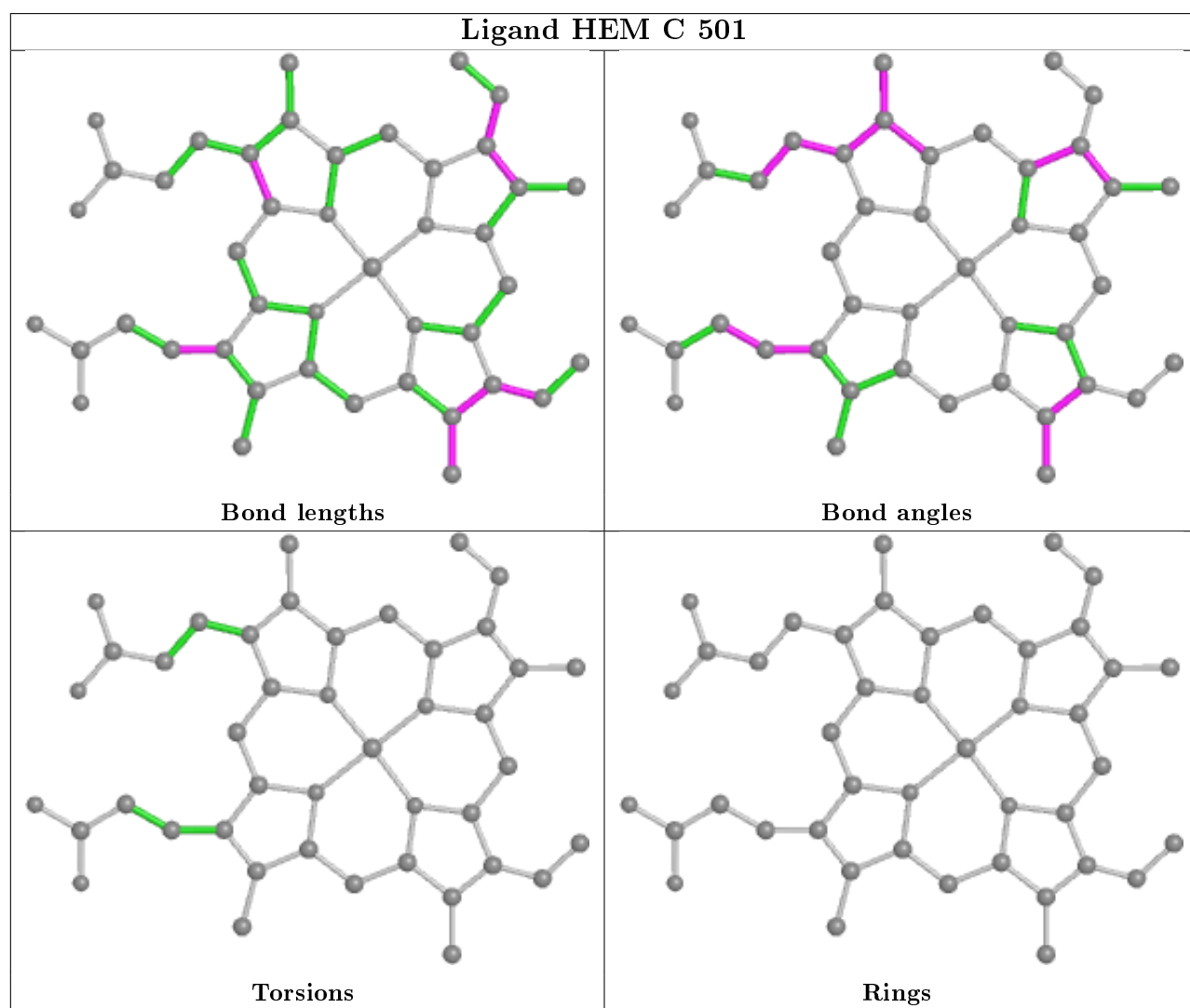
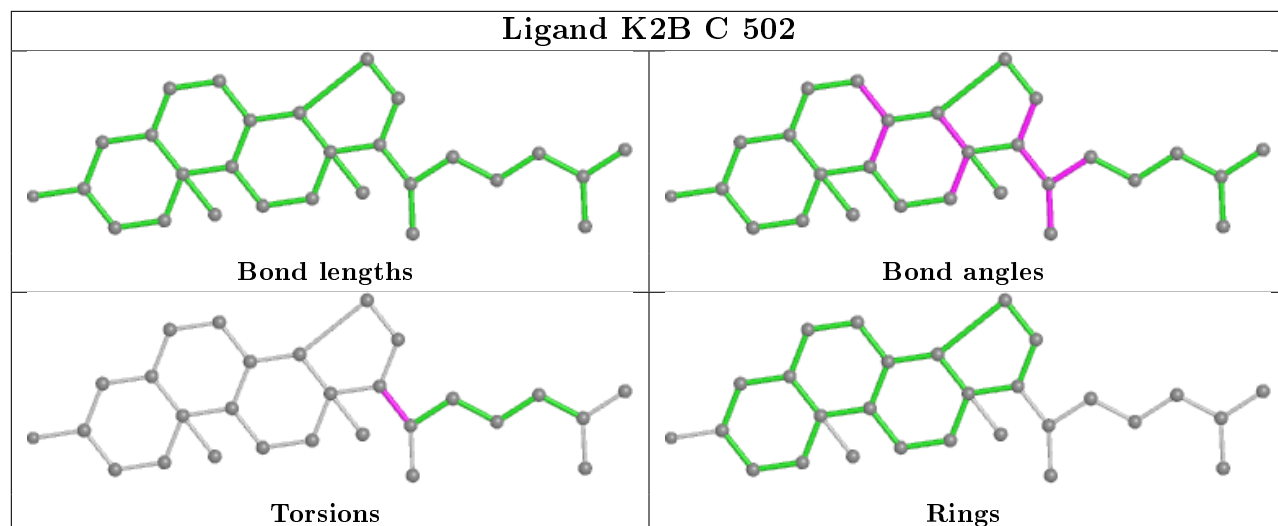
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	504	GOL	2	0
6	C	507[A]	PGE	1	0
5	B	506	GOL	1	0
6	C	506[A]	PGE	11	0
6	A	508[B]	PGE	10	0
5	D	506[A]	GOL	1	0
6	D	504	PGE	2	0
3	A	502	K2B	1	0
2	A	501[B]	HEM	2	0
3	B	502	K2B	1	0
5	A	505	GOL	1	0
6	B	508[A]	PGE	4	0
4	A	503[A]	MPO	1	0
4	D	503[A]	MPO	1	0
4	B	503[A]	MPO	1	0
6	B	507[A]	PGE	4	0
4	B	503[B]	MPO	4	0
5	A	509	GOL	1	0
2	B	501	HEM	1	0

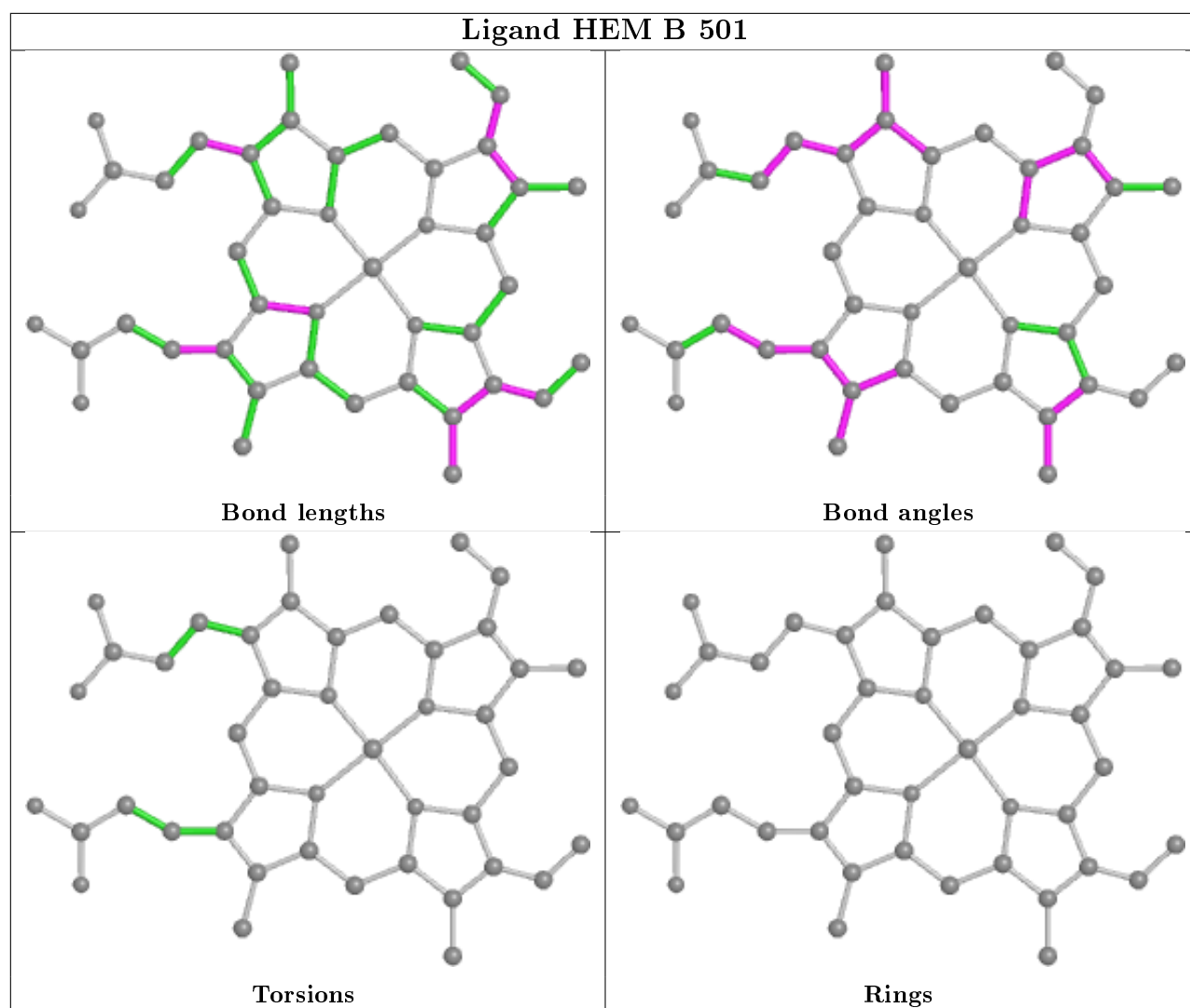
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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











5.7 Other polymers [\(i\)](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [\(i\)](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	431/435 (99%)	-0.89	2 (0%) 91 92	17, 25, 38, 90	1 (0%)
1	B	430/435 (98%)	-0.86	4 (0%) 84 86	18, 25, 41, 89	2 (0%)
1	C	430/435 (98%)	-0.88	3 (0%) 87 89	16, 23, 38, 83	3 (0%)
1	D	431/435 (99%)	-0.94	1 (0%) 95 95	16, 23, 36, 88	1 (0%)
All	All	1722/1740 (98%)	-0.89	10 (0%) 89 90	16, 24, 39, 90	7 (0%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	1[A]	MET	5.2
1	B	-1	HIS	4.3
1	C	-1	HIS	4.2
1	D	-2	HIS	3.1
1	C	0	HIS	3.0
1	A	-2	HIS	2.9
1	A	134[A]	VAL	2.5
1	B	0	HIS	2.3
1	B	4[A]	ASN	2.2
1	B	3[A]	LEU	2.1

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

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

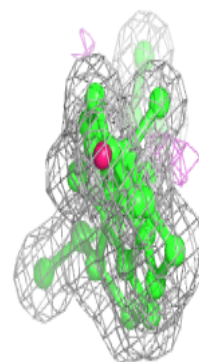
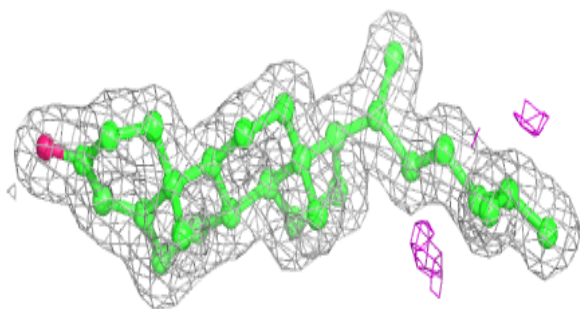
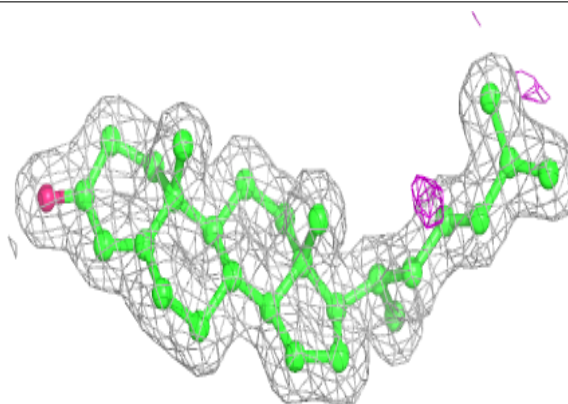
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	GOL	A	506	6/6	0.31	0.20	85,86,87,88	0
5	GOL	C	505	6/6	0.35	0.32	76,78,78,78	0
5	GOL	A	507	6/6	0.36	0.26	94,96,97,98	0
5	GOL	A	505	6/6	0.64	0.18	74,74,76,76	0
5	GOL	B	505	6/6	0.68	0.18	80,81,81,81	0
6	PGE	B	507[A]	7/10	0.71	0.32	40,47,51,51	7
5	GOL	A	509	6/6	0.72	0.11	67,69,69,69	0
6	PGE	C	507[A]	7/10	0.76	0.22	30,33,35,36	7
5	GOL	C	504	6/6	0.80	0.20	64,70,71,71	0
5	GOL	B	506	6/6	0.83	0.29	71,76,78,80	0
6	PGE	B	508[A]	7/10	0.83	0.25	31,34,40,42	7
6	PGE	D	504	10/10	0.83	0.23	43,63,68,69	0
6	PGE	A	508[B]	7/10	0.85	0.25	25,30,40,43	7
6	PGE	C	506[A]	10/10	0.85	0.24	21,32,35,37	10
4	MPO	D	503[A]	13/13	0.88	0.18	52,53,57,57	13
4	MPO	D	503[B]	13/13	0.88	0.18	52,53,57,57	13
4	MPO	A	503[A]	13/13	0.89	0.16	38,40,47,47	13
5	GOL	D	505	6/6	0.89	0.18	37,51,55,59	0
5	GOL	D	506[A]	6/6	0.89	0.20	31,35,38,43	6
4	MPO	A	503[B]	13/13	0.89	0.16	39,40,47,47	13
4	MPO	B	503[A]	13/13	0.91	0.18	38,40,44,44	13
4	MPO	B	503[B]	13/13	0.91	0.18	38,40,43,44	13
5	GOL	A	504	6/6	0.92	0.21	37,54,57,59	0
5	GOL	B	504	6/6	0.94	0.15	36,50,55,57	0
4	MPO	C	503[A]	13/13	0.95	0.10	35,40,47,48	13
3	K2B	C	502	28/28	0.98	0.04	16,18,22,23	0
3	K2B	D	502	28/28	0.98	0.05	14,18,20,21	0
3	K2B	A	502	28/28	0.98	0.05	15,18,21,22	0
3	K2B	B	502	28/28	0.98	0.05	17,19,24,27	0
2	HEM	A	501[A]	43/43	0.99	0.05	17,19,21,22	43
2	HEM	A	501[B]	43/43	0.99	0.05	17,19,21,23	43
2	HEM	D	501	43/43	0.99	0.05	14,17,19,22	0
2	HEM	B	501	43/43	0.99	0.05	15,18,20,22	0
2	HEM	C	501	43/43	0.99	0.05	14,17,19,21	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

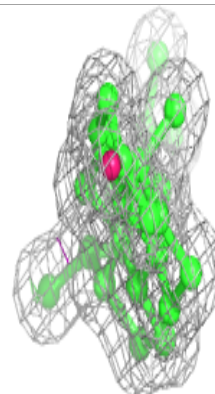
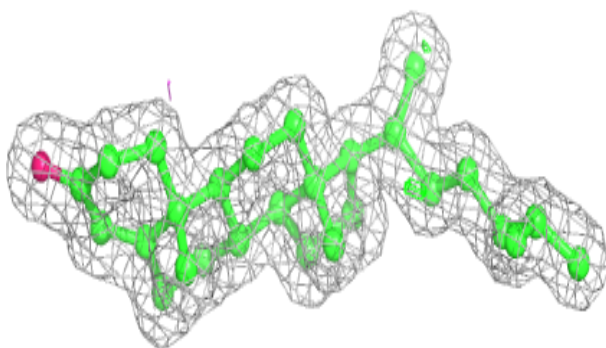
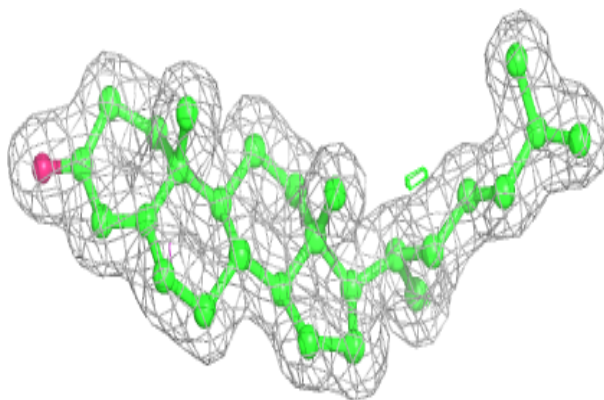
Electron density around K2B C 502:

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



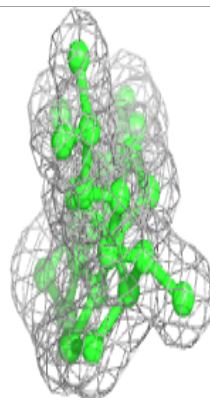
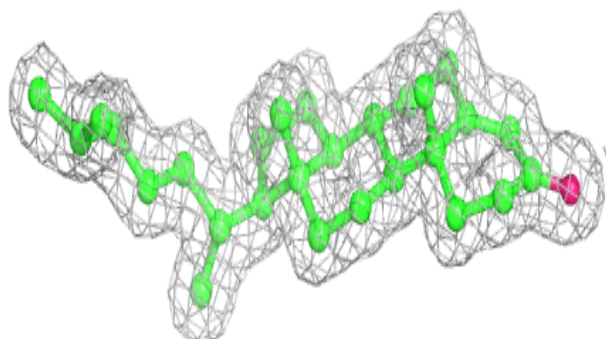
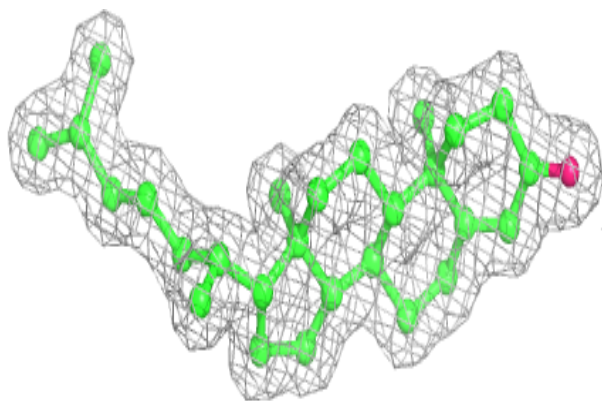
Electron density around K2B D 502:

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

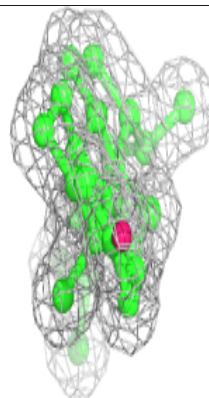
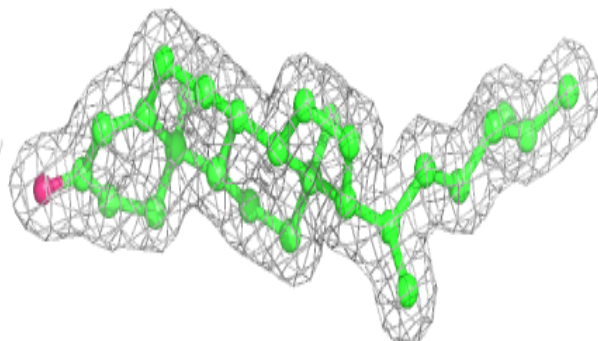
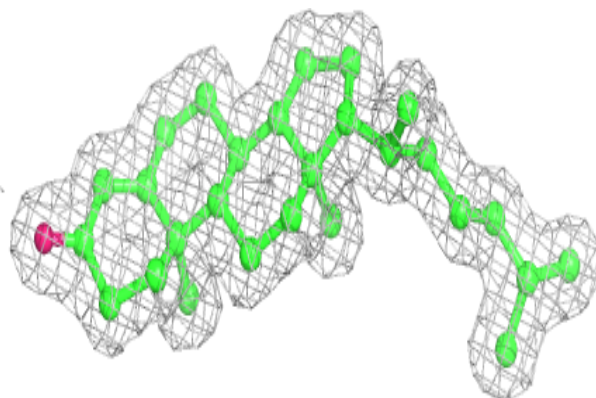


Electron density around K2B A 502:

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

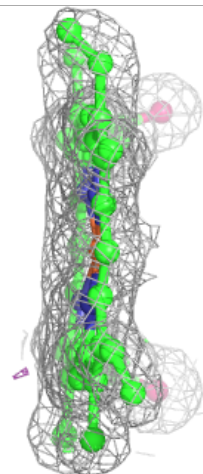
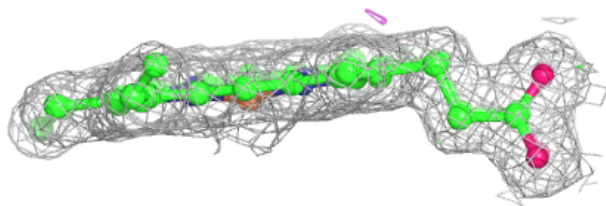
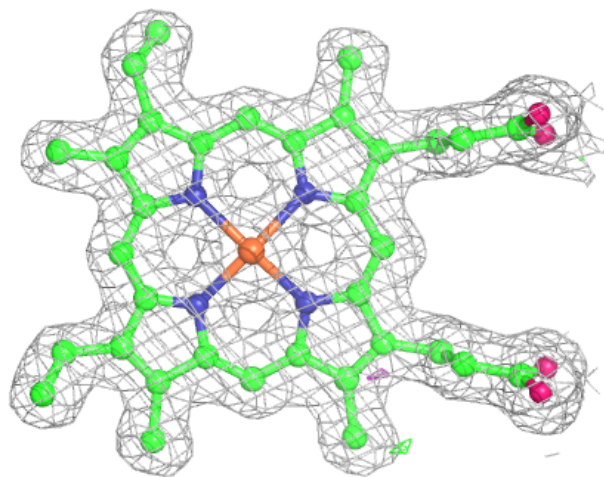
**Electron density around K2B B 502:**

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



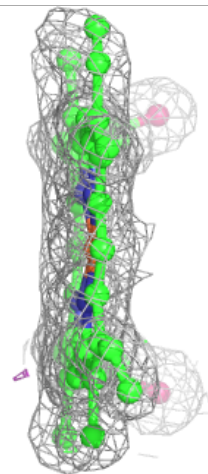
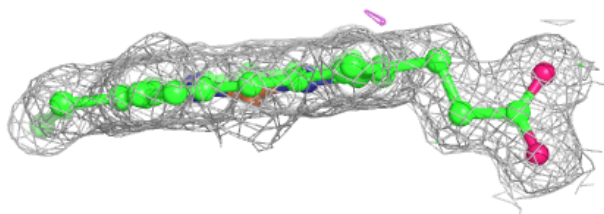
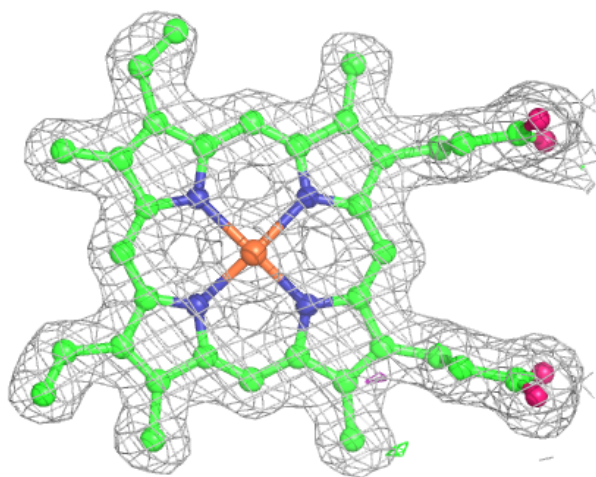
Electron density around HEM A 501 (A):

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



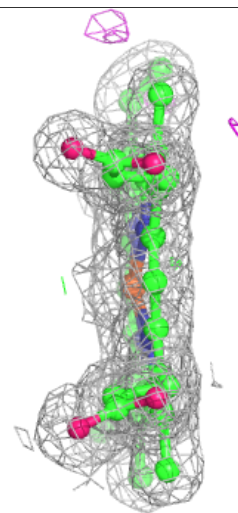
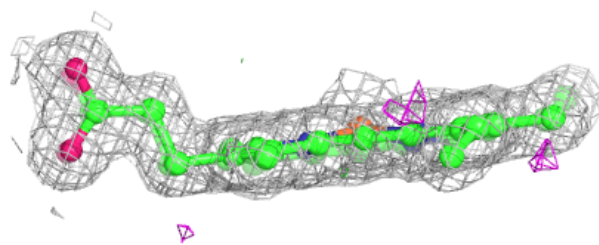
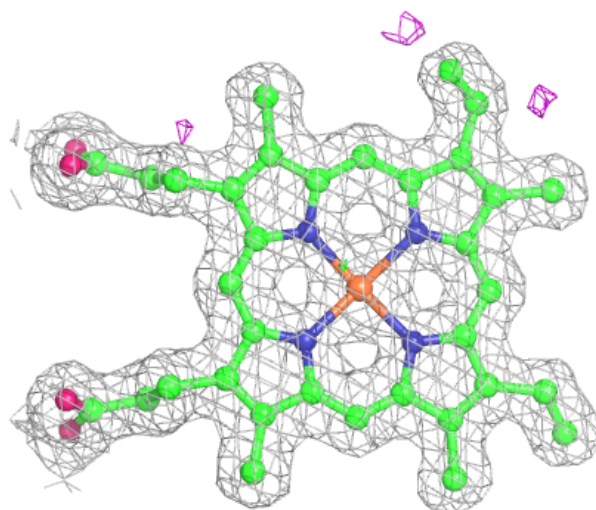
Electron density around HEM A 501 (B):

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



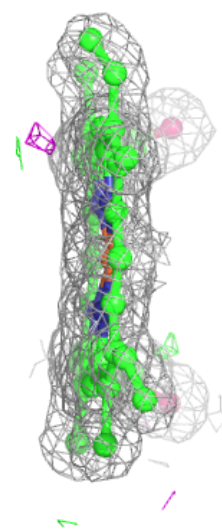
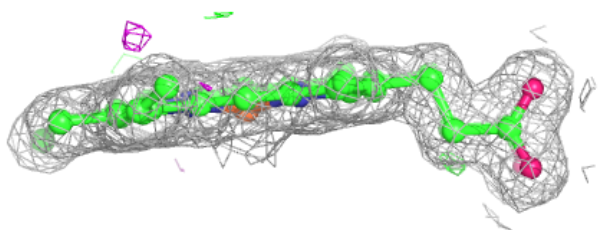
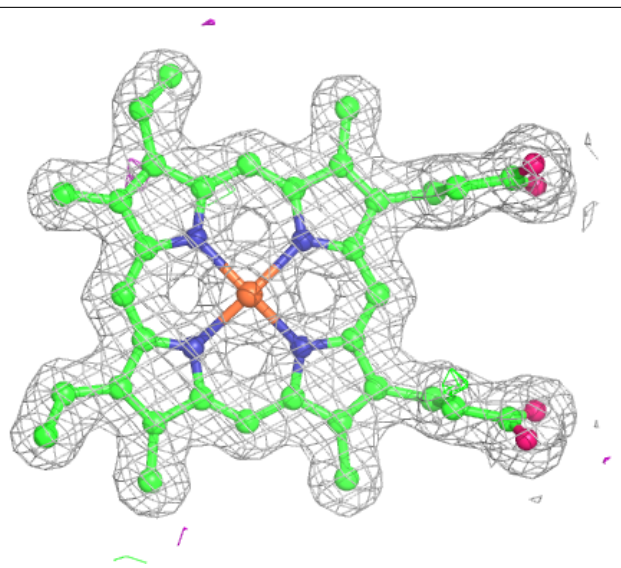
Electron density around HEM D 501:

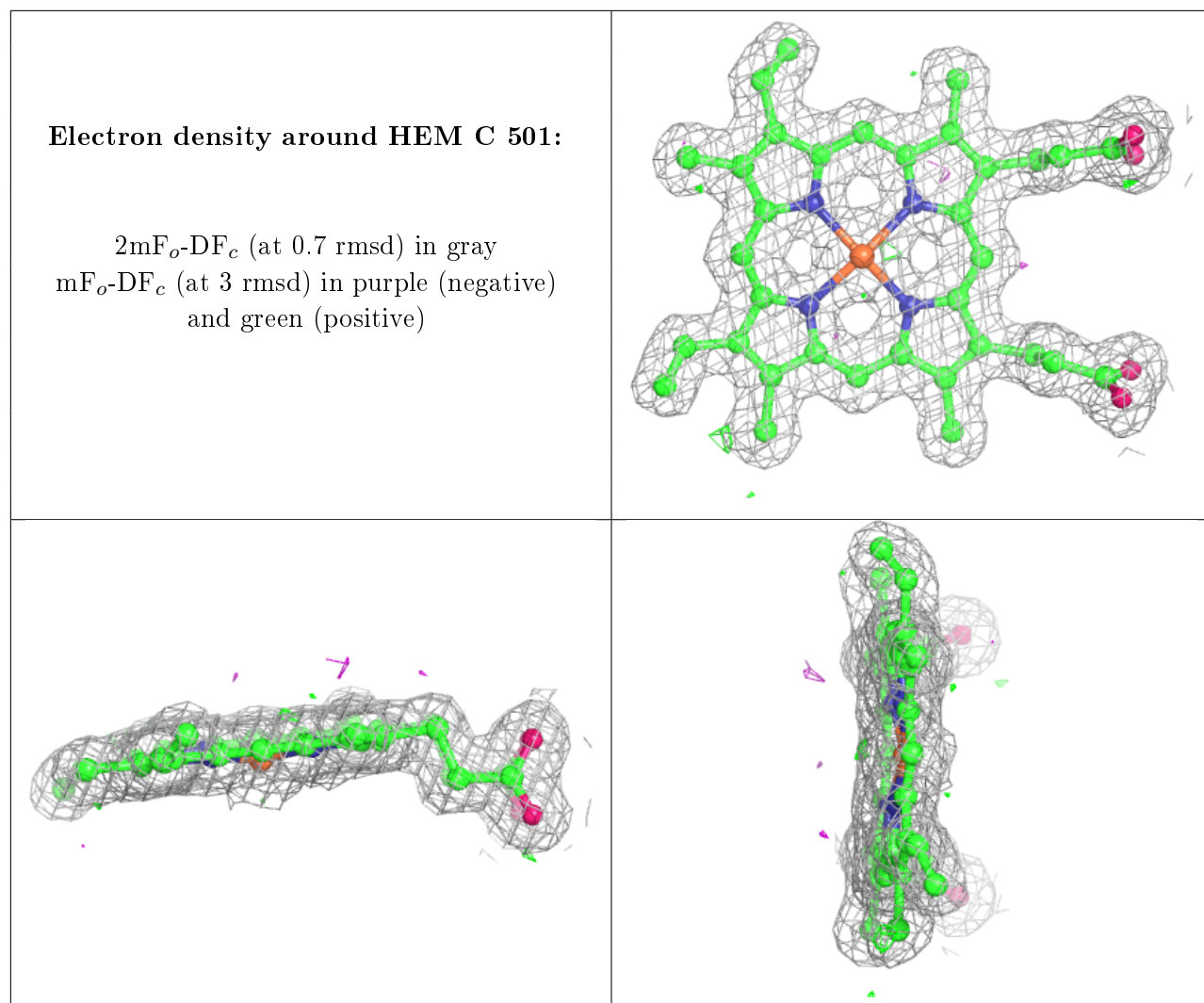
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM B 501:

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





6.5 Other polymers [i](#)

There are no such residues in this entry.