



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 4, 2026 – 09:07 AM EDT

PDB ID : 3T3Z / pdb_00003t3z
Title : Human Cytochrome P450 2E1 in complex with pilocarpine
Authors : Meneely, K.M.; DeVore, N.M.; Scott, E.E.
Deposited on : 2011-07-25
Resolution : 2.35 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 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:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

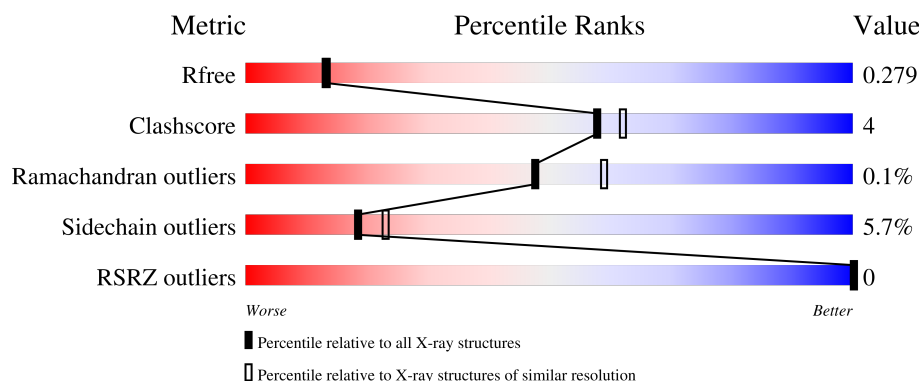
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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.



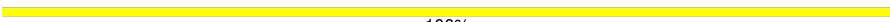
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	476	
1	B	476	
1	C	476	
1	D	476	
2	E	2	

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Mol	Chain	Length	Quality of chain
2	F	2	 100%
2	G	2	 100%
2	H	2	 100%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 16204 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a protein called Cytochrome P450 2E1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	464	Total	C	N	O	S	0	2	0
			3799	2458	651	672	18			
1	B	464	Total	C	N	O	S	0	1	0
			3793	2454	651	670	18			
1	C	464	Total	C	N	O	S	0	0	0
			3787	2450	651	668	18			
1	D	464	Total	C	N	O	S	0	0	0
			3787	2450	651	668	18			

There are 56 discrepancies between the modelled and reference sequences:

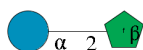
Chain	Residue	Modelled	Actual	Comment	Reference
A	22	MET	-	expression tag	UNP P05181
A	23	ALA	-	expression tag	UNP P05181
A	24	LYS	-	expression tag	UNP P05181
A	25	LYS	-	expression tag	UNP P05181
A	26	THR	-	expression tag	UNP P05181
A	27	SER	-	expression tag	UNP P05181
A	28	SER	-	expression tag	UNP P05181
A	29	LYS	-	expression tag	UNP P05181
A	30	GLY	-	expression tag	UNP P05181
A	31	LYS	-	expression tag	UNP P05181
A	494	HIS	-	expression tag	UNP P05181
A	495	HIS	-	expression tag	UNP P05181
A	496	HIS	-	expression tag	UNP P05181
A	497	HIS	-	expression tag	UNP P05181
B	22	MET	-	expression tag	UNP P05181
B	23	ALA	-	expression tag	UNP P05181
B	24	LYS	-	expression tag	UNP P05181
B	25	LYS	-	expression tag	UNP P05181
B	26	THR	-	expression tag	UNP P05181
B	27	SER	-	expression tag	UNP P05181
B	28	SER	-	expression tag	UNP P05181

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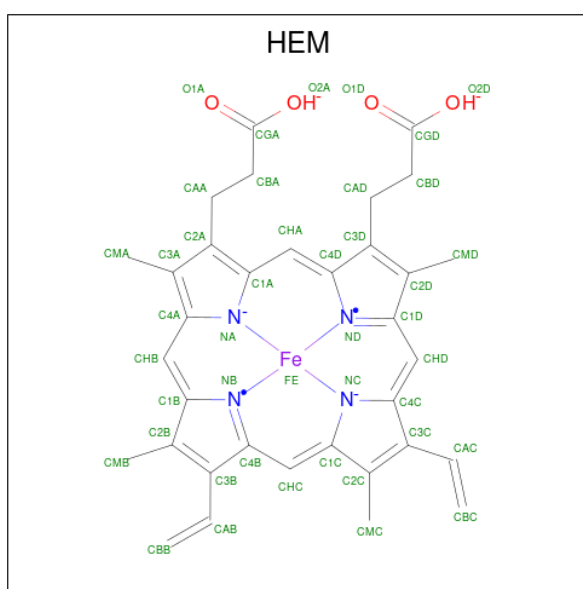
Chain	Residue	Modelled	Actual	Comment	Reference
B	29	LYS	-	expression tag	UNP P05181
B	30	GLY	-	expression tag	UNP P05181
B	31	LYS	-	expression tag	UNP P05181
B	494	HIS	-	expression tag	UNP P05181
B	495	HIS	-	expression tag	UNP P05181
B	496	HIS	-	expression tag	UNP P05181
B	497	HIS	-	expression tag	UNP P05181
C	22	MET	-	expression tag	UNP P05181
C	23	ALA	-	expression tag	UNP P05181
C	24	LYS	-	expression tag	UNP P05181
C	25	LYS	-	expression tag	UNP P05181
C	26	THR	-	expression tag	UNP P05181
C	27	SER	-	expression tag	UNP P05181
C	28	SER	-	expression tag	UNP P05181
C	29	LYS	-	expression tag	UNP P05181
C	30	GLY	-	expression tag	UNP P05181
C	31	LYS	-	expression tag	UNP P05181
C	494	HIS	-	expression tag	UNP P05181
C	495	HIS	-	expression tag	UNP P05181
C	496	HIS	-	expression tag	UNP P05181
C	497	HIS	-	expression tag	UNP P05181
D	22	MET	-	expression tag	UNP P05181
D	23	ALA	-	expression tag	UNP P05181
D	24	LYS	-	expression tag	UNP P05181
D	25	LYS	-	expression tag	UNP P05181
D	26	THR	-	expression tag	UNP P05181
D	27	SER	-	expression tag	UNP P05181
D	28	SER	-	expression tag	UNP P05181
D	29	LYS	-	expression tag	UNP P05181
D	30	GLY	-	expression tag	UNP P05181
D	31	LYS	-	expression tag	UNP P05181
D	494	HIS	-	expression tag	UNP P05181
D	495	HIS	-	expression tag	UNP P05181
D	496	HIS	-	expression tag	UNP P05181
D	497	HIS	-	expression tag	UNP P05181

- Molecule 2 is an oligosaccharide called beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose.



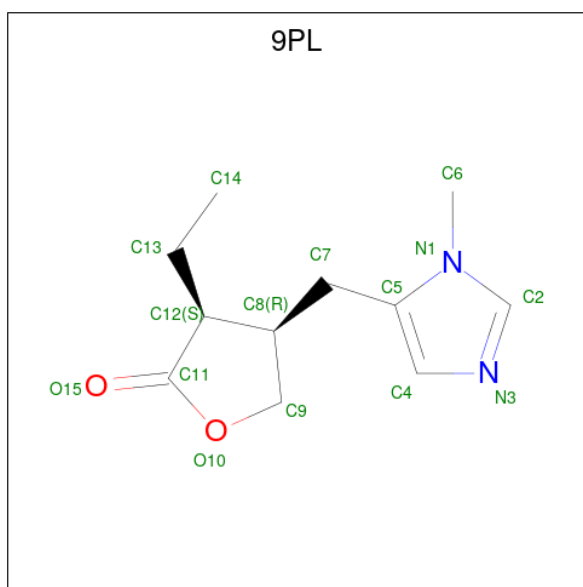
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	E	2	Total 23	C 12	O 11	0	0	0
2	F	2	Total 23	C 12	O 11	0	0	0
2	G	2	Total 23	C 12	O 11	0	0	0
2	H	2	Total 23	C 12	O 11	0	0	0

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
3	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 4 is (3S,4R)-3-ethyl-4-[(1-methyl-1H-imidazol-5-yl)methyl]dihydrofuran-2(3H)-one (CCD ID: 9PL) (formula: C₁₁H₁₆N₂O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			15	11	2	2		
4	B	1	Total	C	N	O	0	0
			15	11	2	2		
4	C	1	Total	C	N	O	0	0
			15	11	2	2		
4	D	1	Total	C	N	O	0	0
			15	11	2	2		

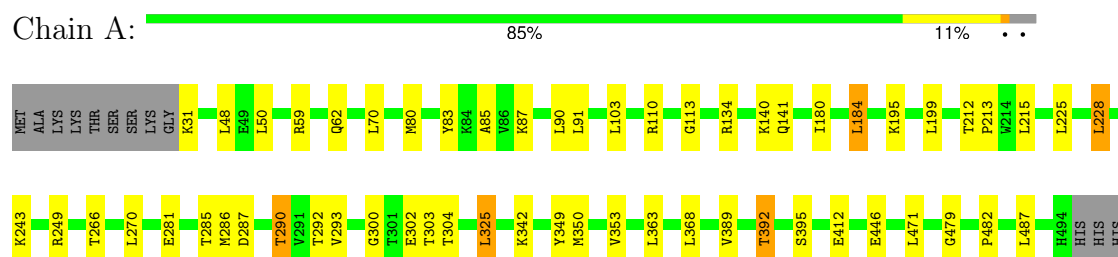
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	202	Total	O	0	0
			202	202		
5	B	202	Total	O	0	0
			202	202		
5	C	151	Total	O	0	0
			151	151		
5	D	159	Total	O	0	0
			159	159		

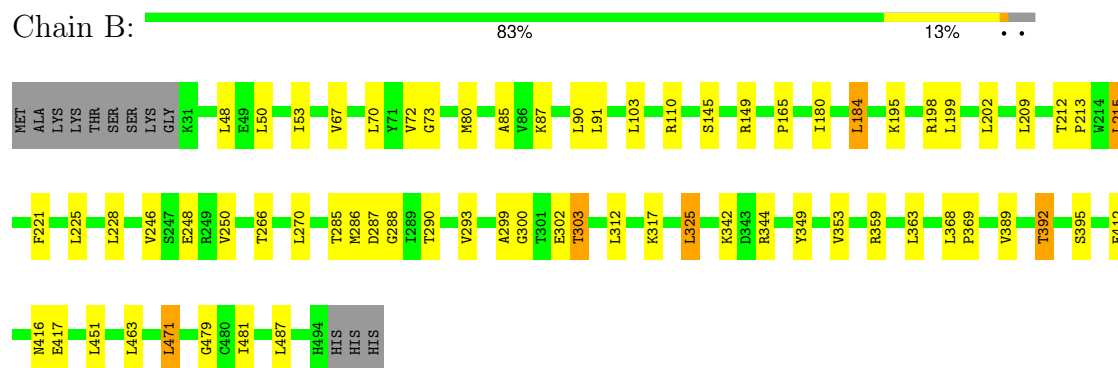
3 Residue-property plots

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

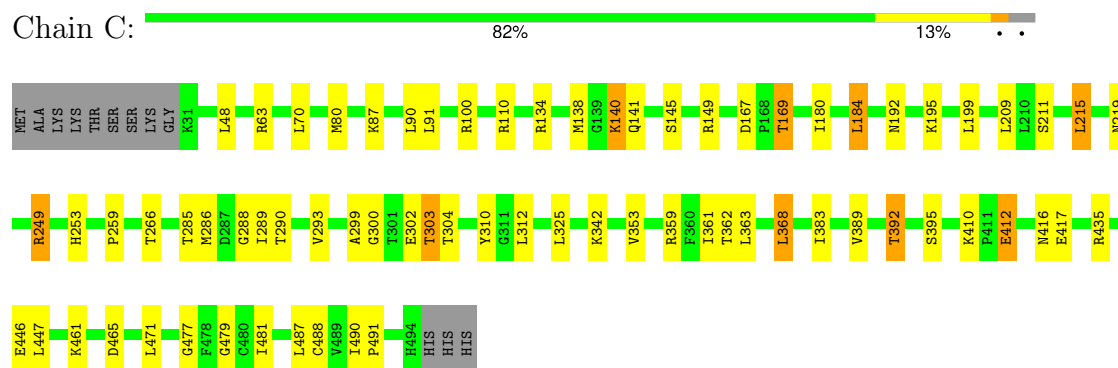
• Molecule 1: Cytochrome P450 2E1



• Molecule 1: Cytochrome P450 2E1

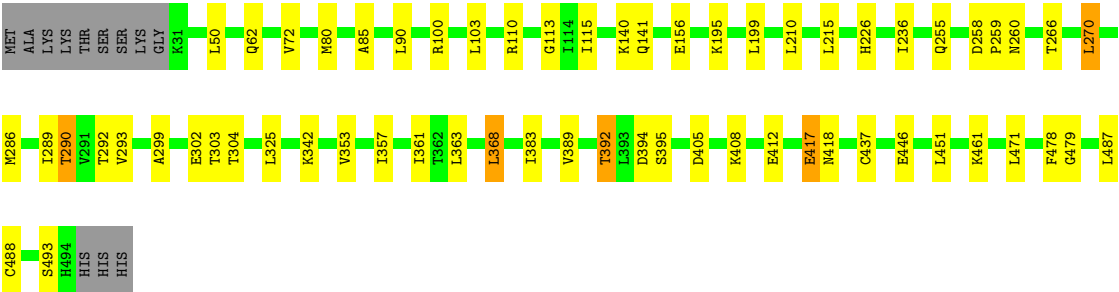


• Molecule 1: Cytochrome P450 2E1

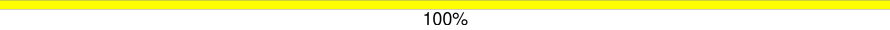


• Molecule 1: Cytochrome P450 2E1

Chain D:  84% 12% ..

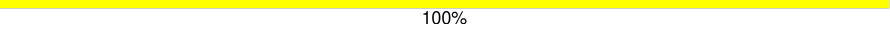


- Molecule 2: beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose

Chain E:  100%

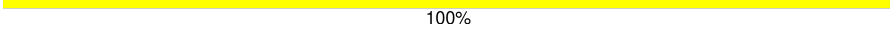


- Molecule 2: beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose

Chain F:  100%



- Molecule 2: beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose

Chain G:  100%



- Molecule 2: beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose

Chain H:  100%



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	100.64Å 100.64Å 259.45Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.98 – 2.35 36.98 – 2.35	Depositor EDS
% Data completeness (in resolution range)	95.9 (36.98-2.35) 95.9 (36.98-2.35)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.21 (at 2.34Å)	Xtriage
Refinement program	REFMAC 6.1.13	Depositor
R, R_{free}	0.217 , 0.281 0.222 , 0.279	Depositor DCC
R_{free} test set	5096 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	33.4	Xtriage
Anisotropy	0.057	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 22.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.448 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	16204	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 72.68 % 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 2.1083e-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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GLC, 9PL, FRU, HEM

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	0.62	0/3911	0.88	4/5296 (0.1%)
1	B	0.62	1/3902 (0.0%)	0.89	2/5284 (0.0%)
1	C	0.61	0/3893	0.90	4/5272 (0.1%)
1	D	0.59	0/3893	0.87	6/5272 (0.1%)
All	All	0.61	1/15599 (0.0%)	0.89	16/21124 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	53	ILE	CA-CB	6.81	1.58	1.54

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	302	GLU	N-CA-C	5.58	117.17	111.14
1	A	103	LEU	CA-C-N	5.50	126.71	119.84
1	A	103	LEU	C-N-CA	5.50	126.71	119.84
1	C	383	ILE	CA-C-N	5.41	125.31	119.85
1	C	383	ILE	C-N-CA	5.41	125.31	119.85
1	D	103	LEU	CA-C-N	5.38	125.46	119.32
1	D	103	LEU	C-N-CA	5.38	125.46	119.32
1	C	481	ILE	N-CA-C	5.38	113.79	107.77
1	B	481	ILE	N-CA-C	5.33	113.14	107.76
1	A	482	PRO	O-C-N	5.23	123.61	121.15
1	D	383	ILE	CA-C-N	5.15	125.13	120.03
1	D	383	ILE	C-N-CA	5.15	125.13	120.03
1	C	302	GLU	N-CA-C	5.09	116.63	111.14
1	B	302	GLU	N-CA-C	5.06	116.61	111.14
1	D	115	ILE	N-CA-C	5.04	115.66	110.36
1	A	302	GLU	N-CA-C	5.01	116.82	111.36

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3799	0	3800	29	0
1	B	3793	0	3794	38	0
1	C	3787	0	3788	36	0
1	D	3787	0	3788	26	0
2	E	23	0	21	0	0
2	F	23	0	21	0	0
2	G	23	0	21	0	0
2	H	23	0	21	1	0
3	A	43	0	30	2	0
3	B	43	0	30	2	0
3	C	43	0	30	7	0
3	D	43	0	30	4	0
4	A	15	0	16	1	0
4	B	15	0	16	2	0
4	C	15	0	16	2	0
4	D	15	0	16	2	0
5	A	202	0	0	2	0
5	B	202	0	0	3	0
5	C	151	0	0	4	0
5	D	159	0	0	0	0
All	All	16204	0	15438	139	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 (139) 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:D:392:THR:HG22	1:D:395:SER:OG	1.86	0.76
1:B:287:ASP:HA	1:B:290:THR:HG22	1.72	0.72
1:A:392:THR:HG22	1:A:395:SER:OG	1.92	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:80:MET:HE1	1:C:389:VAL:HG13	1.72	0.70
1:A:80:MET:HE1	1:A:389:VAL:HG13	1.73	0.69
3:D:500:HEM:HMC2	3:D:500:HEM:HBC2	1.74	0.69
1:D:80:MET:HE1	1:D:389:VAL:HG13	1.75	0.67
1:A:286:MET:O	1:A:290:THR:HG23	1.96	0.65
1:A:87:LYS:HE3	1:A:91:LEU:HD12	1.79	0.65
3:D:500:HEM:HBB2	3:D:500:HEM:HMB1	1.79	0.64
1:C:80:MET:HE1	1:C:389:VAL:CG1	2.28	0.63
1:B:392:THR:HG22	1:B:395:SER:OG	1.97	0.63
1:B:288:GLY:HA3	5:B:609:HOH:O	1.99	0.63
1:A:83:TYR:CZ	1:A:87:LYS:HD3	2.36	0.61
1:A:184:LEU:HG	1:A:293:VAL:HG13	1.83	0.60
1:B:48:LEU:HD21	1:B:70:LEU:HD21	1.82	0.59
1:B:80:MET:O	1:B:392:THR:HB	2.02	0.59
1:B:285:THR:HG22	1:B:287:ASP:H	1.68	0.58
1:D:266:THR:HG22	1:D:293:VAL:HG21	1.86	0.58
1:D:80:MET:HE1	1:D:389:VAL:CG1	2.34	0.57
1:C:87:LYS:HE3	1:C:91:LEU:HD12	1.85	0.57
1:C:416:ASN:ND2	5:C:549:HOH:O	2.37	0.57
1:A:325:LEU:HD13	1:A:349:TYR:HD2	1.70	0.57
1:D:303:THR:HG23	4:D:501:9PL:H6B	1.87	0.56
1:B:300:GLY:HA2	3:B:500:HEM:HMC3	1.86	0.56
1:A:180:ILE:HG13	1:A:184:LEU:HD22	1.88	0.55
3:C:500:HEM:CMC	3:C:500:HEM:HBC2	2.36	0.55
1:B:87:LYS:HE3	1:B:91:LEU:HD12	1.89	0.55
1:B:80:MET:HE1	1:B:389:VAL:HG13	1.89	0.55
1:B:287:ASP:HA	1:B:290:THR:CG2	2.37	0.55
1:C:325:LEU:HD11	1:C:353:VAL:HG11	1.89	0.55
1:B:184:LEU:HG	1:B:293:VAL:HG13	1.90	0.55
1:A:48:LEU:HD21	1:A:70:LEU:HD21	1.90	0.54
1:D:113:GLY:HA2	1:D:292:THR:OG1	2.07	0.54
1:B:248[B]:GLU:OE1	1:B:248[B]:GLU:HA	2.08	0.53
1:C:300:GLY:HA2	3:C:500:HEM:HMC3	1.90	0.53
1:A:266:THR:HG22	1:A:293:VAL:HG21	1.90	0.53
1:B:299:ALA:HA	4:B:501:9PL:H6A	1.91	0.53
1:B:344:ARG:NH1	5:B:2:HOH:O	2.41	0.53
1:C:140:LYS:N	1:C:140:LYS:HD3	2.23	0.53
3:C:500:HEM:HBC2	3:C:500:HEM:HMC1	1.92	0.52
1:C:461:LYS:HB3	1:C:488:CYS:HB2	1.92	0.52
1:B:87:LYS:HE3	1:B:91:LEU:CD1	2.41	0.51
1:A:300:GLY:HA2	3:A:500:HEM:HMC2	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:80:MET:CE	1:A:389:VAL:HG13	2.41	0.51
1:A:80:MET:O	1:A:392:THR:HB	2.11	0.51
1:C:87:LYS:HE3	1:C:91:LEU:CD1	2.42	0.50
1:B:303:THR:HB	3:B:500:HEM:CAB	2.42	0.50
1:C:303:THR:HG23	4:C:501:9PL:H6B	1.94	0.50
1:C:289:ILE:O	1:C:293:VAL:HG23	2.11	0.49
1:A:80:MET:HG3	1:A:85:ALA:HB1	1.94	0.49
3:D:500:HEM:HBC2	3:D:500:HEM:CMC	2.41	0.49
1:B:165:PRO:HB3	1:B:463:LEU:HD11	1.93	0.49
1:B:246:VAL:O	1:B:250:VAL:HG23	2.13	0.49
1:B:287:ASP:CA	1:B:290:THR:HG22	2.41	0.49
1:A:325:LEU:HD13	1:A:349:TYR:CD2	2.48	0.49
1:B:285:THR:HG22	1:B:286:MET:N	2.28	0.49
1:C:140:LYS:HD3	1:C:140:LYS:H	1.78	0.49
1:A:184:LEU:O	1:A:266:THR:HG23	2.13	0.48
1:A:285:THR:HG22	1:A:287:ASP:H	1.77	0.48
1:B:325:LEU:HD13	1:B:349:TYR:CD2	2.49	0.48
1:B:73:GLY:HA2	1:B:221:PHE:CE1	2.49	0.48
1:D:100:ARG:NH2	1:D:368:LEU:HB3	2.28	0.48
1:C:410:LYS:HB3	1:C:412:GLU:HG2	1.96	0.47
1:D:80:MET:HG3	1:D:85:ALA:HB1	1.97	0.47
1:B:325:LEU:HD21	1:B:353:VAL:HG11	1.96	0.47
1:D:299:ALA:HA	4:D:501:9PL:H6A	1.96	0.47
1:A:325:LEU:HD21	1:A:353:VAL:HG11	1.96	0.47
1:C:288:GLY:HA3	5:C:626:HOH:O	2.15	0.47
1:D:80:MET:O	1:D:392:THR:HB	2.15	0.47
1:D:437:CYS:HB2	3:D:500:HEM:NA	2.30	0.47
1:A:113:GLY:HA2	1:A:292:THR:OG1	2.13	0.47
1:B:416:ASN:ND2	5:B:511:HOH:O	2.47	0.47
1:C:134:ARG:O	1:C:138:MET:HG3	2.16	0.46
1:D:72:VAL:O	1:D:72:VAL:HG13	2.14	0.46
1:C:180:ILE:HD13	3:C:500:HEM:HBC1	1.97	0.46
1:C:184:LEU:O	1:C:266:THR:HG23	2.15	0.46
1:D:417:GLU:HG2	1:D:418:ASN:N	2.29	0.46
1:A:303:THR:HG23	4:A:501:9PL:H6A	1.99	0.45
1:A:243:LYS:NZ	5:A:627:HOH:O	2.49	0.45
1:B:285:THR:HB	1:B:288:GLY:H	1.82	0.45
1:C:286:MET:O	1:C:290:THR:HG23	2.16	0.45
1:B:325:LEU:HD13	1:B:349:TYR:HD2	1.81	0.45
1:C:363:LEU:O	1:C:479:GLY:HA2	2.16	0.45
1:C:209:LEU:HB3	1:C:215:LEU:HD13	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:211:SER:HA	1:C:477:GLY:HA3	1.99	0.44
1:C:304:THR:HG21	1:C:446:GLU:OE1	2.17	0.44
1:D:289:ILE:O	1:D:293:VAL:HG23	2.17	0.44
1:B:80:MET:HG3	1:B:85:ALA:HB1	2.00	0.44
1:B:266:THR:HG22	1:B:293:VAL:HG21	2.00	0.44
1:C:48:LEU:HD21	1:C:70:LEU:HD21	1.99	0.44
1:D:392:THR:HG23	1:D:394:ASP:OD1	2.18	0.44
1:D:226:HIS:HE1	1:D:236:ILE:HD13	1.83	0.44
1:B:209:LEU:HB3	1:B:215:LEU:HD13	1.99	0.43
1:D:405:ASP:HB3	1:D:408:LYS:HD2	2.00	0.43
1:B:145:SER:OG	1:B:149:ARG:NH2	2.51	0.43
2:H:1:GLC:O5	2:H:2:FRU:O6	2.36	0.43
1:B:80:MET:CE	1:B:389:VAL:HG13	2.48	0.43
1:C:392:THR:CG2	1:C:395:SER:OG	2.67	0.43
3:C:500:HEM:HMB2	3:C:500:HEM:HBB2	2.00	0.43
1:C:80:MET:HE2	1:C:80:MET:HB2	1.75	0.43
3:C:500:HEM:HBB2	3:C:500:HEM:CMB	2.48	0.43
1:A:363:LEU:O	1:A:479:GLY:HA2	2.19	0.43
1:D:304:THR:HG21	1:D:446:GLU:OE1	2.18	0.43
3:A:500:HEM:HBB2	3:A:500:HEM:HMB2	2.00	0.42
1:A:392:THR:HG22	1:A:395:SER:HG	1.85	0.42
1:B:103:LEU:HD21	1:B:369:PRO:HG3	2.02	0.42
1:D:325:LEU:HD11	1:D:353:VAL:HG11	2.00	0.42
1:A:59:ARG:NH2	5:A:583:HOH:O	2.49	0.42
1:B:180:ILE:HG13	1:B:184:LEU:HD22	2.00	0.42
1:C:249:ARG:O	1:C:253:HIS:ND1	2.48	0.42
1:D:270:LEU:HD12	1:D:289:ILE:HD13	2.01	0.42
1:A:212:THR:HB	1:A:213:PRO:HD2	2.01	0.42
1:D:363:LEU:O	1:D:479:GLY:HA2	2.19	0.42
1:B:198:ARG:O	1:B:202:LEU:HG	2.20	0.42
1:C:100:ARG:NH2	1:C:368:LEU:HB3	2.34	0.42
1:C:435:ARG:O	3:C:500:HEM:HBA2	2.20	0.42
1:C:184:LEU:HG	1:C:293:VAL:HG13	2.02	0.42
1:A:304:THR:HG21	1:A:446:GLU:OE1	2.20	0.41
1:C:490:ILE:HA	1:C:491:PRO:HD2	1.93	0.41
1:D:286:MET:O	1:D:290:THR:HG23	2.20	0.41
1:B:212:THR:HB	1:B:213:PRO:HD2	2.02	0.41
1:B:317:LYS:HD3	1:B:471:LEU:HD22	2.02	0.41
1:D:461:LYS:HB3	1:D:488:CYS:HB2	2.02	0.41
1:C:299:ALA:HA	4:C:501:9PL:H6A	2.03	0.41
1:B:184:LEU:O	1:B:266:THR:HG23	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:63:ARG:NH1	5:C:571:HOH:O	2.54	0.41
1:C:219:ASN:HB2	5:C:604:HOH:O	2.21	0.41
1:C:285:THR:HG22	1:C:286:MET:N	2.36	0.41
1:B:363:LEU:O	1:B:479:GLY:HA2	2.21	0.41
1:C:167:ASP:CG	1:C:169:THR:HG22	2.46	0.41
1:D:357:ILE:O	1:D:361:ILE:HG12	2.21	0.41
1:A:31:LYS:HB2	1:A:31:LYS:HE3	1.92	0.40
1:A:325:LEU:HD11	1:A:350:MET:HA	2.03	0.40
1:D:258:ASP:C	1:D:260:ASN:H	2.29	0.40
4:B:501:9PL:O15	4:B:501:9PL:H14A	2.20	0.40
1:A:225:LEU:HA	1:A:228:LEU:HD22	2.02	0.40
1:D:210:LEU:HD22	1:D:478:PHE:CE2	2.57	0.40
1:C:361:ILE:O	1:C:362:THR:C	2.64	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	464/476 (98%)	447 (96%)	17 (4%)	0	100	100
1	B	463/476 (97%)	443 (96%)	20 (4%)	0	100	100
1	C	462/476 (97%)	445 (96%)	16 (4%)	1 (0%)	43	52
1	D	462/476 (97%)	447 (97%)	14 (3%)	1 (0%)	43	52
All	All	1851/1904 (97%)	1782 (96%)	67 (4%)	2 (0%)	48	59

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	259	PRO
1	D	259	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	420/428 (98%)	397 (94%)	23 (6%)	19	24
1	B	419/428 (98%)	395 (94%)	24 (6%)	18	23
1	C	418/428 (98%)	392 (94%)	26 (6%)	16	19
1	D	418/428 (98%)	396 (95%)	22 (5%)	20	25
All	All	1675/1712 (98%)	1580 (94%)	95 (6%)	18	23

All (95) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	LEU
1	A	62	GLN
1	A	90	LEU
1	A	110	ARG
1	A	134	ARG
1	A	140	LYS
1	A	141	GLN
1	A	184	LEU
1	A	195	LYS
1	A	199	LEU
1	A	215	LEU
1	A	228	LEU
1	A	249	ARG
1	A	270	LEU
1	A	281	GLU
1	A	290	THR
1	A	325	LEU
1	A	342	LYS
1	A	368	LEU
1	A	392	THR
1	A	412	GLU
1	A	471	LEU
1	A	487	LEU
1	B	50	LEU

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Mol	Chain	Res	Type
1	B	67	VAL
1	B	72	VAL
1	B	90	LEU
1	B	110	ARG
1	B	184	LEU
1	B	195	LYS
1	B	199	LEU
1	B	215	LEU
1	B	225	LEU
1	B	228	LEU
1	B	270	LEU
1	B	303	THR
1	B	312	LEU
1	B	325	LEU
1	B	342	LYS
1	B	359	ARG
1	B	368	LEU
1	B	392	THR
1	B	412	GLU
1	B	417	GLU
1	B	451	LEU
1	B	471	LEU
1	B	487	LEU
1	C	90	LEU
1	C	110	ARG
1	C	140	LYS
1	C	141	GLN
1	C	145	SER
1	C	149	ARG
1	C	169	THR
1	C	184	LEU
1	C	192	ASN
1	C	195	LYS
1	C	199	LEU
1	C	215	LEU
1	C	249	ARG
1	C	303	THR
1	C	310	TYR
1	C	312	LEU
1	C	342	LYS
1	C	359	ARG
1	C	368	LEU

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Mol	Chain	Res	Type
1	C	392	THR
1	C	412	GLU
1	C	417	GLU
1	C	447	LEU
1	C	465	ASP
1	C	471	LEU
1	C	487	LEU
1	D	50	LEU
1	D	62	GLN
1	D	90	LEU
1	D	110	ARG
1	D	140	LYS
1	D	141	GLN
1	D	156	GLU
1	D	195	LYS
1	D	199	LEU
1	D	215	LEU
1	D	255	GLN
1	D	270	LEU
1	D	290	THR
1	D	342	LYS
1	D	368	LEU
1	D	392	THR
1	D	412	GLU
1	D	417	GLU
1	D	451	LEU
1	D	471	LEU
1	D	487	LEU
1	D	493	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	GLN
1	A	62	GLN
1	A	75	GLN
1	A	141	GLN
1	A	192	ASN
1	A	278	HIS
1	A	358	GLN
1	A	401	GLN
1	B	75	GLN

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Mol	Chain	Res	Type
1	B	141	GLN
1	B	188	HIS
1	B	192	ASN
1	B	326	HIS
1	B	401	GLN
1	C	75	GLN
1	D	62	GLN
1	D	107	HIS
1	D	192	ASN
1	D	358	GLN

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 ⓘ

8 monosaccharides 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$
2	GLC	E	1	2	11,11,12	0.89	0	15,15,17	1.86	5 (33%)
2	FRU	E	2	2	11,12,12	1.33	2 (18%)	10,18,18	1.80	2 (20%)
2	GLC	F	1	2	11,11,12	1.57	4 (36%)	15,15,17	2.06	5 (33%)
2	FRU	F	2	2	11,12,12	1.33	1 (9%)	10,18,18	1.40	1 (10%)
2	GLC	G	1	2	11,11,12	1.30	2 (18%)	15,15,17	1.67	2 (13%)
2	FRU	G	2	2	11,12,12	1.27	2 (18%)	10,18,18	1.60	1 (10%)
2	GLC	H	1	2	11,11,12	0.99	0	15,15,17	1.51	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	FRU	H	2	2	11,12,12	1.40	1 (9%)	10,18,18	1.27	1 (10%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GLC	E	1	2	-	2/2/19/22	0/1/1/1
2	FRU	E	2	2	-	0/5/24/24	0/1/1/1
2	GLC	F	1	2	-	2/2/19/22	0/1/1/1
2	FRU	F	2	2	-	3/5/24/24	0/1/1/1
2	GLC	G	1	2	-	0/2/19/22	0/1/1/1
2	FRU	G	2	2	-	2/5/24/24	0/1/1/1
2	GLC	H	1	2	-	0/2/19/22	0/1/1/1
2	FRU	H	2	2	-	1/5/24/24	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	2	FRU	O2-C2	3.44	1.46	1.40
2	F	2	FRU	O2-C2	3.10	1.46	1.40
2	G	2	FRU	O2-C2	2.97	1.45	1.40
2	E	2	FRU	O2-C2	2.97	1.45	1.40
2	F	1	GLC	O6-C6	2.54	1.53	1.42
2	E	2	FRU	O5-C2	2.46	1.47	1.43
2	F	1	GLC	O5-C5	2.40	1.48	1.43
2	F	1	GLC	O5-C1	2.26	1.47	1.43
2	G	1	GLC	O5-C5	2.26	1.47	1.43
2	G	1	GLC	O5-C1	2.11	1.47	1.43
2	F	1	GLC	C2-C3	2.03	1.55	1.52
2	G	2	FRU	O5-C5	2.01	1.48	1.43

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	2	FRU	O1-C1-C2	-4.24	102.29	111.67
2	H	1	GLC	O2-C2-C3	-3.95	101.97	110.15
2	E	2	FRU	O1-C1-C2	-3.91	103.02	111.67
2	G	1	GLC	C1-C2-C3	3.87	115.28	109.64
2	F	2	FRU	O1-C1-C2	-3.80	103.27	111.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	1	GLC	O2-C2-C3	-3.66	102.58	110.15
2	F	1	GLC	C1-C2-C3	3.46	114.68	109.64
2	E	1	GLC	O3-C3-C4	-3.22	102.78	110.38
2	F	1	GLC	O3-C3-C4	-3.18	102.88	110.38
2	E	2	FRU	O2-C2-O5	-3.06	103.45	109.33
2	F	1	GLC	O5-C5-C4	-2.92	103.73	110.83
2	E	1	GLC	O2-C2-C3	-2.89	104.16	110.15
2	E	1	GLC	C1-O5-C5	2.81	115.95	112.19
2	H	2	FRU	O3-C3-C4	-2.73	103.60	113.25
2	G	1	GLC	O2-C2-C3	-2.54	104.90	110.15
2	E	1	GLC	O3-C3-C2	-2.44	105.07	110.05
2	E	1	GLC	O5-C1-C2	-2.40	105.08	110.79
2	H	1	GLC	O3-C3-C2	-2.27	105.43	110.05
2	F	1	GLC	O4-C4-C3	-2.15	105.31	110.38

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	F	2	FRU	O1-C1-C2-C3
2	F	2	FRU	O1-C1-C2-O2
2	G	2	FRU	O1-C1-C2-O2
2	F	2	FRU	O1-C1-C2-O5
2	F	1	GLC	O5-C5-C6-O6
2	G	2	FRU	O1-C1-C2-O5
2	F	1	GLC	C4-C5-C6-O6
2	H	2	FRU	O5-C5-C6-O6
2	E	1	GLC	O5-C5-C6-O6
2	E	1	GLC	C4-C5-C6-O6

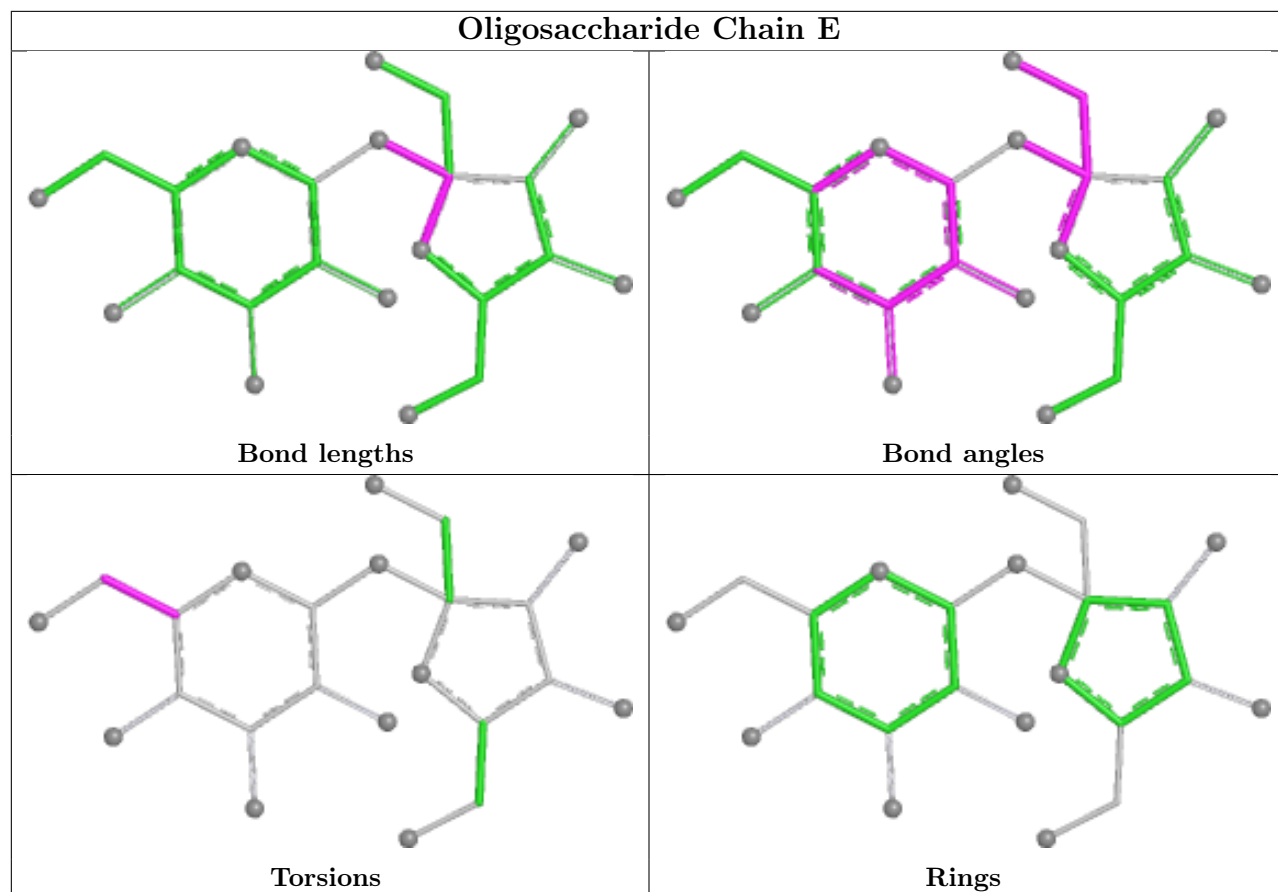
There are no ring outliers.

2 monomers are involved in 1 short contact:

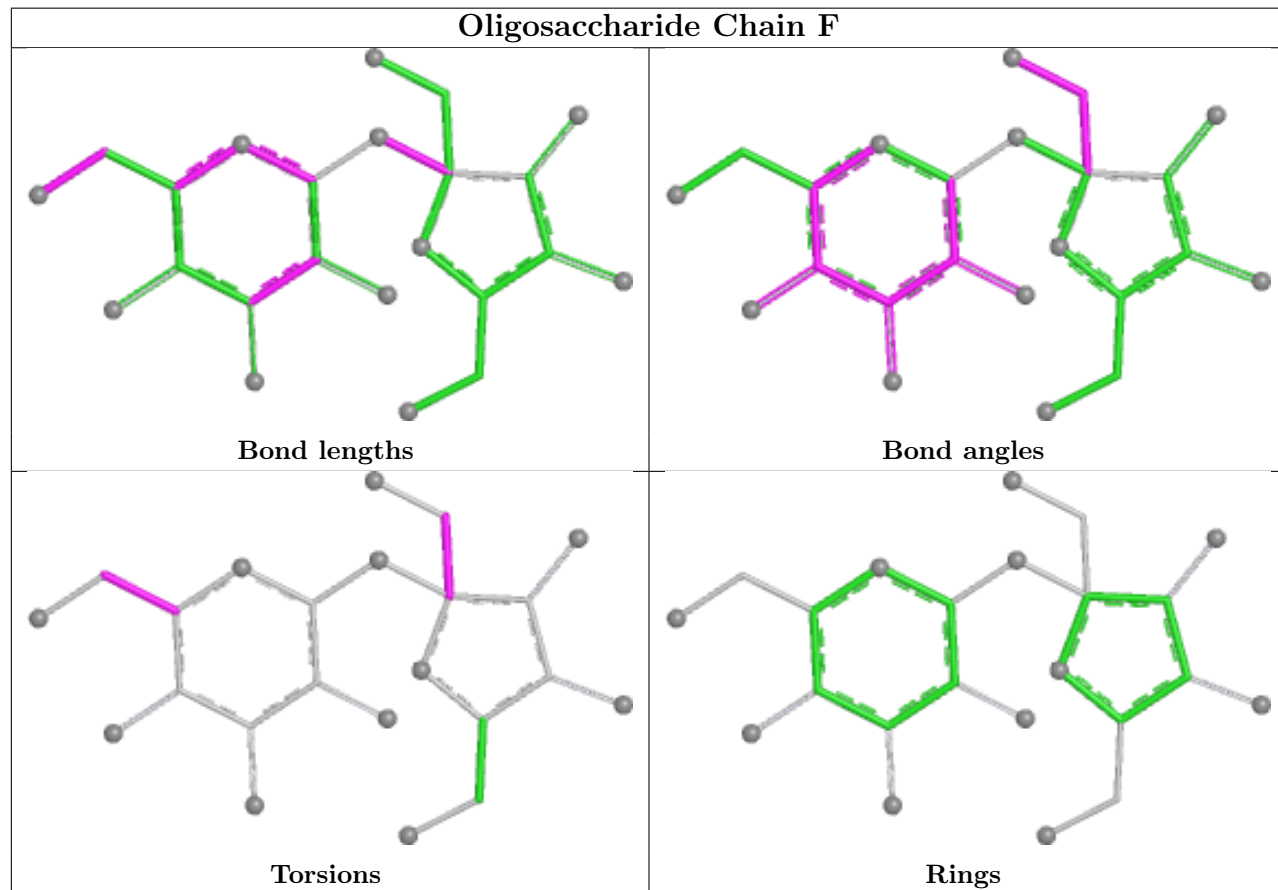
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	H	1	GLC	1	0
2	H	2	FRU	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

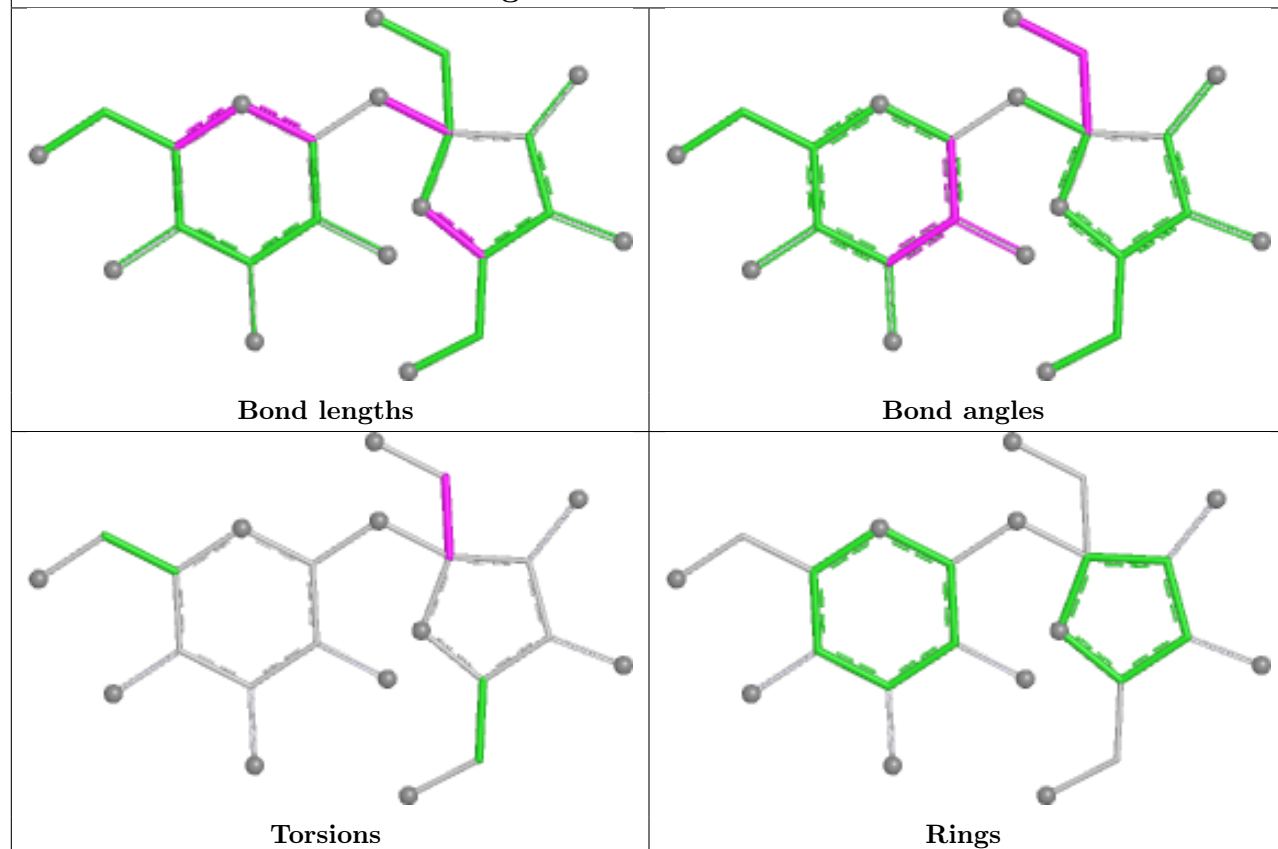
Oligosaccharide Chain E



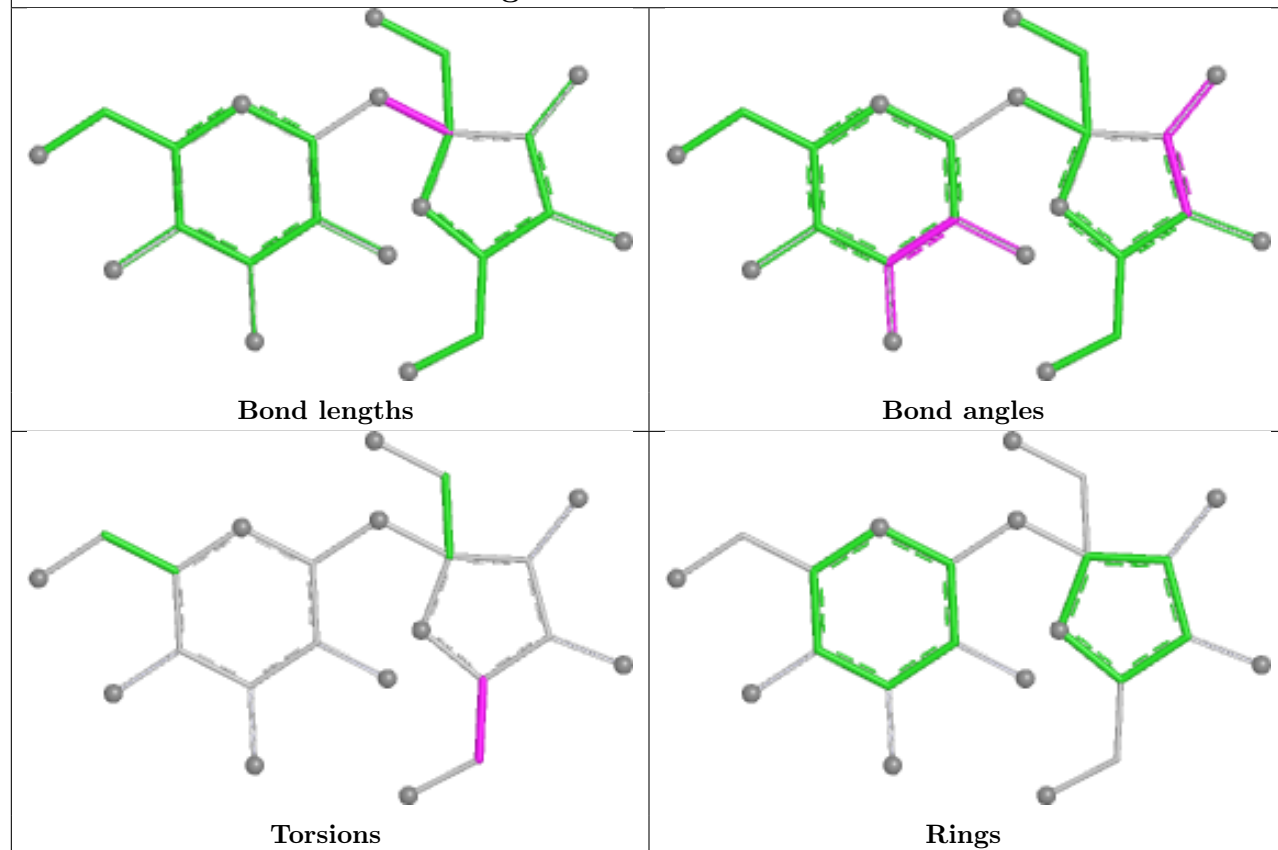
Oligosaccharide Chain F



Oligosaccharide Chain G



Oligosaccharide Chain H



5.6 Ligand geometry

8 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
4	9PL	B	501	3	16,16,16	1.72	3 (18%)	14,22,22	3.38	4 (28%)
3	HEM	C	500	1,4	50,50,50	1.95	12 (24%)	67,82,82	1.31	8 (11%)
4	9PL	D	501	3	16,16,16	1.69	3 (18%)	14,22,22	2.94	6 (42%)
4	9PL	A	501	3	16,16,16	1.55	3 (18%)	14,22,22	2.98	5 (35%)
3	HEM	A	500	1,4	50,50,50	2.02	11 (22%)	67,82,82	1.20	6 (8%)
3	HEM	B	500	1,4	50,50,50	1.79	9 (18%)	67,82,82	1.01	3 (4%)
4	9PL	C	501	3	16,16,16	1.77	4 (25%)	14,22,22	2.56	4 (28%)
3	HEM	D	500	1,4	50,50,50	2.05	12 (24%)	67,82,82	1.29	6 (8%)

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
4	9PL	B	501	3	-	0/6/19/19	0/2/2/2
3	HEM	C	500	1,4	-	2/14/54/54	-
4	9PL	D	501	3	-	3/6/19/19	0/2/2/2
4	9PL	A	501	3	-	0/6/19/19	0/2/2/2
3	HEM	A	500	1,4	-	0/14/54/54	-
3	HEM	B	500	1,4	-	2/14/54/54	-
4	9PL	C	501	3	-	3/6/19/19	0/2/2/2
3	HEM	D	500	1,4	-	0/14/54/54	-

All (57) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	500	HEM	C3D-C2D	7.74	1.53	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	500	HEM	C3D-C2D	6.96	1.51	1.36
3	B	500	HEM	C3D-C2D	6.59	1.51	1.36
3	C	500	HEM	C3D-C2D	6.25	1.50	1.36
3	A	500	HEM	FE-ND	5.18	2.10	1.94
3	D	500	HEM	FE-ND	5.01	2.10	1.94
3	D	500	HEM	FE-NA	4.61	2.10	1.95
4	B	501	9PL	O10-C11	4.53	1.44	1.35
4	C	501	9PL	O10-C11	4.38	1.44	1.35
3	C	500	HEM	FE-NC	4.15	2.08	1.95
4	D	501	9PL	O10-C11	4.13	1.44	1.35
3	C	500	HEM	FE-NA	4.12	2.08	1.95
3	D	500	HEM	CMC-C2C	3.96	1.58	1.50
4	A	501	9PL	O10-C11	3.85	1.43	1.35
3	B	500	HEM	FE-ND	3.75	2.06	1.94
3	A	500	HEM	CAC-C3C	3.68	1.57	1.47
3	D	500	HEM	CAC-C3C	3.62	1.57	1.47
3	C	500	HEM	CAC-C3C	3.61	1.57	1.47
3	C	500	HEM	FE-ND	3.59	2.05	1.94
3	B	500	HEM	CAC-C3C	3.54	1.56	1.47
3	A	500	HEM	C2A-C3A	-3.50	1.29	1.38
3	A	500	HEM	FE-NC	3.39	2.06	1.95
4	D	501	9PL	C4-C5	3.35	1.41	1.36
3	C	500	HEM	CMC-C2C	3.31	1.57	1.50
4	C	501	9PL	C4-C5	3.18	1.41	1.36
3	D	500	HEM	CMD-C2D	3.14	1.57	1.50
3	D	500	HEM	C1A-NA	-3.09	1.33	1.39
3	B	500	HEM	CMC-C2C	3.09	1.57	1.50
3	D	500	HEM	FE-NC	3.07	2.05	1.95
3	C	500	HEM	CMD-C2D	3.04	1.57	1.50
3	C	500	HEM	CAB-C3B	2.99	1.55	1.47
3	A	500	HEM	CAB-C3B	2.90	1.55	1.47
3	D	500	HEM	CAB-C3B	2.86	1.55	1.47
3	A	500	HEM	CMC-C2C	2.85	1.56	1.50
3	C	500	HEM	C4C-NC	-2.82	1.34	1.39
3	B	500	HEM	CAB-C3B	2.81	1.54	1.47
4	A	501	9PL	C6-N1	2.74	1.52	1.46
3	B	500	HEM	FE-NC	2.68	2.04	1.95
3	A	500	HEM	FE-NA	2.63	2.03	1.95
4	C	501	9PL	C5-N1	-2.62	1.33	1.38
4	B	501	9PL	C4-N3	-2.59	1.33	1.37
3	D	500	HEM	CMB-C2B	2.53	1.55	1.50
4	B	501	9PL	C6-N1	2.52	1.51	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	500	HEM	CMB-C2B	2.46	1.55	1.50
4	C	501	9PL	C4-N3	-2.41	1.33	1.37
3	B	500	HEM	FE-NA	2.30	2.02	1.95
4	A	501	9PL	C4-N3	-2.26	1.33	1.37
4	D	501	9PL	C12-C11	-2.25	1.48	1.51
3	D	500	HEM	C4C-NC	-2.23	1.35	1.39
3	A	500	HEM	C4A-C3A	2.22	1.48	1.43
3	B	500	HEM	C1B-NB	-2.21	1.36	1.40
3	C	500	HEM	C1A-NA	-2.18	1.35	1.39
3	B	500	HEM	FE-NB	2.18	2.01	1.94
3	A	500	HEM	CMD-C2D	2.17	1.55	1.50
3	A	500	HEM	CMA-C3A	2.08	1.55	1.50
3	C	500	HEM	C3B-C2B	-2.01	1.33	1.37
3	D	500	HEM	C2A-C3A	-2.00	1.33	1.38

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	501	9PL	N1-C2-N3	-10.62	106.62	112.94
4	A	501	9PL	N1-C2-N3	-9.15	107.49	112.94
4	D	501	9PL	N1-C2-N3	-7.84	108.27	112.94
4	C	501	9PL	N1-C2-N3	-7.24	108.63	112.94
3	A	500	HEM	C4D-ND-C1D	4.05	110.01	105.21
3	D	500	HEM	C4D-ND-C1D	3.97	109.91	105.21
4	D	501	9PL	O10-C11-O15	3.95	125.63	121.43
4	B	501	9PL	O10-C11-O15	3.78	125.44	121.43
4	C	501	9PL	O10-C11-O15	3.58	125.23	121.43
4	A	501	9PL	C4-N3-C2	3.51	109.83	105.24
4	B	501	9PL	C4-N3-C2	3.45	109.75	105.24
3	C	500	HEM	CMD-C2D-C1D	3.37	130.30	125.03
4	D	501	9PL	C4-N3-C2	3.22	109.45	105.24
4	B	501	9PL	O15-C11-C12	-3.21	124.40	128.81
3	B	500	HEM	C4D-ND-C1D	3.19	108.99	105.21
3	A	500	HEM	C1B-NB-C4B	3.05	108.82	105.21
3	C	500	HEM	C4D-ND-C1D	3.02	108.79	105.21
3	C	500	HEM	CAA-CBA-CGA	-2.95	105.84	113.67
4	D	501	9PL	C14-C13-C12	-2.95	107.16	113.41
3	D	500	HEM	CAD-C3D-C4D	2.87	129.71	124.70
4	D	501	9PL	O10-C9-C8	-2.77	100.68	105.40
3	D	500	HEM	C3B-C2B-C1B	2.72	108.45	106.41
3	C	500	HEM	C1D-C2D-C3D	-2.66	104.18	106.98
4	A	501	9PL	O10-C11-O15	2.65	124.25	121.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	500	HEM	C3B-C2B-C1B	2.65	108.40	106.41
4	C	501	9PL	C4-N3-C2	2.64	108.70	105.24
4	A	501	9PL	O15-C11-C12	-2.63	125.20	128.81
3	C	500	HEM	C4C-C3C-C2C	2.58	109.05	106.81
3	B	500	HEM	O2A-CGA-CBA	2.53	122.01	114.00
3	C	500	HEM	CAD-C3D-C4D	2.51	129.07	124.70
3	A	500	HEM	C4D-C3D-C2D	-2.37	103.44	106.89
3	B	500	HEM	C4C-C3C-C2C	2.32	108.83	106.81
3	D	500	HEM	CAA-CBA-CGA	-2.29	107.58	113.67
3	A	500	HEM	CMB-C2B-C1B	-2.26	121.50	125.03
4	C	501	9PL	O10-C11-C12	-2.26	107.27	110.19
3	A	500	HEM	O2A-CGA-CBA	2.14	120.75	114.00
3	D	500	HEM	CMD-C2D-C1D	2.12	128.35	125.03
3	C	500	HEM	C4A-NA-C1A	2.11	109.27	105.82
4	D	501	9PL	C9-C8-C12	2.07	104.63	101.42
3	D	500	HEM	CAD-CBD-CGD	-2.06	108.19	113.67
4	A	501	9PL	C7-C8-C9	-2.04	108.57	112.39
3	C	500	HEM	CBD-CAD-C3D	-2.03	106.92	112.53

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	501	9PL	C5-C7-C8-C9
4	D	501	9PL	C5-C7-C8-C9
4	D	501	9PL	C8-C12-C13-C14
4	D	501	9PL	C11-C12-C13-C14
4	C	501	9PL	C11-C12-C13-C14
4	C	501	9PL	C8-C12-C13-C14
3	C	500	HEM	CAA-CBA-CGA-O1A
3	C	500	HEM	CAA-CBA-CGA-O2A
3	B	500	HEM	CAA-CBA-CGA-O1A
3	B	500	HEM	CAA-CBA-CGA-O2A

There are no ring outliers.

8 monomers are involved in 22 short contacts:

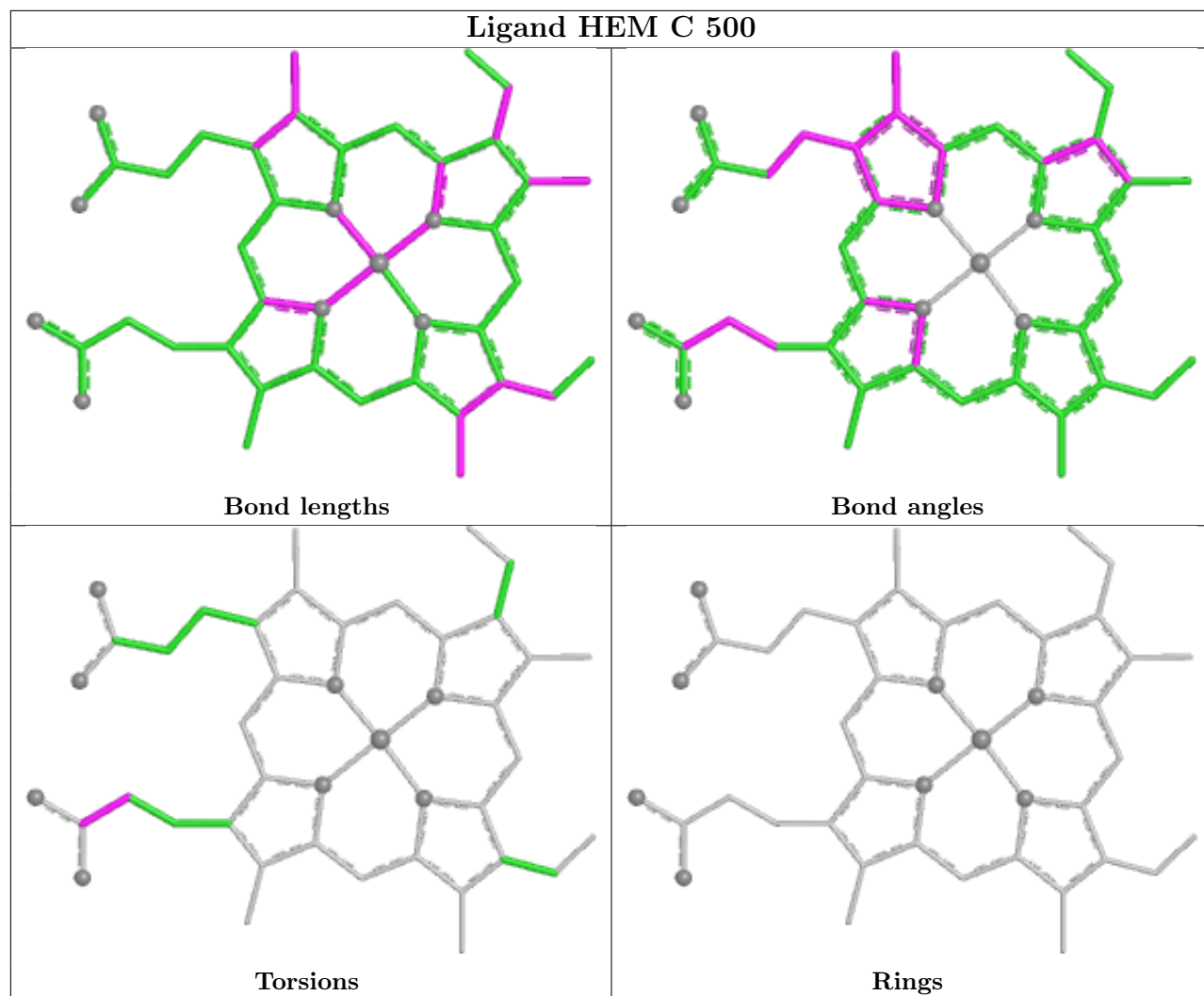
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	501	9PL	2	0
3	C	500	HEM	7	0
4	D	501	9PL	2	0

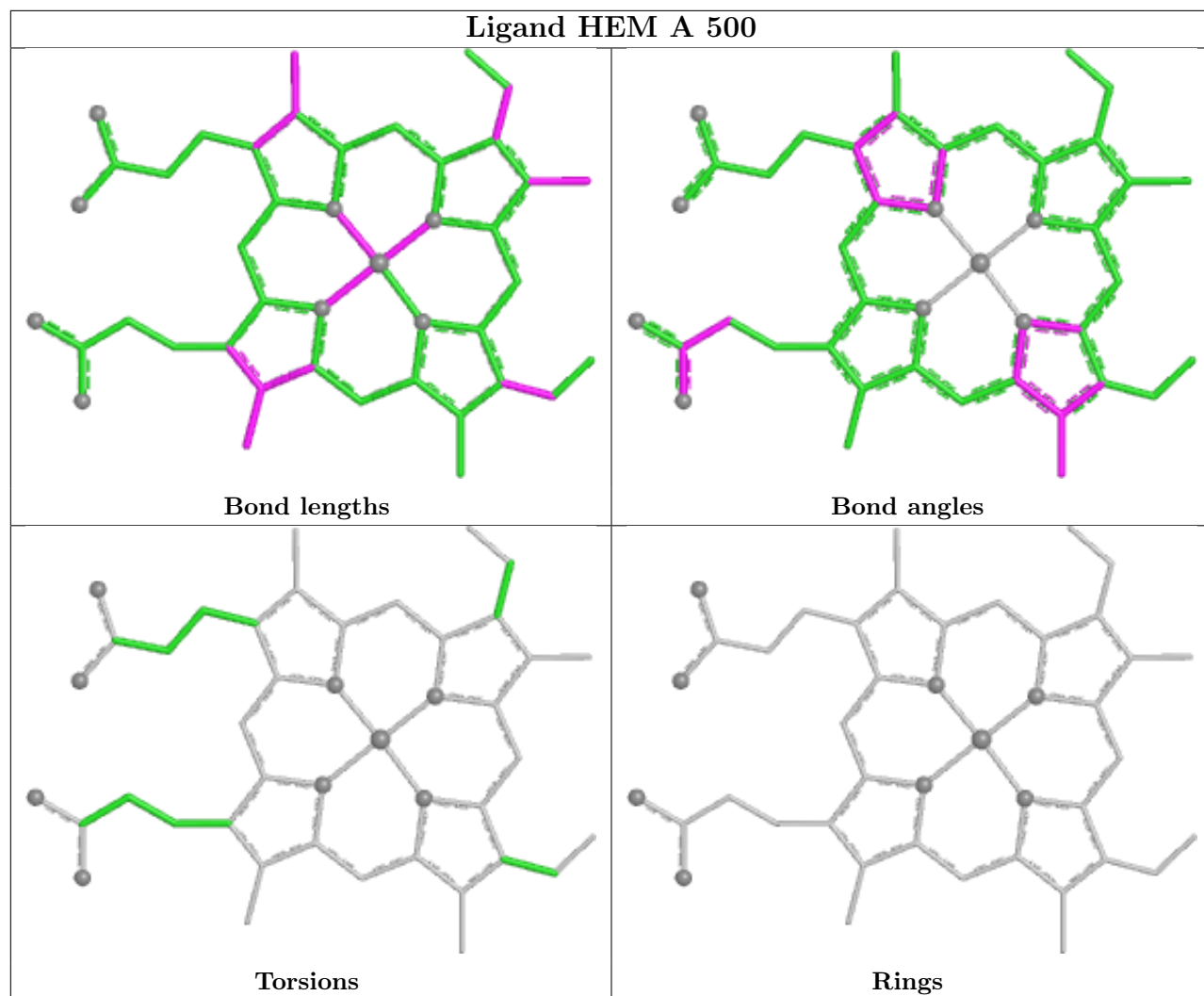
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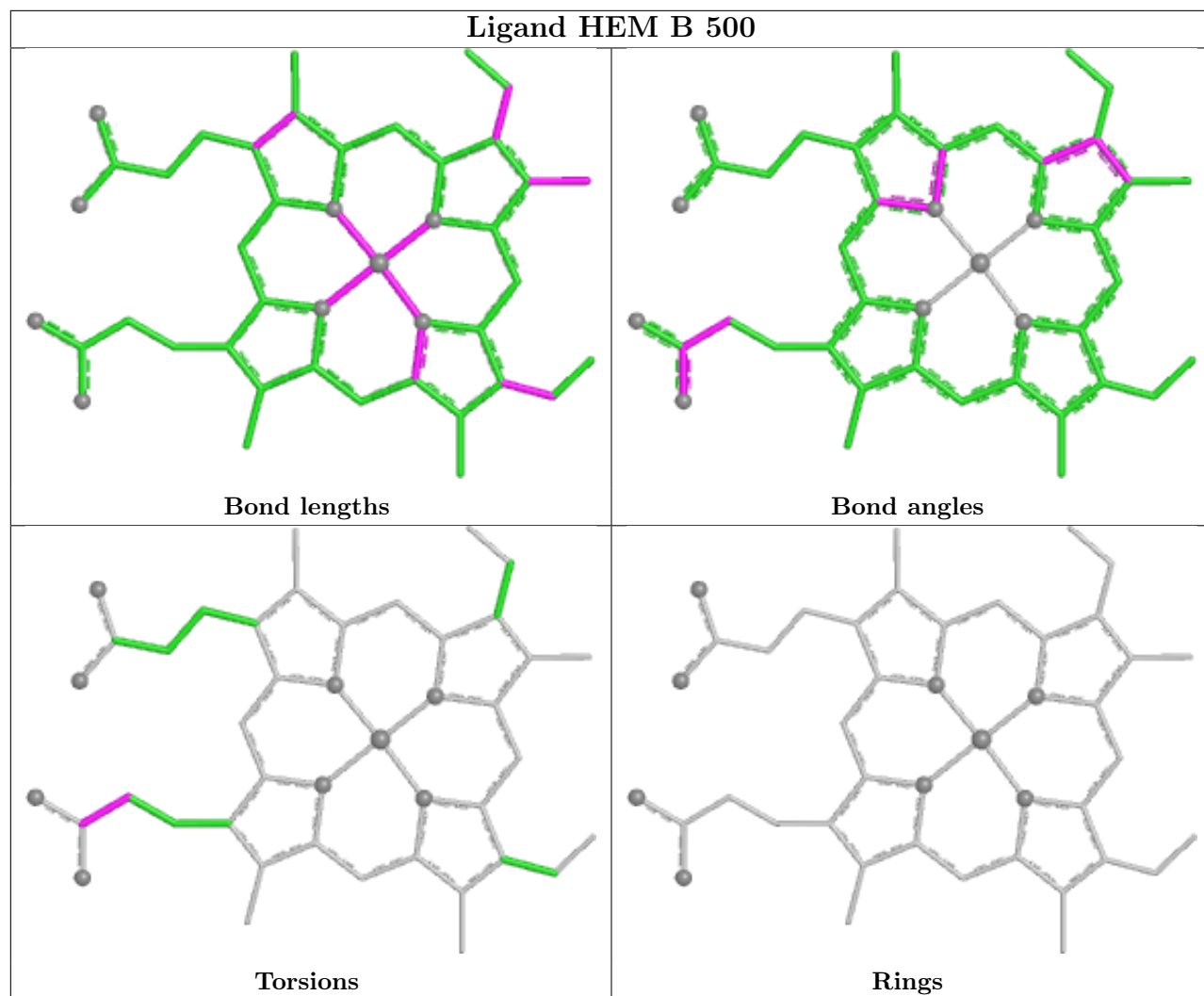
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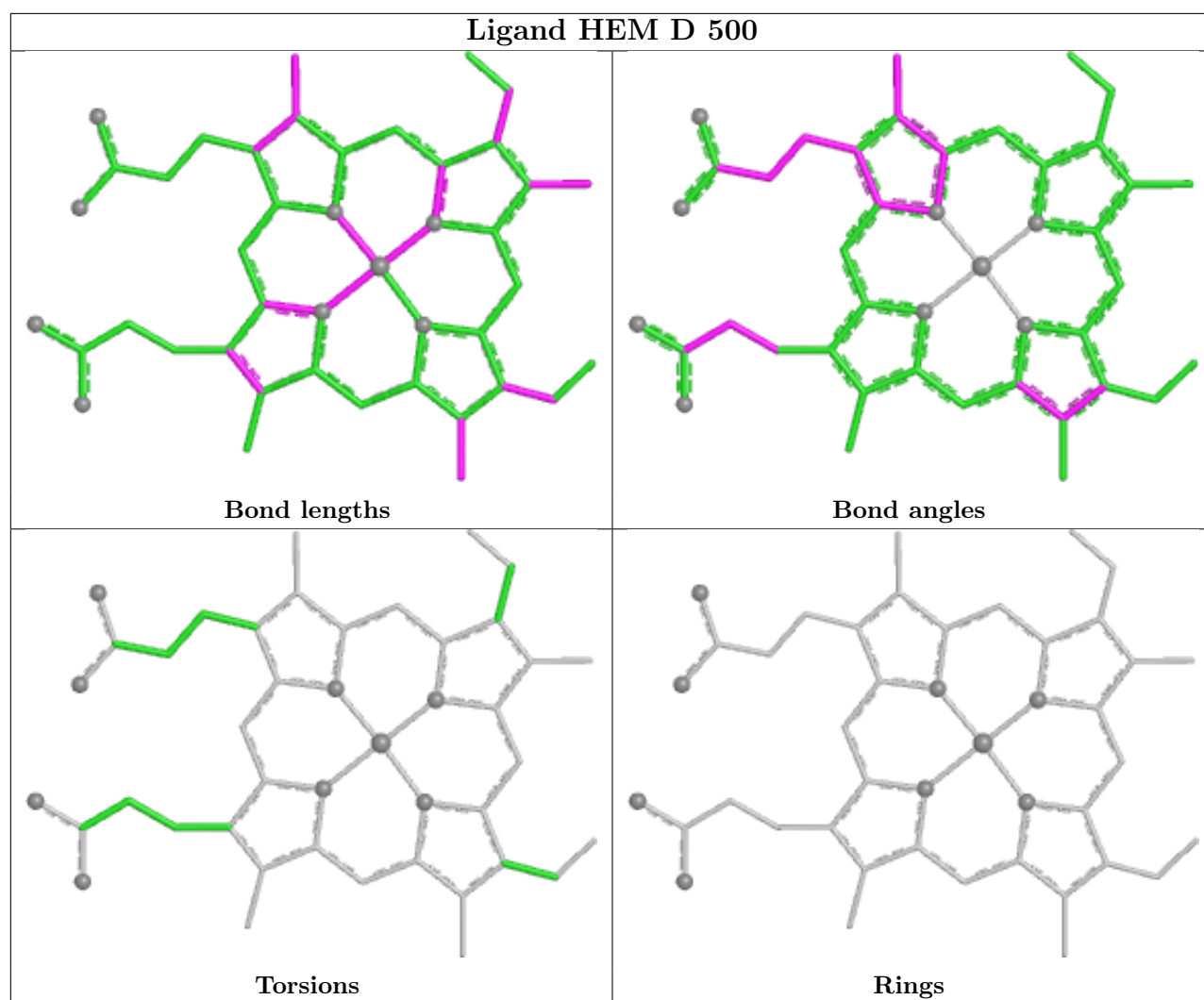
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	501	9PL	1	0
3	A	500	HEM	2	0
3	B	500	HEM	2	0
4	C	501	9PL	2	0
3	D	500	HEM	4	0

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	464/476 (97%)	-1.44	0 100 100	17, 28, 50, 71	2 (0%)
1	B	464/476 (97%)	-1.43	0 100 100	17, 28, 51, 69	1 (0%)
1	C	464/476 (97%)	-1.33	0 100 100	17, 33, 62, 75	0
1	D	464/476 (97%)	-1.33	0 100 100	17, 32, 61, 76	0
All	All	1856/1904 (97%)	-1.38	0 100 100	17, 30, 57, 76	3 (0%)

There are no RSRZ outliers to report.

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

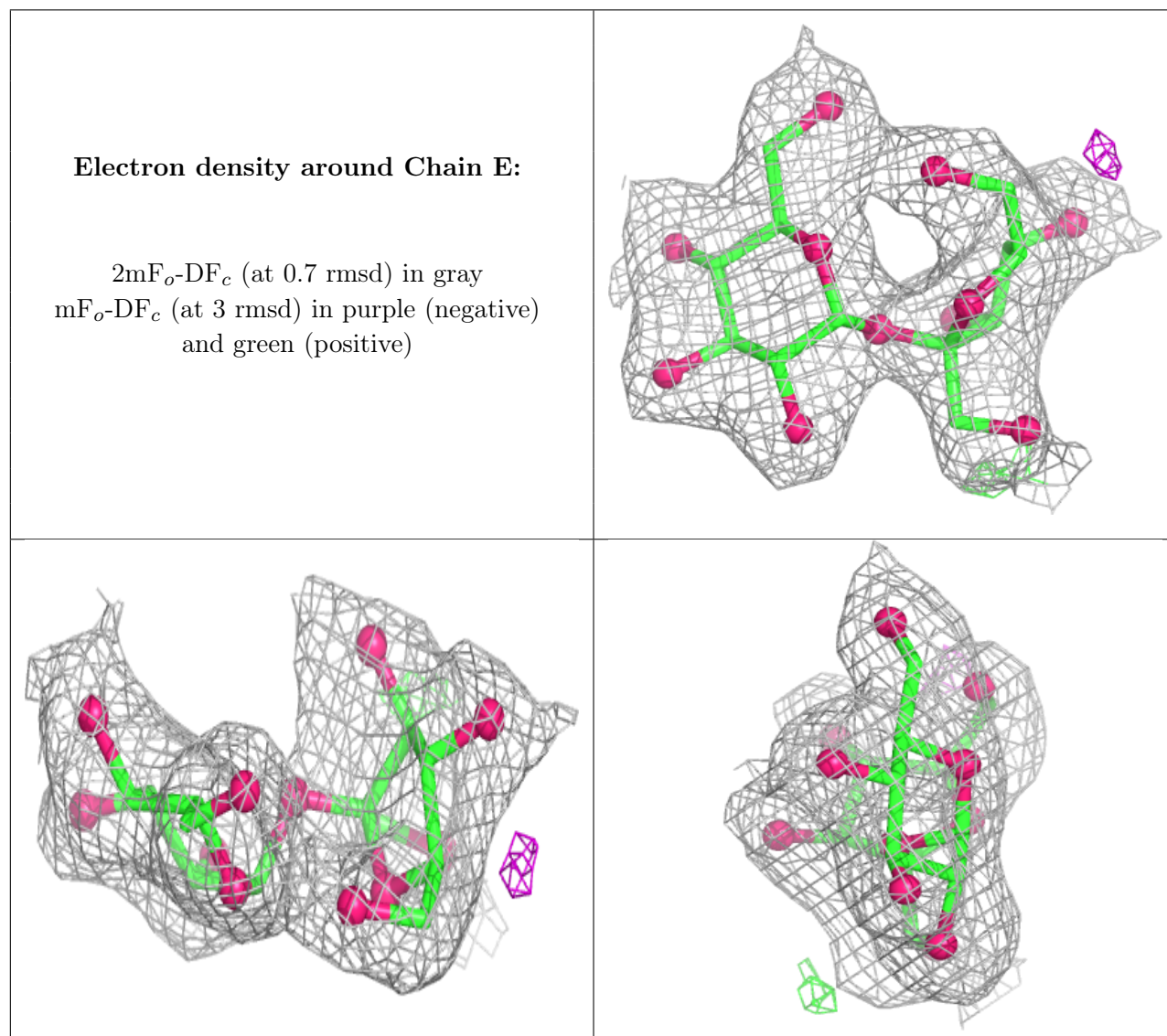
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

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.

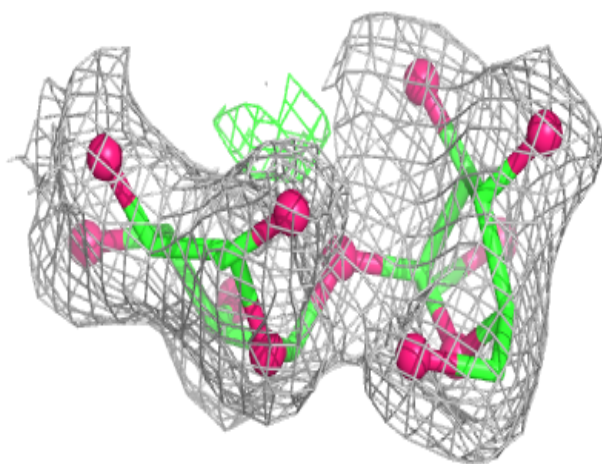
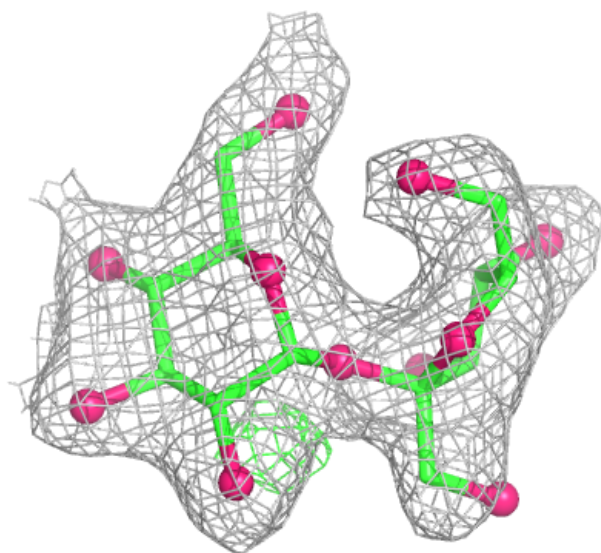
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GLC	F	1	11/12	0.98	0.03	25,27,32,38	0
2	FRU	F	2	12/12	0.98	0.03	26,29,31,34	0
2	GLC	E	1	11/12	0.99	0.03	24,26,28,29	0
2	FRU	E	2	12/12	0.99	0.03	24,29,30,31	0
2	GLC	G	1	11/12	0.99	0.03	26,29,32,32	0
2	FRU	G	2	12/12	0.99	0.03	26,29,32,33	0
2	GLC	H	1	11/12	0.99	0.03	25,28,29,30	0
2	FRU	H	2	12/12	0.99	0.03	27,28,32,35	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



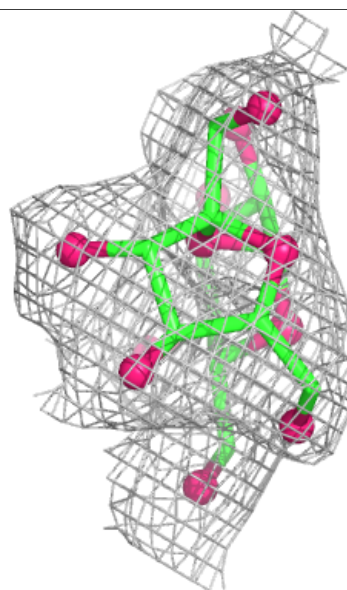
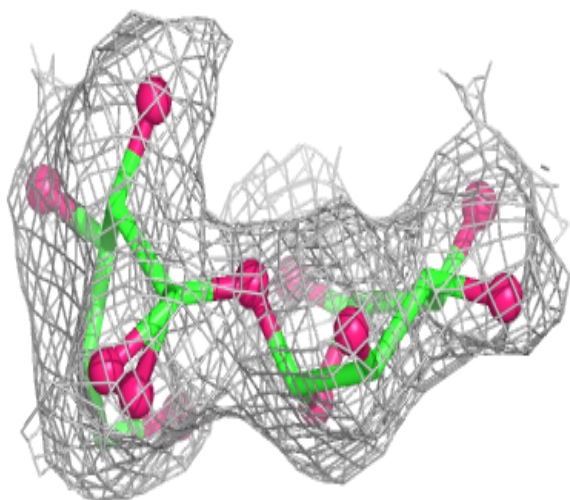
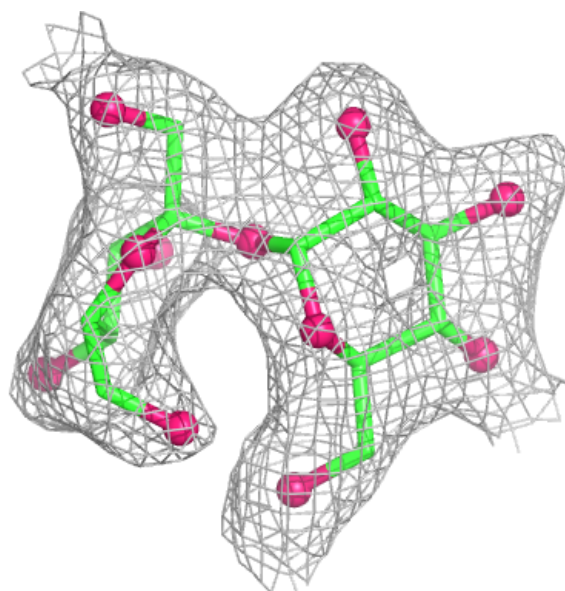
Electron density around Chain F:

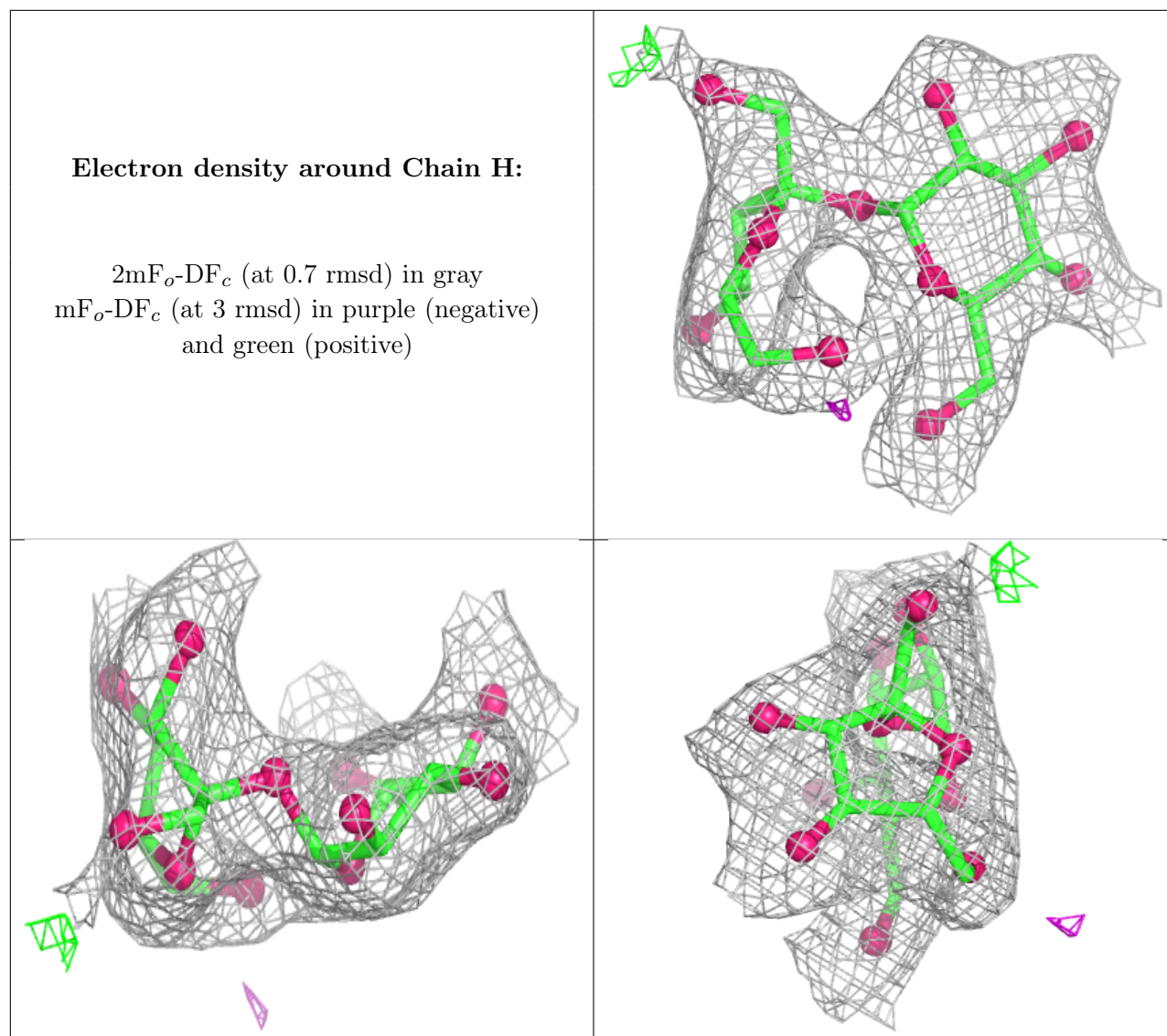
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain G:

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





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.

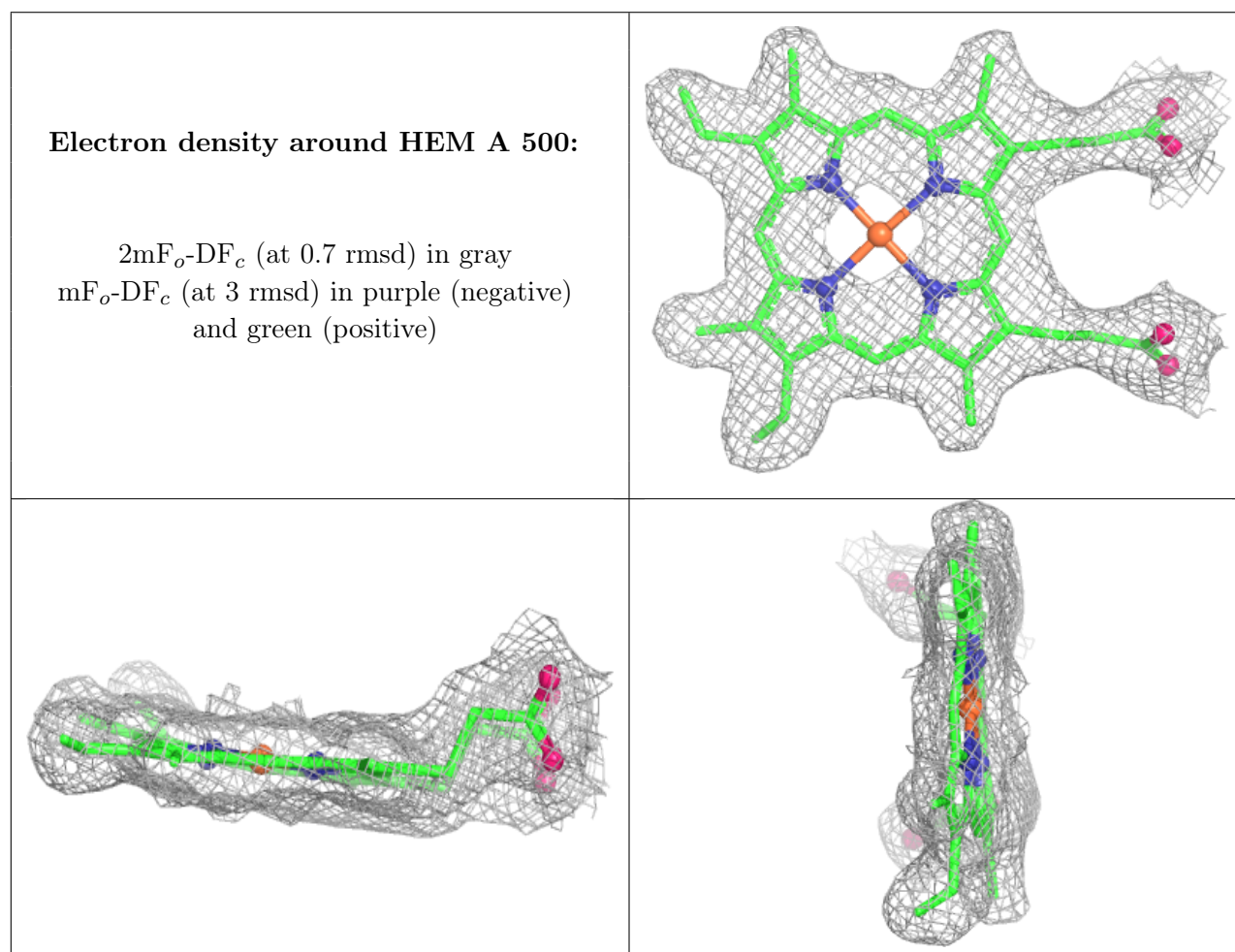
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	9PL	A	501	15/15	0.99	0.03	18,25,31,32	0
4	9PL	B	501	15/15	0.99	0.04	20,24,26,27	0
4	9PL	C	501	15/15	0.99	0.03	24,28,34,35	0
4	9PL	D	501	15/15	0.99	0.03	19,25,34,36	0
3	HEM	A	500	43/43	1.00	0.02	14,22,26,26	0
3	HEM	B	500	43/43	1.00	0.03	14,21,24,27	0
3	HEM	C	500	43/43	1.00	0.03	13,21,24,29	0

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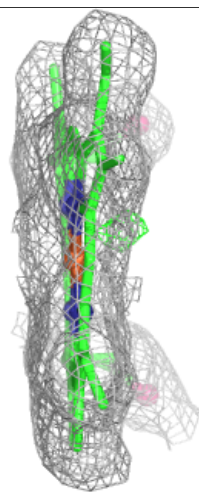
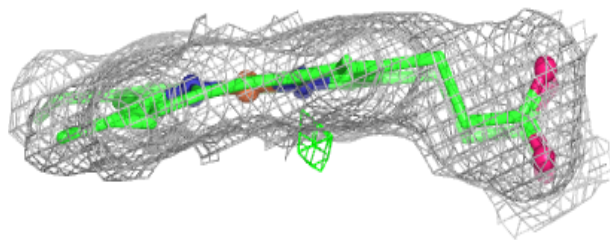
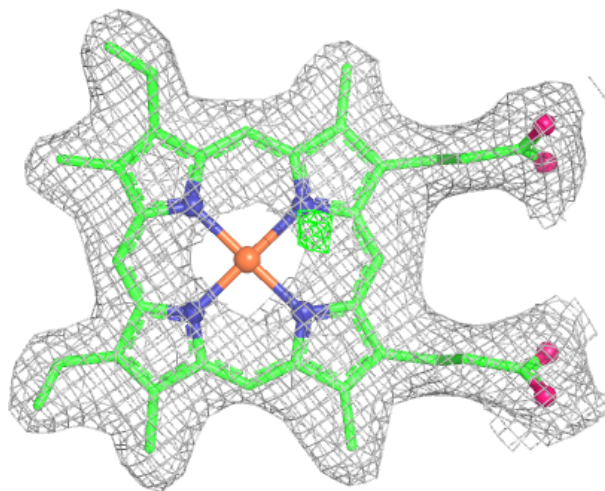
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	HEM	D	500	43/43	1.00	0.02	17,23,27,31	0

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.



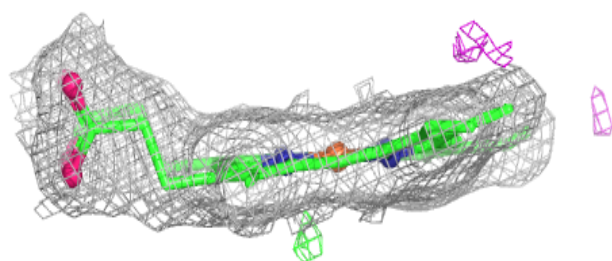
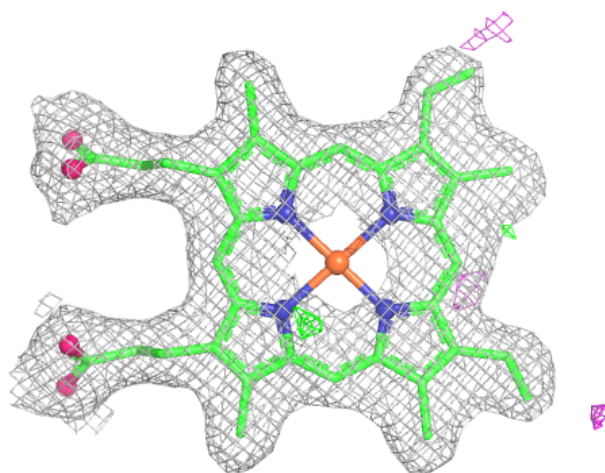
Electron density around HEM B 500:

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



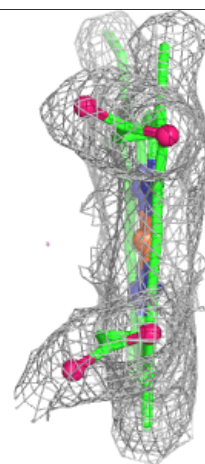
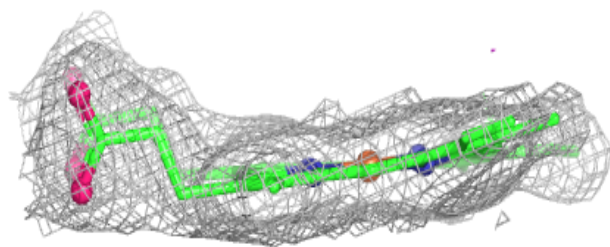
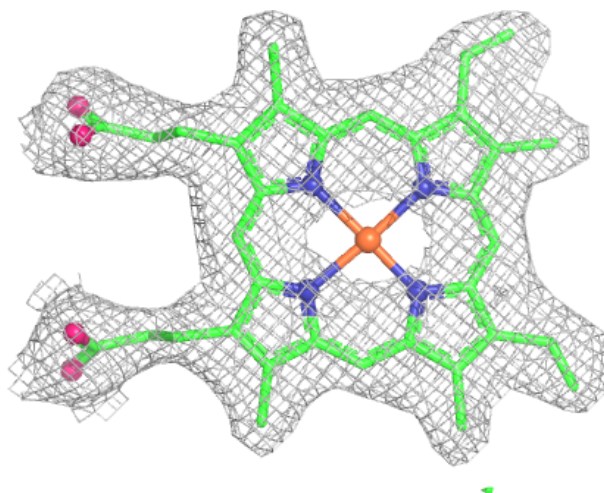
Electron density around HEM C 500:

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



Electron density around HEM D 500:

$2mF_o-DF_c$ (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.