Full wwPDB X-ray Structure Validation Report (i)

Nov 2, 2023 – 06:35 AM EDT

PDB ID	:	3SXV
Title	:	Crystal structure of the complex of goat lactoperoxidase with amitrole at 2.1 A resolution
Authors	:	Pandey, N.; Singh, R.P.; Singh, A.K.; Sinha, M.; Kaur, P.; Sharma, S.; Singh, T.P.
Deposited on	:	2011-07-15
Resolution	:	2.10 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 (i) 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 (i)) were used in the production of this report:

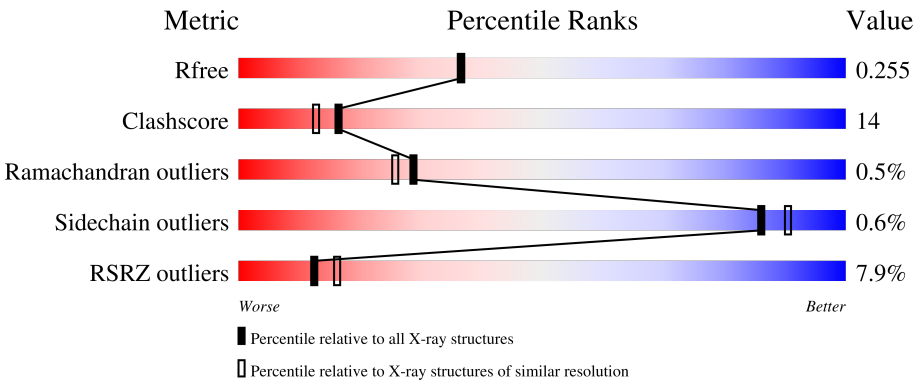
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 <=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	595	8% 81% 18% •
2	B	2	100%

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
12	MPD	A	621	-	-	X	-
9	3TR	A	597	-	-	X	-
9	3TR	A	751	-	-	X	-

2 Entry composition (i)

There are 14 unique types of molecules in this entry. The entry contains 5252 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.

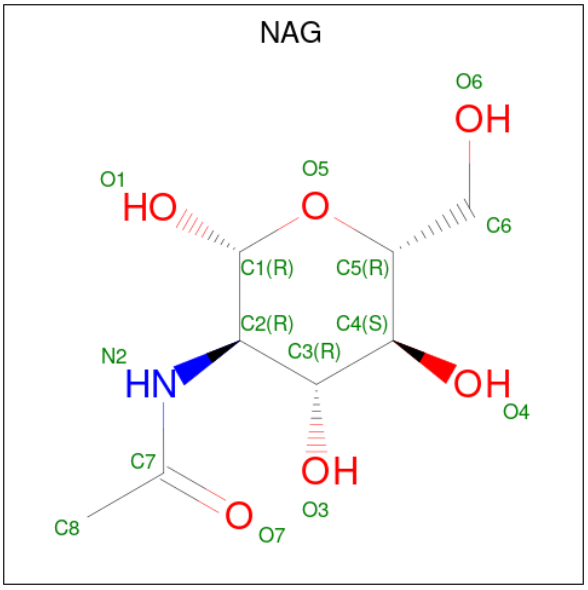
- Molecule 1 is a protein called Lactoperoxidase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	A	595	4758	3021	844	866	1	26	0	0	0

- Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

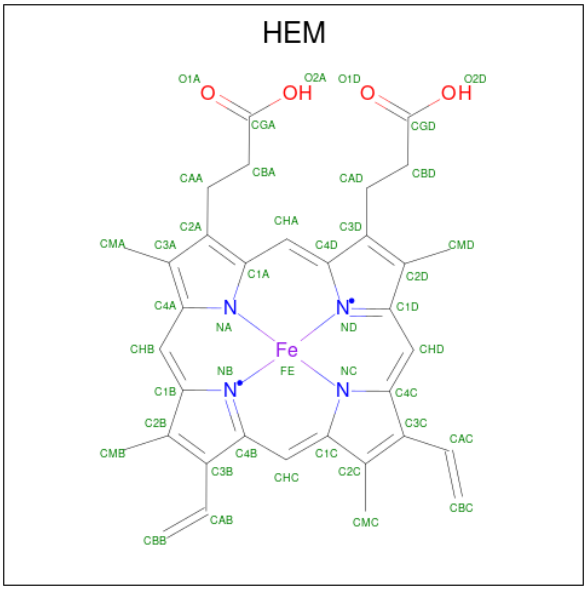
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	B	2	28	16	2	10	0	0	0

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	14	8	1	5	0	0
			Total	C	N	O		
3	A	1	14	8	1	5	0	0
			Total	C	N	O		
3	A	1	14	8	1	5	0	0

- Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Ca		
5	A	1	1	1	0	0

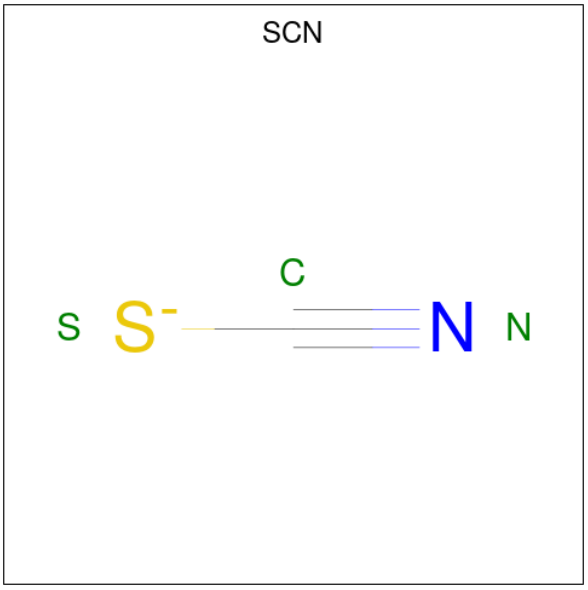
- Molecule 6 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
6	A	1	1	1	0	0

- Molecule 7 is IODIDE ION (three-letter code: IOD) (formula: I).

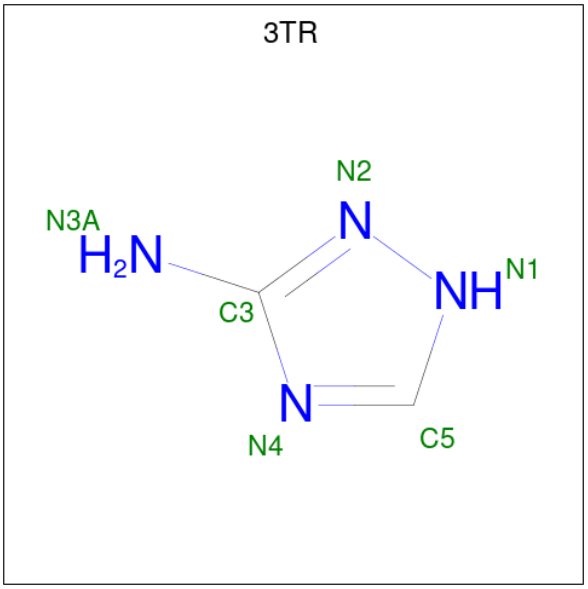
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	12	Total I 12 12	0	0

- Molecule 8 is THIOCYANATE ION (three-letter code: SCN) (formula: CNS).



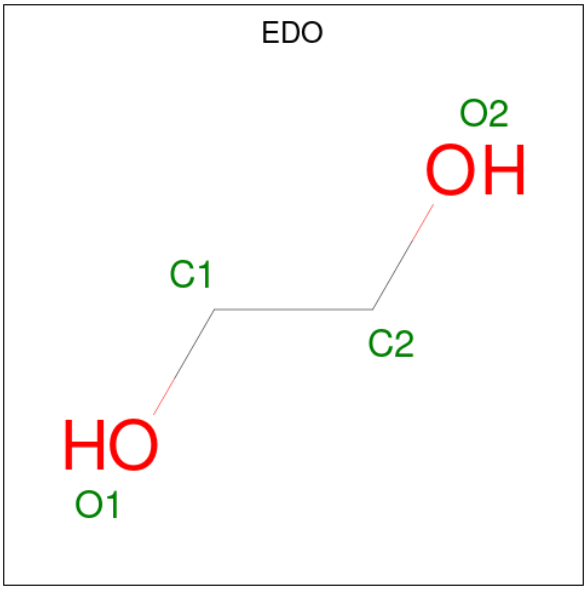
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total C N S 3 1 1 1	0	0

- Molecule 9 is 3-AMINO-1,2,4-TRIAZOLE (three-letter code: 3TR) (formula: $C_2H_4N_4$).



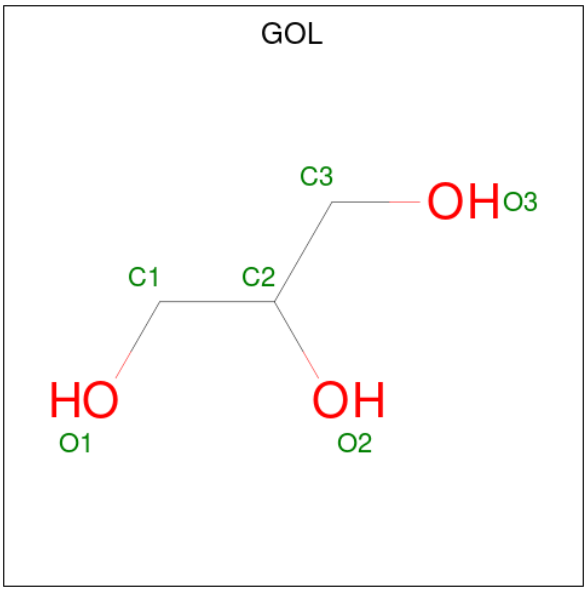
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	N		
9	A	1	6	2	4	0	0
			Total	C	N		
9	A	1	6	2	4	0	0

- Molecule 10 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



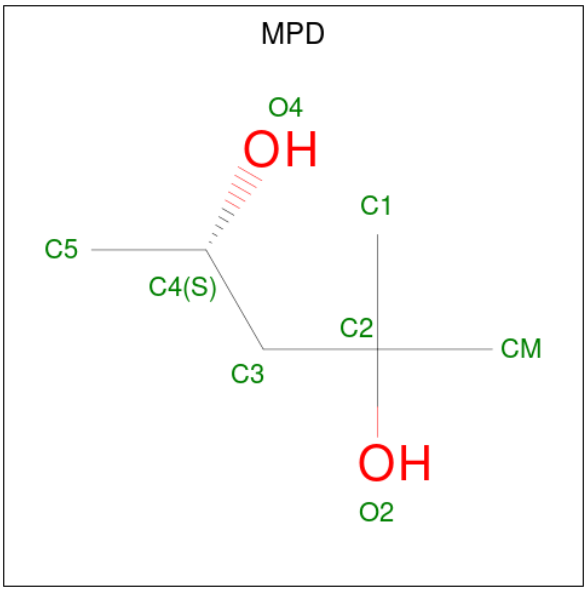
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
10	A	1	4	2	2	0	0
			Total	C	O		
10	A	1	4	2	2	0	0
			Total	C	O		
10	A	1	4	2	2	0	0
			Total	C	O		
10	A	1	4	2	2	0	0
			Total	C	O		
10	A	1	4	2	2	0	0

- Molecule 11 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



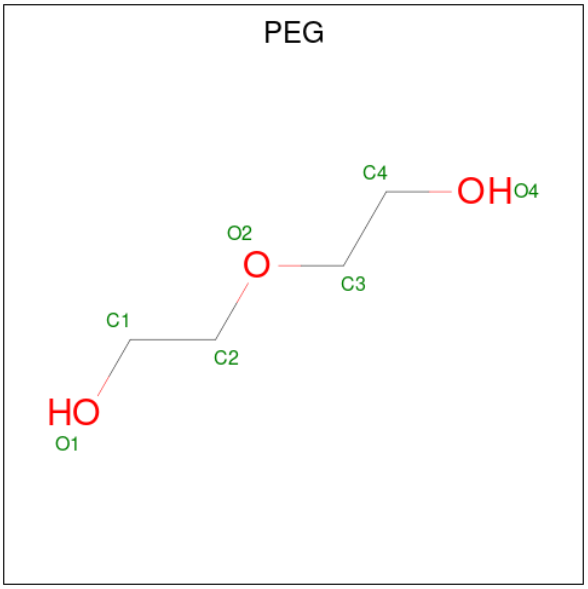
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
11	A	1	6	3	3	0	0
			Total	C	O		
11	A	1	6	3	3	0	0

- Molecule 12 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
12	A	1	8	6	2	0	0

- Molecule 13 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
13	A	1	7	4	3	0	0

- Molecule 14 is water.

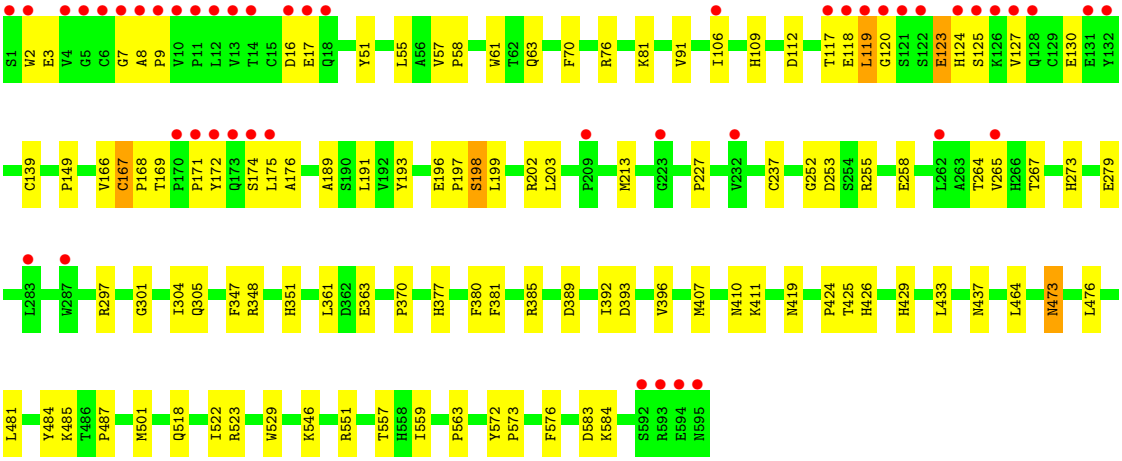
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
14	A	305	305	305	0	0

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.

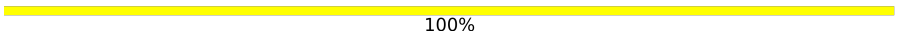
• Molecule 1: Lactoperoxidase

Chain A:



• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	53.99Å 79.62Å 77.52Å 90.00° 102.24° 90.00°	Depositor
Resolution (Å)	75.76 – 2.10 43.98 – 2.10	Depositor EDS
% Data completeness (in resolution range)	98.6 (75.76-2.10) 98.6 (43.98-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	3.58 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, R_{free}	0.212 , 0.257 0.210 , 0.255	Depositor DCC
R_{free} test set	1865 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å^2)	42.1	Xtriage
Anisotropy	0.315	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.28 , 45.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	5252	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.53% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: GOL, CA, EDO, 3TR, MPD, SCN, ZN, SEP, HEM, NAG, PEG, IOD

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	0.44	0/4876	0.60	0/6621

There are no bond length outliers.

There are no bond angle outliers.

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	4758	0	4644	128	0
2	B	28	0	25	0	0
3	A	42	0	39	0	0
4	A	43	0	30	6	0
5	A	1	0	0	0	0
6	A	1	0	0	0	0
7	A	12	0	0	0	0
8	A	3	0	0	0	0
9	A	12	0	8	9	0
10	A	20	0	30	6	0
11	A	12	0	16	4	0
12	A	8	0	14	16	0
13	A	7	0	10	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
14	A	305	0	0	14	0
All	All	5252	0	4816	140	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 14.

All (140) 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:167:CYS:HB2	1:A:168:PRO:CD	1.69	1.21
1:A:167:CYS:CB	1:A:168:PRO:CD	2.37	1.02
1:A:196:GLU:HB3	1:A:198:SEP:O3P	1.58	1.01
1:A:167:CYS:HB2	1:A:168:PRO:HD2	1.41	0.97
1:A:167:CYS:HB2	1:A:168:PRO:HD3	1.45	0.97
1:A:120:GLY:O	1:A:123:GLU:HB2	1.66	0.96
1:A:167:CYS:CB	1:A:168:PRO:HD3	1.96	0.95
1:A:202:ARG:HD3	12:A:621:MPD:H11	1.48	0.95
1:A:81:LYS:HD3	13:A:622:PEG:H42	1.49	0.94
1:A:172:TYR:HE2	1:A:174:SER:HB2	1.30	0.94
1:A:106:ILE:HG13	1:A:265:VAL:HG11	1.57	0.86
9:A:751:3TR:H3A1	9:A:597:3TR:C5	1.89	0.86
1:A:123:GLU:HG2	1:A:125:SER:OG	1.78	0.83
1:A:120:GLY:HA3	1:A:123:GLU:CD	1.99	0.82
1:A:363:GLU:HG3	14:A:829:HOH:O	1.79	0.82
1:A:175:LEU:HD12	1:A:176:ALA:H	1.44	0.82
1:A:172:TYR:CE2	1:A:174:SER:HB2	2.14	0.82
1:A:2:TRP:CD1	1:A:175:LEU:HD13	2.13	0.81
1:A:9:PRO:HB3	14:A:931:HOH:O	1.81	0.80
1:A:407:MET:HB3	1:A:501:MET:CE	2.13	0.78
1:A:2:TRP:HB2	1:A:175:LEU:HD22	1.64	0.77
9:A:751:3TR:N3A	9:A:597:3TR:H5	1.98	0.77
1:A:264:THR:HG23	1:A:392:ILE:HG23	1.69	0.75
1:A:258:GLU:HG3	9:A:597:3TR:HN1	1.50	0.75
1:A:424:PRO:O	1:A:425:THR:HB	1.84	0.74
1:A:279:GLU:HG3	14:A:724:HOH:O	1.87	0.74
1:A:464:LEU:HA	1:A:481:LEU:HD12	1.70	0.71
1:A:199:LEU:CD1	12:A:621:MPD:H13	2.22	0.70
9:A:751:3TR:H3A1	9:A:597:3TR:H5	1.58	0.68
1:A:109:HIS:NE2	9:A:597:3TR:N3A	2.41	0.68
1:A:167:CYS:HB3	1:A:168:PRO:HD3	1.76	0.68
1:A:202:ARG:HD2	12:A:621:MPD:H31	1.75	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:199:LEU:HA	12:A:621:MPD:C1	2.24	0.67
1:A:258:GLU:HG3	9:A:597:3TR:N1	2.10	0.67
4:A:618:HEM:HMC1	4:A:618:HEM:HBC2	1.76	0.66
1:A:119:LEU:HD21	1:A:169:THR:HG22	1.78	0.66
1:A:199:LEU:HA	12:A:621:MPD:H13	1.78	0.65
1:A:63:GLN:HG2	14:A:686:HOH:O	1.98	0.63
1:A:130:GLU:HB2	14:A:908:HOH:O	1.97	0.63
1:A:117:THR:HG21	1:A:119:LEU:HD11	1.81	0.62
1:A:407:MET:HB3	1:A:501:MET:HE3	1.81	0.62
1:A:117:THR:HG21	1:A:119:LEU:CD1	2.30	0.60
1:A:393:ASP:OD1	1:A:557:THR:HB	2.02	0.60
1:A:2:TRP:CB	1:A:175:LEU:HD22	2.30	0.59
1:A:120:GLY:HA3	1:A:123:GLU:CG	2.32	0.59
1:A:551:ARG:HD3	1:A:583:ASP:O	2.03	0.59
1:A:199:LEU:HD12	12:A:621:MPD:H13	1.85	0.59
1:A:473:ASN:ND2	1:A:476:LEU:H	2.01	0.59
1:A:124:HIS:HD2	1:A:127:VAL:HB	1.68	0.58
12:A:621:MPD:H52	12:A:621:MPD:H12	1.85	0.58
1:A:429:HIS:CE1	10:A:598:EDO:H22	2.39	0.58
1:A:139:CYS:HB2	11:A:619:GOL:O1	2.04	0.57
1:A:125:SER:HB2	11:A:619:GOL:H12	1.85	0.57
1:A:169:THR:HG23	14:A:790:HOH:O	2.03	0.57
1:A:213:MET:HG2	1:A:273:HIS:CD2	2.41	0.56
1:A:202:ARG:HH11	12:A:621:MPD:C3	2.18	0.56
1:A:425:THR:HG22	1:A:425:THR:O	2.07	0.55
1:A:118:GLU:HG3	1:A:119:LEU:H	1.72	0.55
1:A:171:PRO:HG3	14:A:921:HOH:O	2.08	0.54
1:A:255:ARG:HD2	9:A:597:3TR:N4	2.22	0.54
1:A:199:LEU:HD13	12:A:621:MPD:H13	1.90	0.54
1:A:361:LEU:HD21	10:A:606:EDO:H21	1.90	0.53
1:A:119:LEU:HD21	1:A:169:THR:CG2	2.39	0.52
1:A:106:ILE:CG1	1:A:265:VAL:HG11	2.35	0.52
1:A:117:THR:CG2	1:A:119:LEU:CD1	2.88	0.52
1:A:377:HIS:CD2	10:A:598:EDO:H11	2.45	0.52
9:A:751:3TR:H3A1	9:A:597:3TR:HN1	1.55	0.51
10:A:600:EDO:H11	14:A:793:HOH:O	2.10	0.51
1:A:202:ARG:CD	12:A:621:MPD:H11	2.29	0.51
4:A:618:HEM:HMB2	4:A:618:HEM:HBB2	1.93	0.51
1:A:202:ARG:HB3	12:A:621:MPD:H11	1.92	0.51
1:A:551:ARG:HB3	14:A:821:HOH:O	2.11	0.51
1:A:2:TRP:O	1:A:3:GLU:C	2.47	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:197:PRO:HA	10:A:603:EDO:H12	1.94	0.50
1:A:149:PRO:HG2	11:A:614:GOL:H2	1.93	0.50
1:A:91:VAL:HG12	1:A:411:LYS:HD2	1.94	0.49
1:A:348:ARG:HH11	1:A:437:ASN:HD22	1.60	0.49
9:A:751:3TR:N3A	9:A:597:3TR:C5	2.57	0.49
1:A:557:THR:OG1	1:A:559:ILE:HG12	2.12	0.49
1:A:407:MET:CB	1:A:501:MET:CE	2.88	0.49
1:A:370:PRO:HD2	14:A:777:HOH:O	2.13	0.49
1:A:523:ARG:HG3	1:A:529:TRP:CE2	2.48	0.48
4:A:618:HEM:HBB2	4:A:618:HEM:CMB	2.43	0.48
1:A:237:CYS:HA	1:A:381:PHE:O	2.13	0.48
1:A:2:TRP:CD2	1:A:175:LEU:HB2	2.48	0.48
1:A:57:VAL:HG13	1:A:58:PRO:HD2	1.95	0.48
1:A:203:LEU:HD11	1:A:252:GLY:HA2	1.95	0.48
1:A:481:LEU:HD21	1:A:487:PRO:HD3	1.96	0.48
1:A:202:ARG:HH11	12:A:621:MPD:H32	1.79	0.48
1:A:377:HIS:HA	1:A:380:PHE:CE2	2.49	0.47
1:A:348:ARG:HH11	1:A:437:ASN:ND2	2.10	0.47
1:A:112:ASP:OD2	4:A:618:HEM:O1D	2.32	0.47
1:A:481:LEU:HA	1:A:484:TYR:O	2.14	0.47
1:A:175:LEU:CD1	1:A:176:ALA:H	2.23	0.47
1:A:119:LEU:HD11	1:A:169:THR:HG22	1.96	0.47
1:A:410:ASN:OD1	14:A:781:HOH:O	2.20	0.46
1:A:426:HIS:HE1	14:A:908:HOH:O	1.98	0.46
1:A:202:ARG:CD	12:A:621:MPD:H31	2.42	0.46
1:A:419:ASN:ND2	11:A:614:GOL:H11	2.31	0.46
1:A:189:ALA:HB2	1:A:304:ILE:HD12	1.98	0.46
1:A:407:MET:HB3	1:A:501:MET:HE1	1.93	0.45
1:A:7:GLY:C	1:A:9:PRO:HD3	2.37	0.45
1:A:51:TYR:CD2	1:A:55:LEU:O	2.69	0.45
1:A:202:ARG:HD3	12:A:621:MPD:C1	2.34	0.45
1:A:301:GLY:O	1:A:305:GLN:HG3	2.17	0.45
1:A:76:ARG:HH22	1:A:419:ASN:HD21	1.63	0.45
1:A:193:TYR:OH	1:A:297:ARG:HA	2.17	0.45
1:A:425:THR:O	1:A:425:THR:CG2	2.64	0.45
1:A:76:ARG:HH12	1:A:419:ASN:ND2	2.14	0.44
1:A:70:PHE:CG	1:A:485:LYS:HB2	2.53	0.44
1:A:546:LYS:NZ	14:A:823:HOH:O	2.50	0.44
1:A:57:VAL:HG13	1:A:61:TRP:CG	2.54	0.43
1:A:166:VAL:O	1:A:168:PRO:HD2	2.18	0.43
1:A:392:ILE:HG22	1:A:396:VAL:HG23	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:227:PRO:HD3	1:A:267:THR:HG23	2.01	0.43
1:A:117:THR:HG23	1:A:169:THR:HG23	2.01	0.43
1:A:361:LEU:CD2	10:A:606:EDO:H21	2.49	0.43
1:A:16:ASP:O	1:A:17:GLU:HB3	2.19	0.42
1:A:174:SER:HA	14:A:923:HOH:O	2.19	0.42
1:A:546:LYS:NZ	1:A:546:LYS:HB2	2.34	0.42
1:A:123:GLU:HG2	1:A:125:SER:HG	1.81	0.42
12:A:621:MPD:H12	12:A:621:MPD:C5	2.50	0.42
1:A:199:LEU:HA	12:A:621:MPD:H12	1.98	0.42
1:A:385:ARG:O	1:A:389:ASP:HB3	2.19	0.42
1:A:191:LEU:HA	1:A:253:ASP:HB2	2.01	0.42
4:A:618:HEM:HBC2	4:A:618:HEM:CMC	2.48	0.42
1:A:464:LEU:CA	1:A:481:LEU:HD12	2.44	0.42
1:A:518:GLN:O	1:A:522:ILE:HG23	2.20	0.42
1:A:76:ARG:HH22	1:A:419:ASN:ND2	2.18	0.41
1:A:264:THR:HA	1:A:392:ILE:HD12	2.02	0.41
1:A:348:ARG:NH2	4:A:618:HEM:HAD1	2.36	0.41
1:A:16:ASP:O	1:A:17:GLU:CB	2.69	0.41
1:A:117:THR:CG2	1:A:119:LEU:HD12	2.51	0.41
1:A:572:TYR:CD1	1:A:573:PRO:HA	2.56	0.40
1:A:473:ASN:C	1:A:473:ASN:HD22	2.24	0.40
1:A:120:GLY:HA3	1:A:123:GLU:OE1	2.18	0.40
1:A:175:LEU:HG	1:A:176:ALA:N	2.36	0.40
1:A:351:HIS:CE1	1:A:433:LEU:HD21	2.56	0.40
1:A:551:ARG:CZ	1:A:584:LYS:HG2	2.52	0.40
1:A:563:PRO:HD3	1:A:576:PHE:CE2	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	592/595 (100%)	569 (96%)	20 (3%)	3 (0%)	29 26

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	167	CYS
1	A	119	LEU
1	A	8	ALA

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	516/516 (100%)	513 (99%)	3 (1%)	86 90

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	123	GLU
1	A	347	PHE
1	A	473	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	124	HIS
1	A	128	GLN
1	A	147	ASN
1	A	329	GLN
1	A	419	ASN
1	A	437	ASN
1	A	468	GLN
1	A	473	ASN
1	A	497	ASN

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Mol	Chain	Res	Type
1	A	570	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	$\#\|Z\| > 2$
1	SEP	A	198	1	8,9,10	1.16	1 (12%)	8,12,14	1.17	1 (12%)

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
1	SEP	A	198	1	-	4/5/8/10	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	198	SEP	P-O1P	2.22	1.57	1.50

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	198	SEP	O2P-P-OG	2.09	112.30	106.73

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	198	SEP	CB-OG-P-O2P
1	A	198	SEP	CB-OG-P-O3P
1	A	198	SEP	CB-OG-P-O1P
1	A	198	SEP	CA-CB-OG-P

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	198	SEP	1	0

5.5 Carbohydrates (i)

2 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	NAG	B	1	2,1	14,14,15	0.73	0	17,19,21	1.10	1 (5%)
2	NAG	B	2	2	14,14,15	0.69	0	17,19,21	2.49	5 (29%)

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	NAG	B	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	B	2	2	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	2	NAG	C1-O5-C5	6.78	121.38	112.19
2	B	2	NAG	C4-C3-C2	-5.15	103.47	111.02
2	B	2	NAG	C1-C2-N2	3.34	116.20	110.49
2	B	1	NAG	O4-C4-C3	-2.25	105.14	110.35
2	B	2	NAG	C3-C4-C5	2.20	114.17	110.24
2	B	2	NAG	O5-C5-C4	2.14	116.03	110.83

There are no chirality outliers.

All (1) torsion outliers are listed below:

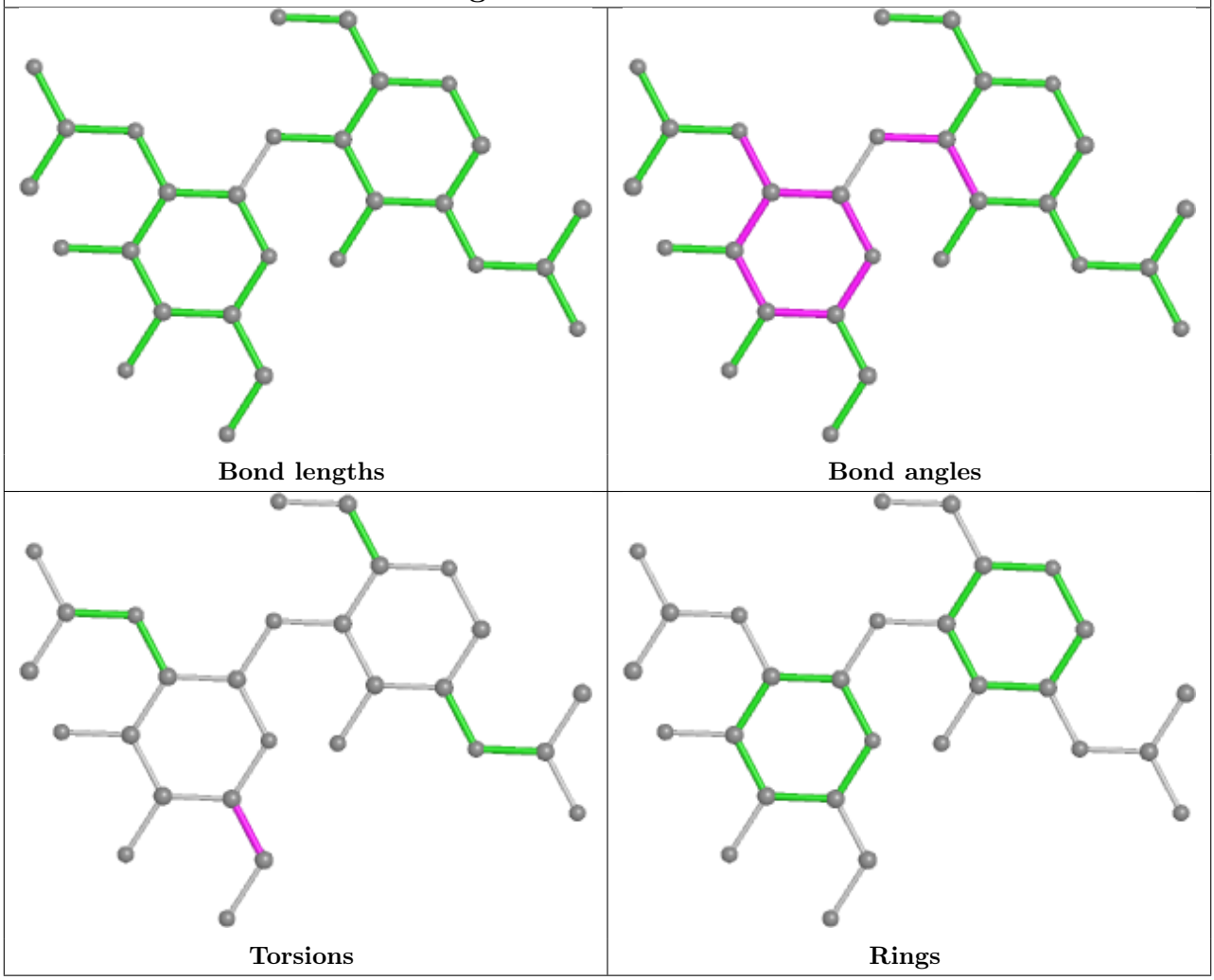
Mol	Chain	Res	Type	Atoms
2	B	2	NAG	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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 B



5.6 Ligand geometry (i)

Of 30 ligands modelled in this entry, 14 are monoatomic - leaving 16 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
8	SCN	A	700	-	1,2,2	10.62	1 (100%)	0,1,1	-	-
9	3TR	A	751	-	4,6,6	2.69	2 (50%)	0,7,7	-	-
11	GOL	A	614	-	5,5,5	0.76	0	5,5,5	1.06	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
10	EDO	A	598	-	3,3,3	0.28	0	2,2,2	0.53	0
10	EDO	A	600	-	3,3,3	0.46	0	2,2,2	0.37	0
13	PEG	A	622	-	6,6,6	0.52	0	5,5,5	0.32	0
10	EDO	A	606	-	3,3,3	0.42	0	2,2,2	0.45	0
3	NAG	A	599	1	14,14,15	0.53	0	17,19,21	0.97	0
12	MPD	A	621	-	7,7,7	1.16	1 (14%)	9,10,10	0.90	0
9	3TR	A	597	-	4,6,6	3.93	3 (75%)	0,7,7	-	-
3	NAG	A	604	1	14,14,15	0.55	0	17,19,21	1.66	3 (17%)
10	EDO	A	605	-	3,3,3	0.39	0	2,2,2	0.39	0
3	NAG	A	596	1	14,14,15	0.53	0	17,19,21	0.93	2 (11%)
11	GOL	A	619	-	5,5,5	0.32	0	5,5,5	0.54	0
4	HEM	A	618	1	41,50,50	1.94	8 (19%)	45,82,82	1.86	9 (20%)
10	EDO	A	603	-	3,3,3	0.48	0	2,2,2	0.26	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
11	GOL	A	614	-	-	4/4/4/4	-
9	3TR	A	751	-	-	-	0/1/1/1
10	EDO	A	598	-	-	1/1/1/1	-
10	EDO	A	600	-	-	1/1/1/1	-
13	PEG	A	622	-	-	3/4/4/4	-
10	EDO	A	606	-	-	1/1/1/1	-
3	NAG	A	599	1	-	0/6/23/26	0/1/1/1
12	MPD	A	621	-	-	2/5/5/5	-
9	3TR	A	597	-	-	-	0/1/1/1
3	NAG	A	604	1	-	2/6/23/26	0/1/1/1
10	EDO	A	605	-	-	0/1/1/1	-
3	NAG	A	596	1	-	0/6/23/26	0/1/1/1
11	GOL	A	619	-	-	3/4/4/4	-
4	HEM	A	618	1	-	4/12/54/54	-
10	EDO	A	603	-	-	0/1/1/1	-

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	A	700	SCN	C-N	10.62	1.51	1.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	618	HEM	C3D-C2D	6.66	1.50	1.36
9	A	597	3TR	C3-N3A	5.02	1.43	1.33
4	A	618	HEM	C3C-C2C	-4.74	1.33	1.40
9	A	597	3TR	C3-N4	-4.64	1.27	1.34
9	A	751	3TR	C3-N3A	4.62	1.43	1.33
9	A	597	3TR	C5-N1	-3.82	1.25	1.32
4	A	618	HEM	C3C-CAC	3.66	1.55	1.47
4	A	618	HEM	CAB-C3B	2.94	1.55	1.47
12	A	621	MPD	O2-C2	-2.88	1.37	1.44
9	A	751	3TR	C3-N4	-2.58	1.30	1.34
4	A	618	HEM	FE-ND	2.55	2.09	1.96
4	A	618	HEM	CAA-C2A	2.52	1.55	1.52
4	A	618	HEM	FE-NB	2.27	2.08	1.96
4	A	618	HEM	C1B-NB	-2.27	1.36	1.40

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	618	HEM	C4D-ND-C1D	6.04	111.31	105.07
3	A	604	NAG	O5-C5-C6	4.22	113.82	107.20
4	A	618	HEM	C3B-C2B-C1B	3.82	109.32	106.49
4	A	618	HEM	CMD-C2D-C1D	3.70	130.67	125.04
4	A	618	HEM	C1B-NB-C4B	3.37	108.56	105.07
4	A	618	HEM	C4C-CHD-C1D	3.06	126.60	122.56
4	A	618	HEM	CAD-C3D-C4D	3.01	129.91	124.66
3	A	604	NAG	O5-C1-C2	-2.64	107.11	111.29
4	A	618	HEM	C2B-C1B-NB	-2.53	106.84	109.84
3	A	604	NAG	O3-C3-C4	-2.48	104.61	110.35
4	A	618	HEM	CBD-CAD-C3D	-2.45	105.82	112.63
4	A	618	HEM	C4B-CHC-C1C	2.44	125.78	122.56
3	A	596	NAG	C1-O5-C5	2.24	115.23	112.19
3	A	596	NAG	O5-C1-C2	-2.20	107.81	111.29

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
11	A	619	GOL	C1-C2-C3-O3
13	A	622	PEG	O1-C1-C2-O2
3	A	604	NAG	O5-C5-C6-O6
3	A	604	NAG	C4-C5-C6-O6
11	A	614	GOL	C1-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
11	A	619	GOL	O2-C2-C3-O3
10	A	598	EDO	O1-C1-C2-O2
10	A	600	EDO	O1-C1-C2-O2
11	A	614	GOL	O1-C1-C2-O2
11	A	614	GOL	O1-C1-C2-C3
12	A	621	MPD	C2-C3-C4-C5
13	A	622	PEG	C4-C3-O2-C2
11	A	614	GOL	O2-C2-C3-O3
4	A	618	HEM	CAD-CBD-CGD-O2D
12	A	621	MPD	O2-C2-C3-C4
4	A	618	HEM	CAD-CBD-CGD-O1D
4	A	618	HEM	CAA-CBA-CGA-O2A
10	A	606	EDO	O1-C1-C2-O2
4	A	618	HEM	CAA-CBA-CGA-O1A
11	A	619	GOL	O1-C1-C2-C3
13	A	622	PEG	O2-C3-C4-O4

There are no ring outliers.

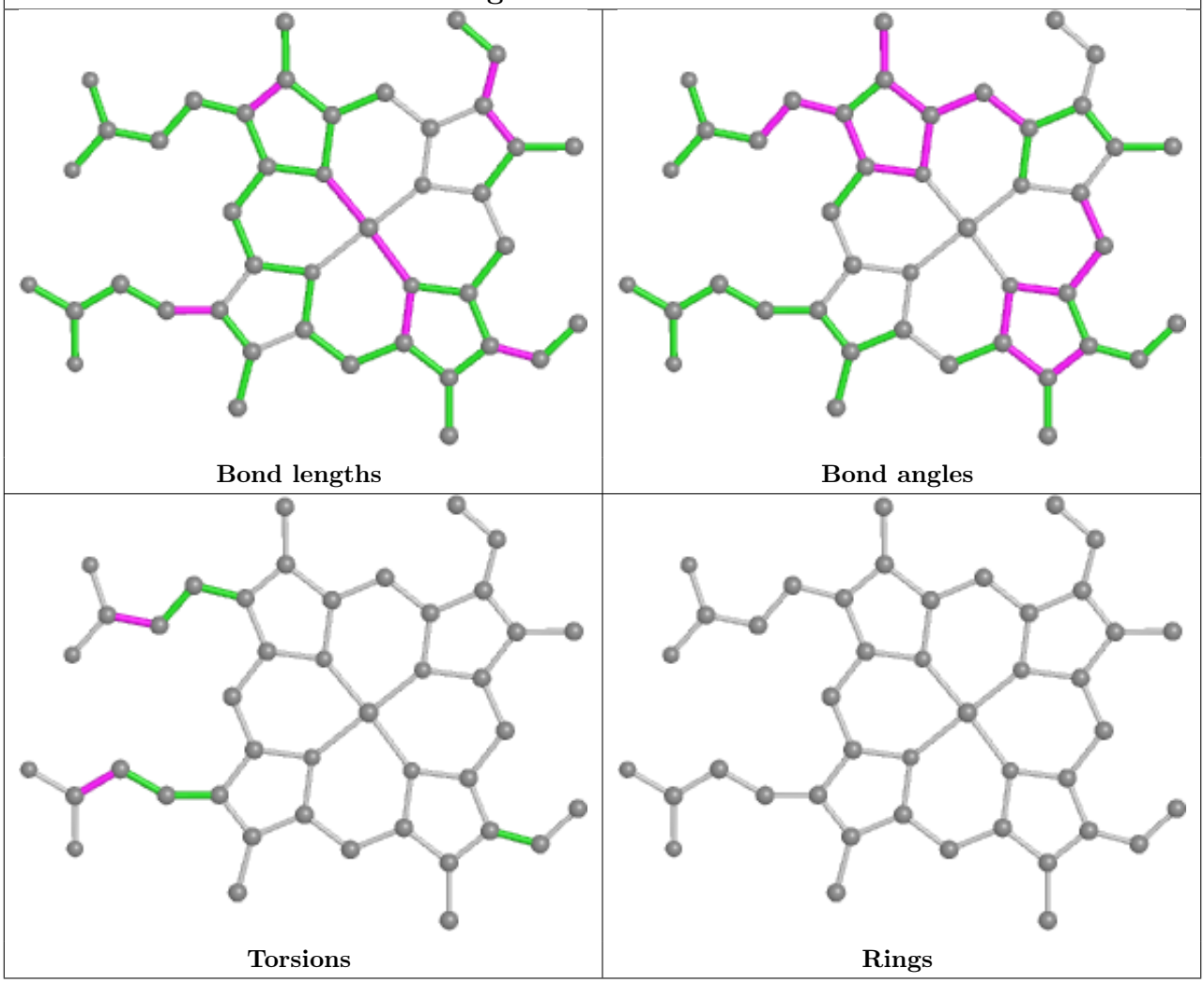
11 monomers are involved in 42 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	A	751	3TR	5	0
11	A	614	GOL	2	0
10	A	598	EDO	2	0
10	A	600	EDO	1	0
13	A	622	PEG	1	0
10	A	606	EDO	2	0
12	A	621	MPD	16	0
9	A	597	3TR	9	0
11	A	619	GOL	2	0
4	A	618	HEM	6	0
10	A	603	EDO	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and

any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

Ligand HEM A 618



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	594/595 (99%)	0.48	47 (7%) 12 16	29, 42, 72, 88	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	12	LEU	13.2
1	A	172	TYR	10.8
1	A	2	TRP	9.4
1	A	1	SER	9.1
1	A	121	SER	9.1
1	A	173	GLN	8.5
1	A	11	PRO	8.3
1	A	595	ASN	7.3
1	A	119	LEU	7.2
1	A	4	VAL	7.2
1	A	13	VAL	6.2
1	A	174	SER	6.2
1	A	122	SER	6.0
1	A	10	VAL	5.9
1	A	120	GLY	5.6
1	A	171	PRO	5.4
1	A	593	ARG	5.3
1	A	7	GLY	5.1
1	A	5	GLY	5.0
1	A	14	THR	4.9
1	A	283	LEU	4.7
1	A	170	PRO	4.6
1	A	124	HIS	4.3
1	A	9	PRO	4.2
1	A	18	GLN	4.1
1	A	265	VAL	3.9
1	A	594	GLU	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	118	GLU	3.8
1	A	16	ASP	3.7
1	A	125	SER	3.7
1	A	6	CYS	3.6
1	A	132	TYR	3.5
1	A	223	GLY	3.2
1	A	209	PRO	3.1
1	A	175	LEU	3.1
1	A	126	LYS	3.0
1	A	128	GLN	2.9
1	A	106	ILE	2.6
1	A	117	THR	2.5
1	A	287	TRP	2.5
1	A	262	LEU	2.4
1	A	127	VAL	2.2
1	A	232	VAL	2.2
1	A	592	SER	2.2
1	A	17	GLU	2.1
1	A	131	GLU	2.0
1	A	8	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (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
1	SEP	A	198	10/11	0.89	0.17	40,44,47,49	0

6.3 Carbohydrates (i)

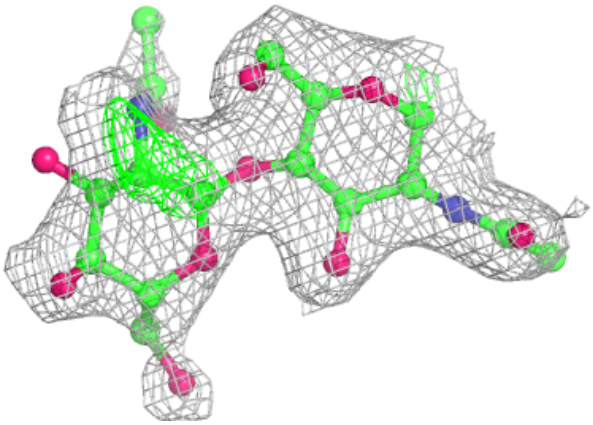
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

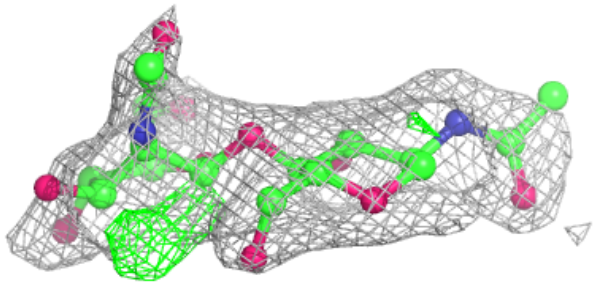
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	NAG	B	2	14/15	0.77	0.21	61,66,67,71	14
2	NAG	B	1	14/15	0.92	0.13	49,52,55,56	0

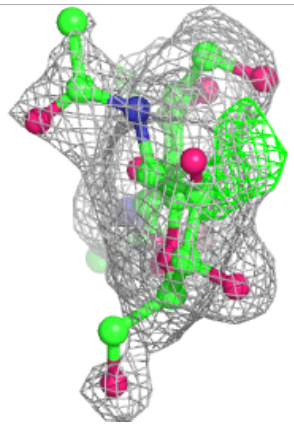
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain B:

$2mF_o$-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
11	GOL	A	619	6/6	0.74	0.22	67,68,68,68	0
10	EDO	A	603	4/4	0.77	0.24	33,37,39,40	0
3	NAG	A	599	14/15	0.81	0.14	57,63,65,68	0
10	EDO	A	606	4/4	0.83	0.18	34,34,38,41	0
3	NAG	A	596	14/15	0.83	0.33	57,61,63,63	0
3	NAG	A	604	14/15	0.84	0.15	59,63,67,68	14
12	MPD	A	621	8/8	0.86	0.34	18,34,38,39	0
13	PEG	A	622	7/7	0.86	0.17	60,60,62,62	7
10	EDO	A	598	4/4	0.87	0.25	36,38,40,42	0
9	3TR	A	751	6/6	0.87	0.21	43,44,45,45	0

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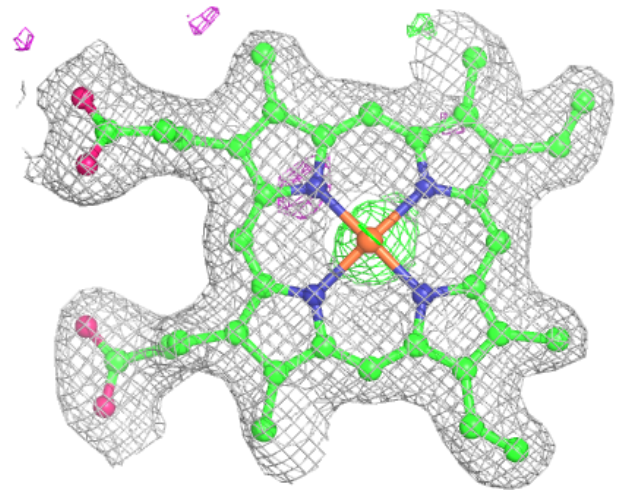
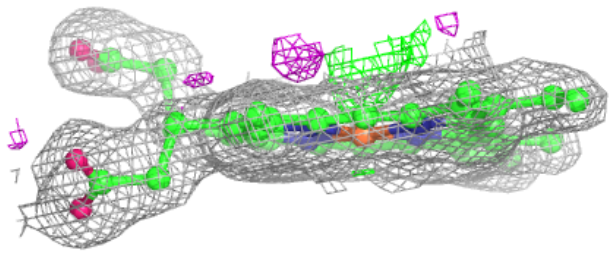
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
11	GOL	A	614	6/6	0.90	0.20	41,44,44,46	0
10	EDO	A	605	4/4	0.91	0.18	47,47,49,51	0
10	EDO	A	600	4/4	0.92	0.13	46,48,49,50	0
9	3TR	A	597	6/6	0.92	0.30	29,32,33,34	0
7	IOD	A	627	1/1	0.93	0.07	59,59,59,59	1
6	ZN	A	608	1/1	0.94	0.04	63,63,63,63	0
8	SCN	A	700	3/3	0.95	0.19	59,59,59,60	0
7	IOD	A	620	1/1	0.96	0.07	60,60,60,60	1
7	IOD	A	613	1/1	0.96	0.07	67,67,67,67	1
5	CA	A	607	1/1	0.97	0.12	35,35,35,35	0
4	HEM	A	618	43/43	0.97	0.18	25,28,33,35	0
7	IOD	A	611	1/1	0.97	0.04	60,60,60,60	0
7	IOD	A	624	1/1	0.98	0.08	55,55,55,55	1
7	IOD	A	626	1/1	0.98	0.04	57,57,57,57	1
7	IOD	A	610	1/1	0.98	0.06	58,58,58,58	1
7	IOD	A	617	1/1	0.98	0.04	58,58,58,58	1
7	IOD	A	612	1/1	0.98	0.04	50,50,50,50	0
7	IOD	A	616	1/1	0.99	0.04	57,57,57,57	1
7	IOD	A	615	1/1	0.99	0.08	46,46,46,46	0
7	IOD	A	609	1/1	1.00	0.09	34,34,34,34	0

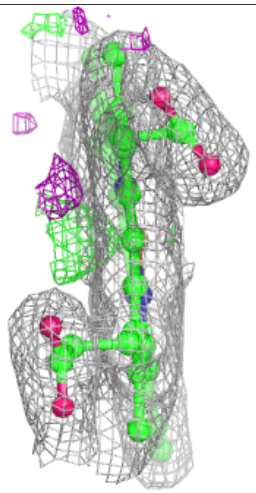
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around HEM A 618:

$2mF_o$-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.