



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 20, 2020 – 02:05 PM BST

PDB ID : 4GM7
Title : Structure of cinnamic acid bound bovine lactoperoxidase at 2.6Å resolution.
Authors : Pandey, N.; Singh, A.K.; Sinha, M.; Kaur, P.; Sharma, S.; Singh, T.P.
Deposited on : 2012-08-15
Resolution : 2.60 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 : 2.13.1
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) : 2.13.1

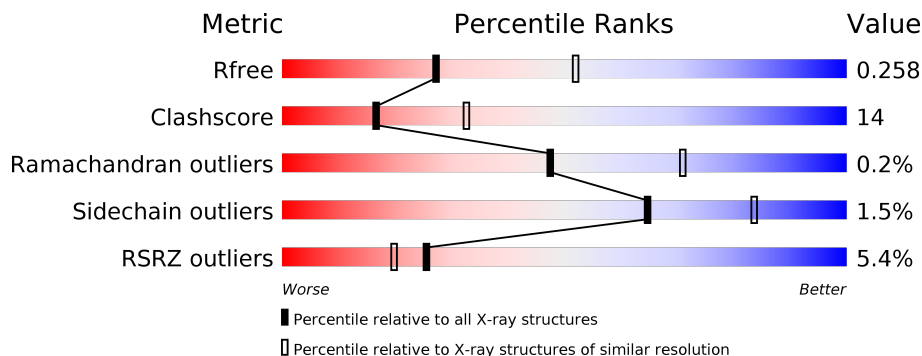
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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 on 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	595	
2	B	2	

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
11	HCl	A	622	-	-	X	-

2 Entry composition [i](#)

There are 12 unique types of molecules in this entry. The entry contains 5078 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.

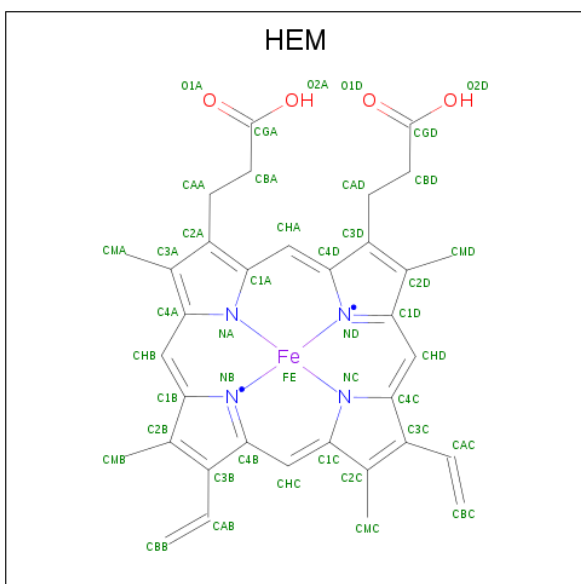
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	A	595	4775	3037	847	864	1	26	0	0	0

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



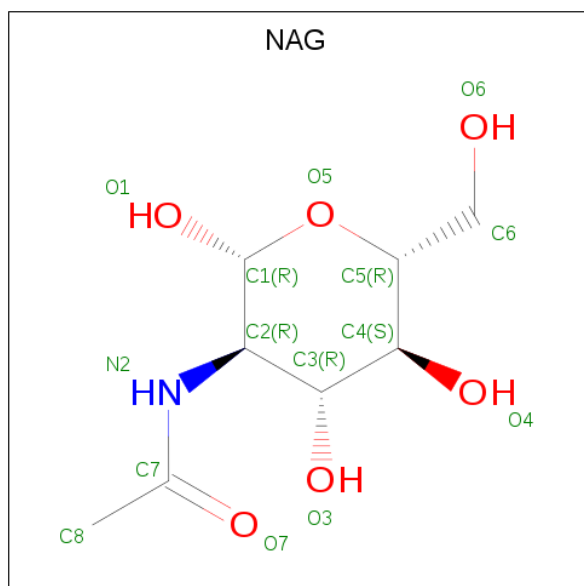
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	B	2	28	16	2	10	0	0	0

- Molecule 3 is 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
3	A	1	43	34	1	4	4	0	0

- Molecule 4 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		
4	A	1	14	8	1	5	0	0
4	A	1	14	8	1	5	0	0
4	A	1	14	8	1	5	0	0

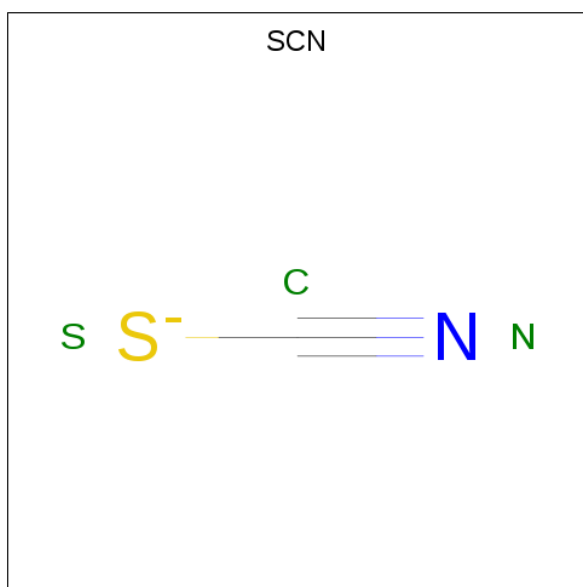
- Molecule 5 is IODIDE ION (three-letter code: IOD) (formula: I).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	8	Total I 8 8	0	0

- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 7 is THIOCYANATE ION (three-letter code: SCN) (formula: CNS).



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

- Molecule 8 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



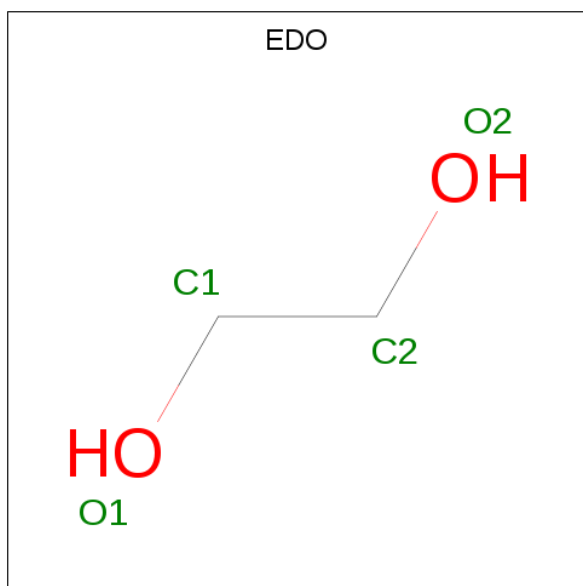
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
8	A	1	7	4	3	0	0

- Molecule 9 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



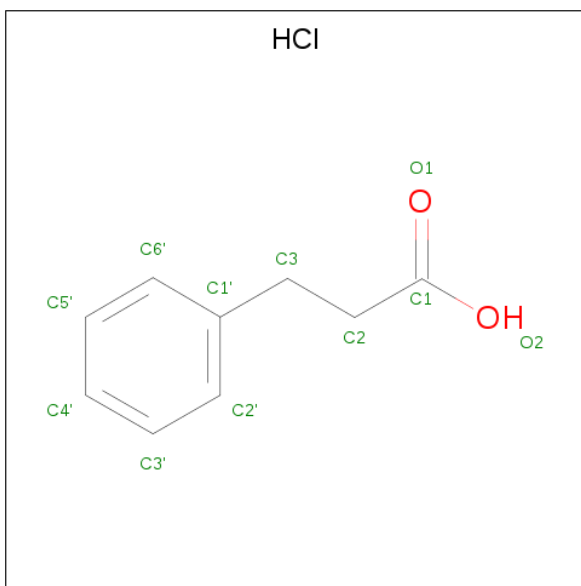
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 10 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

- Molecule 11 is HYDROCINNAMIC ACID (three-letter code: HCI) (formula: C₉H₁₀O₂).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	1	Total	C O	0	0
			11	9 2		

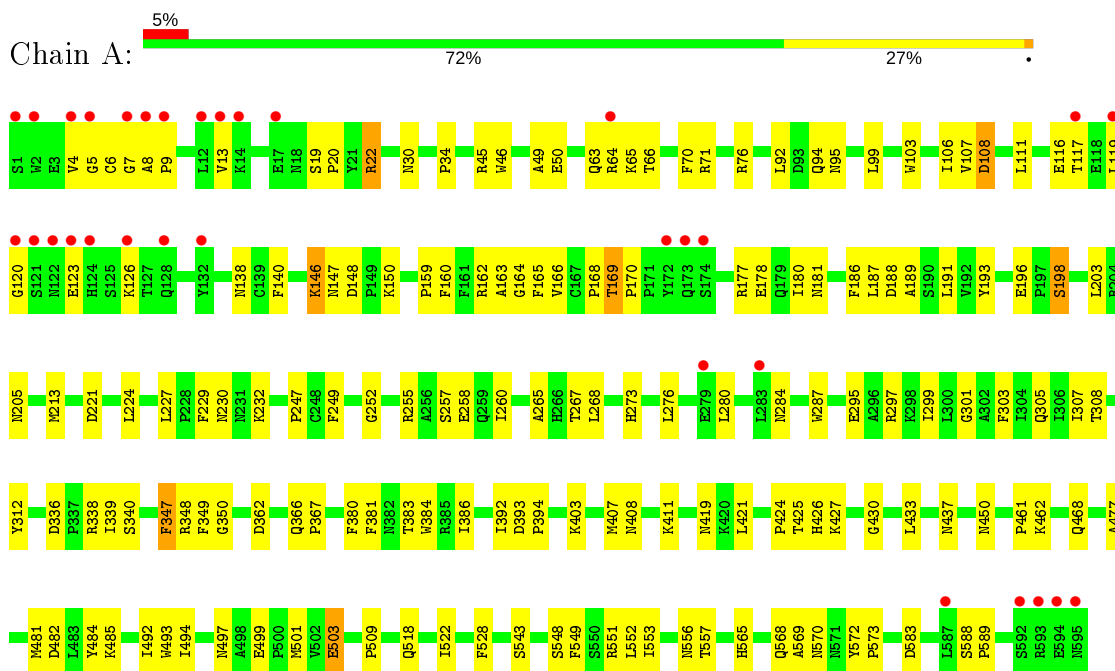
- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	142	Total	O	0	0
			142	142		

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: Lactoperoxidase



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	54.05Å 80.23Å 76.04Å 90.00° 103.30° 90.00°	Depositor
Resolution (Å)	43.99 – 2.60 43.99 – 2.60	Depositor EDS
% Data completeness (in resolution range)	97.8 (43.99-2.60) 97.9 (43.99-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.73 (at 2.61Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.221 , 0.235 0.218 , 0.258	Depositor DCC
R_{free} test set	983 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	43.9	Xtrriage
Anisotropy	0.901	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 35.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5078	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.98% of the height of the origin peak. No significant pseudotranslation is detected.*

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: GOL, SCN, NAG, SEP, CA, EDO, PEG, HEM, HCI, 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	0.42	0/4892	0.67	0/6634

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	4775	0	4685	132	0
2	B	28	0	25	1	0
3	A	43	0	30	6	0
4	A	42	0	39	1	0
5	A	8	0	0	0	0
6	A	1	0	0	0	0
7	A	3	0	0	1	0
8	A	7	0	9	0	0
9	A	6	0	8	3	0
10	A	12	0	15	0	0
11	A	11	0	9	8	0
12	A	142	0	0	5	0
All	All	5078	0	4820	138	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 (138) 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:168:PRO:HB2	1:A:169:THR:OG1	1.71	0.91
1:A:482:ASP:O	1:A:485:LYS:HE2	1.69	0.91
1:A:350:GLY:HA3	3:A:601:HEM:CBC	2.02	0.89
1:A:381:PHE:HZ	11:A:622:HCl:H21	1.38	0.87
1:A:518:GLN:O	1:A:522:ILE:HG12	1.77	0.84
1:A:213:MET:HG2	1:A:273:HIS:CD2	2.17	0.79
1:A:150:LYS:HE2	1:A:419:ASN:HD22	1.48	0.79
1:A:381:PHE:CZ	11:A:622:HCl:H21	2.18	0.79
1:A:551:ARG:HD3	1:A:583:ASP:O	1.84	0.78
1:A:350:GLY:HA3	3:A:601:HEM:HBC2	1.67	0.77
1:A:20:PRO:O	1:A:528:PHE:HE2	1.68	0.76
1:A:258:GLU:HB2	11:A:622:HCl:H6'	1.68	0.74
1:A:339:ILE:HD13	1:A:522:ILE:HD11	1.68	0.73
1:A:95:ASN:HD22	9:A:618:GOL:H12	1.51	0.73
11:A:622:HCl:H4'	12:A:840:HOH:O	1.89	0.72
1:A:9:PRO:HG3	1:A:168:PRO:HD3	1.72	0.71
1:A:348:ARG:HH11	1:A:437:ASN:ND2	1.89	0.70
1:A:117:THR:CG2	1:A:164:GLY:HA2	2.21	0.69
3:A:601:HEM:HMB1	3:A:601:HEM:HBB2	1.73	0.69
1:A:120:GLY:HA2	1:A:126:LYS:HG2	1.75	0.68
1:A:339:ILE:CD1	1:A:522:ILE:HD11	2.24	0.68
1:A:588:SER:N	1:A:589:PRO:HD2	2.09	0.68
1:A:462:LYS:HD2	12:A:734:HOH:O	1.93	0.67
1:A:503:GLU:O	1:A:503:GLU:HG3	1.93	0.67
1:A:543:SER:OG	1:A:589:PRO:HG2	1.95	0.65
1:A:63:GLN:HG3	1:A:71:ARG:NH2	2.12	0.64
1:A:117:THR:HG22	1:A:164:GLY:HA2	1.80	0.63
1:A:588:SER:N	1:A:589:PRO:CD	2.60	0.63
1:A:169:THR:H	1:A:170:PRO:CD	2.13	0.61
1:A:425:THR:O	1:A:425:THR:HG22	2.00	0.61
1:A:588:SER:HB2	1:A:589:PRO:HD3	1.83	0.60
1:A:117:THR:OG1	1:A:138:ASN:ND2	2.35	0.59
1:A:366:GLN:O	1:A:367:PRO:C	2.40	0.58
1:A:419:ASN:O	1:A:430:GLY:HA2	2.03	0.58
4:A:602:NAG:O7	9:A:618:GOL:H11	2.04	0.58
1:A:20:PRO:O	1:A:528:PHE:CE2	2.55	0.58
1:A:146:LYS:HG3	1:A:147:ASN:N	2.16	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:424:PRO:O	1:A:425:THR:HB	2.04	0.57
1:A:116:GLU:HA	1:A:163:ALA:HA	1.86	0.57
1:A:392:ILE:HG13	1:A:393:ASP:N	2.17	0.57
1:A:295:GLU:O	1:A:299:ILE:HG12	2.04	0.57
1:A:8:ALA:HB3	1:A:9:PRO:HD3	1.87	0.57
1:A:150:LYS:CE	1:A:419:ASN:HD22	2.18	0.56
1:A:301:GLY:O	1:A:305:GLN:HG3	2.05	0.56
1:A:393:ASP:HB2	12:A:740:HOH:O	2.05	0.56
1:A:30:ASN:O	1:A:34:PRO:HA	2.06	0.55
1:A:393:ASP:HB2	1:A:394:PRO:HD3	1.89	0.55
1:A:66:THR:HB	1:A:70:PHE:O	2.07	0.55
1:A:383:THR:HB	12:A:721:HOH:O	2.06	0.55
1:A:407:MET:HB3	1:A:501:MET:CE	2.36	0.55
1:A:568:GLN:HE21	1:A:570:ASN:HD21	1.53	0.55
1:A:106:ILE:HG23	1:A:191:LEU:HD11	1.88	0.54
1:A:258:GLU:O	1:A:380:PHE:HA	2.07	0.53
1:A:63:GLN:HG3	1:A:71:ARG:HH22	1.74	0.53
1:A:159:PRO:HG3	1:A:426:HIS:CE1	2.44	0.53
1:A:481:MET:O	1:A:484:TYR:O	2.26	0.52
1:A:76:ARG:HH22	1:A:419:ASN:HD21	1.56	0.52
1:A:268:LEU:HB3	1:A:552:LEU:HD21	1.90	0.52
1:A:108:ASP:OD1	1:A:108:ASP:C	2.48	0.52
1:A:188:ASP:O	1:A:189:ALA:HB3	2.09	0.51
1:A:276:LEU:O	1:A:280:LEU:HG	2.11	0.51
1:A:205:ASN:OD1	1:A:205:ASN:C	2.49	0.50
1:A:249:PHE:CZ	1:A:383:THR:HG23	2.47	0.50
1:A:150:LYS:HE2	1:A:419:ASN:ND2	2.22	0.50
1:A:92:LEU:HD13	1:A:403:LYS:HB2	1.93	0.50
1:A:549:PHE:O	1:A:553:ILE:HG12	2.12	0.49
3:A:601:HEM:C1A	11:A:622:HCl:H3'	2.48	0.49
1:A:120:GLY:CA	1:A:126:LYS:HG2	2.42	0.49
1:A:169:THR:H	1:A:170:PRO:HD2	1.77	0.49
1:A:117:THR:HG23	1:A:164:GLY:HA2	1.96	0.48
1:A:380:PHE:CE2	1:A:421:LEU:HA	2.49	0.48
1:A:257:SER:O	1:A:381:PHE:HA	2.14	0.48
1:A:499:GLU:OE1	1:A:509:PRO:HD2	2.13	0.48
1:A:255:ARG:HB3	11:A:622:HCl:C6'	2.43	0.48
1:A:168:PRO:HB2	1:A:169:THR:HG1	1.77	0.48
1:A:230:ASN:ND2	7:A:616:SCN:S	2.88	0.47
1:A:148:ASP:OD1	1:A:150:LYS:HB2	2.14	0.47
1:A:227:LEU:CD2	1:A:267:THR:HA	2.43	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:A:622:HCl:H6'	11:A:622:HCl:H22	1.43	0.47
1:A:19:SER:HA	1:A:20:PRO:HD3	1.70	0.47
1:A:303:PHE:O	1:A:307:ILE:HG12	2.15	0.47
1:A:260:ILE:HD11	1:A:386:ILE:HG13	1.96	0.47
1:A:6:CYS:SG	1:A:7:GLY:N	2.88	0.47
1:A:99:LEU:HD21	1:A:549:PHE:CD2	2.50	0.46
1:A:362:ASP:OD2	1:A:366:GLN:HB2	2.15	0.46
1:A:177:ARG:HA	12:A:713:HOH:O	2.15	0.45
1:A:196:GLU:HB3	1:A:198:SEP:O3P	2.15	0.45
1:A:203:LEU:HD11	1:A:252:GLY:HA2	1.98	0.45
1:A:76:ARG:HH22	1:A:419:ASN:ND2	2.14	0.45
1:A:107:VAL:O	1:A:111:LEU:HG	2.17	0.45
1:A:229:PHE:CD1	1:A:247:PRO:HD2	2.51	0.45
1:A:13:VAL:HG11	1:A:34:PRO:HB2	1.99	0.45
1:A:492:ILE:HG23	1:A:493:TRP:N	2.32	0.45
1:A:565:HIS:O	1:A:568:GLN:HG2	2.18	0.44
1:A:4:VAL:HG13	1:A:5:GLY:N	2.33	0.44
1:A:425:THR:CG2	1:A:425:THR:O	2.65	0.44
1:A:138:ASN:O	1:A:162:ARG:HG3	2.18	0.44
1:A:347:PHE:O	1:A:347:PHE:CD1	2.70	0.44
1:A:119:LEU:O	1:A:123:GLU:HB2	2.18	0.44
1:A:186:PHE:O	1:A:188:ASP:N	2.51	0.44
1:A:103:TRP:O	1:A:107:VAL:HG23	2.18	0.43
1:A:193:TYR:OH	1:A:297:ARG:HA	2.17	0.43
1:A:95:ASN:ND2	9:A:618:GOL:H12	2.26	0.43
1:A:187:LEU:HA	1:A:187:LEU:HD12	1.76	0.43
1:A:336:ASP:OD2	1:A:338:ARG:NH2	2.38	0.43
1:A:468:GLN:HG2	1:A:477:ALA:HB3	2.00	0.43
1:A:45:ARG:CZ	1:A:49:ALA:HB2	2.48	0.43
1:A:349:PHE:CB	1:A:497:ASN:HD21	2.30	0.43
1:A:140:PHE:O	1:A:160:PHE:HB3	2.18	0.43
1:A:572:TYR:HA	1:A:573:PRO:HA	1.79	0.43
1:A:284:ASN:ND2	1:A:287:TRP:CZ2	2.87	0.42
1:A:180:ILE:CG2	1:A:181:ASN:N	2.82	0.42
1:A:384:TRP:CH2	2:B:1:NAG:H2	2.55	0.42
1:A:548:SER:OG	1:A:551:ARG:HB2	2.19	0.42
3:A:601:HEM:C2A	11:A:622:HCl:H3'	2.54	0.42
1:A:494:ILE:HD12	1:A:494:ILE:HA	1.87	0.42
1:A:549:PHE:CE2	1:A:553:ILE:HD11	2.55	0.42
1:A:166:VAL:HG22	1:A:178:GLU:O	2.20	0.41
1:A:308:THR:O	1:A:312:TYR:HB3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:433:LEU:HD12	1:A:433:LEU:HA	1.72	0.41
1:A:94:GLN:O	1:A:569:ALA:CB	2.68	0.41
1:A:64:ARG:NH1	1:A:65:LYS:HE2	2.35	0.41
1:A:221:ASP:O	1:A:224:LEU:HB2	2.20	0.41
1:A:165:PHE:CZ	1:A:169:THR:O	2.73	0.41
1:A:408:ASN:HB3	1:A:411:LYS:HB2	2.02	0.41
1:A:551:ARG:HD2	1:A:551:ARG:HA	1.71	0.41
1:A:180:ILE:HG22	1:A:181:ASN:N	2.34	0.41
1:A:556:ASN:O	1:A:557:THR:HG23	2.21	0.41
1:A:22:ARG:H	1:A:22:ARG:HG3	1.59	0.41
1:A:70:PHE:CG	1:A:485:LYS:HB2	2.56	0.41
1:A:362:ASP:OD1	1:A:362:ASP:C	2.59	0.41
1:A:450:ASN:OD1	1:A:461:PRO:HD2	2.20	0.40
1:A:350:GLY:CA	3:A:601:HEM:HBC2	2.46	0.40
1:A:106:ILE:HD11	1:A:265:ALA:HB3	2.03	0.40
1:A:46:TRP:CE2	1:A:340:SER:HB3	2.56	0.40
1:A:227:LEU:HD21	1:A:267:THR:HA	2.02	0.40
1:A:393:ASP:CB	1:A:394:PRO:HD3	2.52	0.40
1:A:349:PHE:HB2	1:A:497:ASN:HD21	1.86	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%)	550 (93%)	41 (7%)	1 (0%)	47 71

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	169	THR

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	517/517 (100%)	509 (98%)	8 (2%)	65 83

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	ARG
1	A	50	GLU
1	A	108	ASP
1	A	146	LYS
1	A	232	LYS
1	A	347	PHE
1	A	427	LYS
1	A	503	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	138	ASN
1	A	364	ASN
1	A	419	ASN
1	A	426	HIS
1	A	437	ASN
1	A	468	GLN
1	A	497	ASN
1	A	521	GLN
1	A	570	ASN
1	A	574	HIS

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	0.81	0	8,12,14	1.50	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	-	0/5/8/10	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	198	SEP	OG-CB-CA	3.10	111.16	108.14

There are no chirality outliers.

There are no torsion outliers.

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	1,2	14,14,15	2.25	5 (35%)	17,19,21	1.98	6 (35%)
2	NAG	B	2	2	14,14,15	1.05	1 (7%)	17,19,21	2.17	4 (23%)

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

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	NAG	O4-C4	-5.96	1.29	1.43
2	B	1	NAG	O5-C5	-4.08	1.35	1.43
2	B	1	NAG	C4-C5	-2.69	1.47	1.53
2	B	1	NAG	C3-C2	-2.38	1.47	1.52
2	B	2	NAG	C3-C2	-2.27	1.47	1.52
2	B	1	NAG	O5-C1	-2.00	1.40	1.43

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	NAG	C4-C3-C2	-6.17	101.97	111.02
2	B	1	NAG	O4-C4-C5	-4.03	99.29	109.30
2	B	2	NAG	C1-O5-C5	4.02	117.64	112.19
2	B	1	NAG	C3-C4-C5	3.58	116.62	110.24
2	B	1	NAG	C1-O5-C5	-3.12	107.97	112.19
2	B	2	NAG	C3-C4-C5	3.06	115.70	110.24
2	B	1	NAG	O4-C4-C3	-2.37	104.87	110.35
2	B	1	NAG	C4-C3-C2	-2.10	107.94	111.02
2	B	2	NAG	O5-C1-C2	2.05	114.53	111.29
2	B	1	NAG	O5-C5-C4	-2.04	105.85	110.83

There are no chirality outliers.

All (3) torsion outliers are listed below:

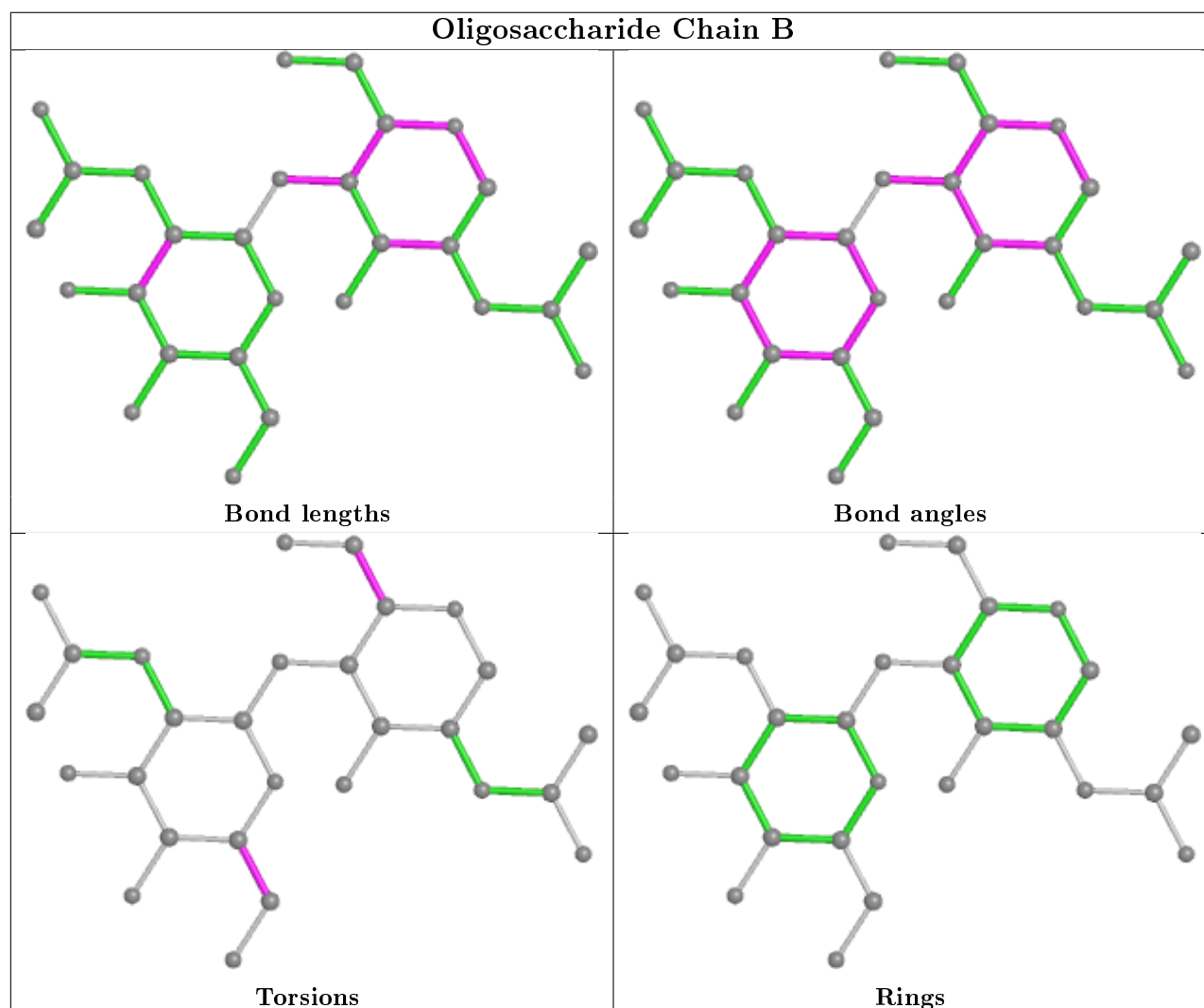
Mol	Chain	Res	Type	Atoms
2	B	2	NAG	C4-C5-C6-O6
2	B	1	NAG	C4-C5-C6-O6
2	B	1	NAG	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1	NAG	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.



5.6 Ligand geometry [i](#)

Of 20 ligands modelled in this entry, 9 are monoatomic - leaving 11 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
11	HCI	A	622	-	8,11,11	0.14	0	10,13,13	0.77	0
4	NAG	A	602	1	14,14,15	1.17	2 (14%)	17,19,21	1.12	1 (5%)
10	EDO	A	621	-	3,3,3	1.67	1 (33%)	2,2,2	0.95	0
3	HEM	A	601	1,12	27,50,50	2.36	12 (44%)	17,82,82	1.80	6 (35%)
4	NAG	A	606	1	14,14,15	1.10	2 (14%)	17,19,21	0.99	1 (5%)
8	PEG	A	617	-	6,6,6	1.29	1 (16%)	5,5,5	1.11	1 (20%)
10	EDO	A	620	-	3,3,3	1.57	1 (33%)	2,2,2	0.88	0
9	GOL	A	618	-	5,5,5	0.75	0	5,5,5	0.43	0
7	SCN	A	616	-	1,2,2	3.54	1 (100%)	0,1,1	0.00	-
4	NAG	A	603	1	14,14,15	1.15	2 (14%)	17,19,21	1.04	1 (5%)
10	EDO	A	619	-	3,3,3	1.45	1 (33%)	2,2,2	0.65	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	HCI	A	622	-	-	3/3/5/5	0/1/1/1
4	NAG	A	602	1	-	0/6/23/26	0/1/1/1
10	EDO	A	621	-	-	1/1/1/1	-
3	HEM	A	601	1,12	-	0/6/54/54	-
4	NAG	A	606	1	-	0/6/23/26	0/1/1/1
8	PEG	A	617	-	-	2/4/4/4	-
10	EDO	A	620	-	-	1/1/1/1	-
9	GOL	A	618	-	-	1/4/4/4	-
4	NAG	A	603	1	-	0/6/23/26	0/1/1/1
10	EDO	A	619	-	-	1/1/1/1	-

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	HEM	C3B-CAB	5.20	1.58	1.47
3	A	601	HEM	C3C-CAC	5.17	1.58	1.47
3	A	601	HEM	CAD-C3D	3.61	1.58	1.52
7	A	616	SCN	C-N	3.54	1.27	1.15
3	A	601	HEM	C1A-CHA	-3.52	1.31	1.41
3	A	601	HEM	C3B-C2B	-3.35	1.35	1.40
3	A	601	HEM	CAA-C2A	3.27	1.56	1.52
3	A	601	HEM	C3C-C2C	-3.03	1.36	1.40
10	A	621	EDO	O2-C2	-2.84	1.27	1.42
4	A	603	NAG	O5-C5	-2.82	1.37	1.43
8	A	617	PEG	O4-C4	-2.75	1.27	1.42
10	A	620	EDO	O2-C2	-2.70	1.28	1.42
4	A	606	NAG	O5-C5	-2.68	1.38	1.43
4	A	602	NAG	O5-C5	-2.62	1.38	1.43
3	A	601	HEM	C3D-C2D	2.58	1.45	1.37
4	A	602	NAG	O5-C1	-2.51	1.39	1.43
4	A	606	NAG	O5-C1	-2.48	1.39	1.43
3	A	601	HEM	C4B-CHC	-2.48	1.34	1.41
10	A	619	EDO	O2-C2	-2.41	1.29	1.42
3	A	601	HEM	C4D-C3D	2.23	1.47	1.42
4	A	603	NAG	O5-C1	-2.20	1.40	1.43
3	A	601	HEM	C1C-C2C	2.20	1.47	1.42
3	A	601	HEM	CMC-C2C	2.11	1.56	1.51

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	HEM	CMD-C2D-C1D	4.25	135.00	128.46
4	A	602	NAG	C4-C3-C2	-2.97	106.67	111.02
3	A	601	HEM	CMD-C2D-C3D	-2.91	119.46	124.94
4	A	603	NAG	O5-C1-C2	2.28	114.89	111.29
8	A	617	PEG	O2-C3-C4	2.26	120.02	110.07
4	A	606	NAG	C2-N2-C7	-2.22	119.73	122.90
3	A	601	HEM	C4C-C3C-C2C	-2.21	105.36	106.90
3	A	601	HEM	CAD-CBD-CGD	2.10	116.20	112.67
3	A	601	HEM	C1D-C2D-C3D	-2.09	105.54	107.00
3	A	601	HEM	CAD-C3D-C2D	-2.08	121.28	127.25

There are no chirality outliers.

All (9) torsion outliers are listed below:

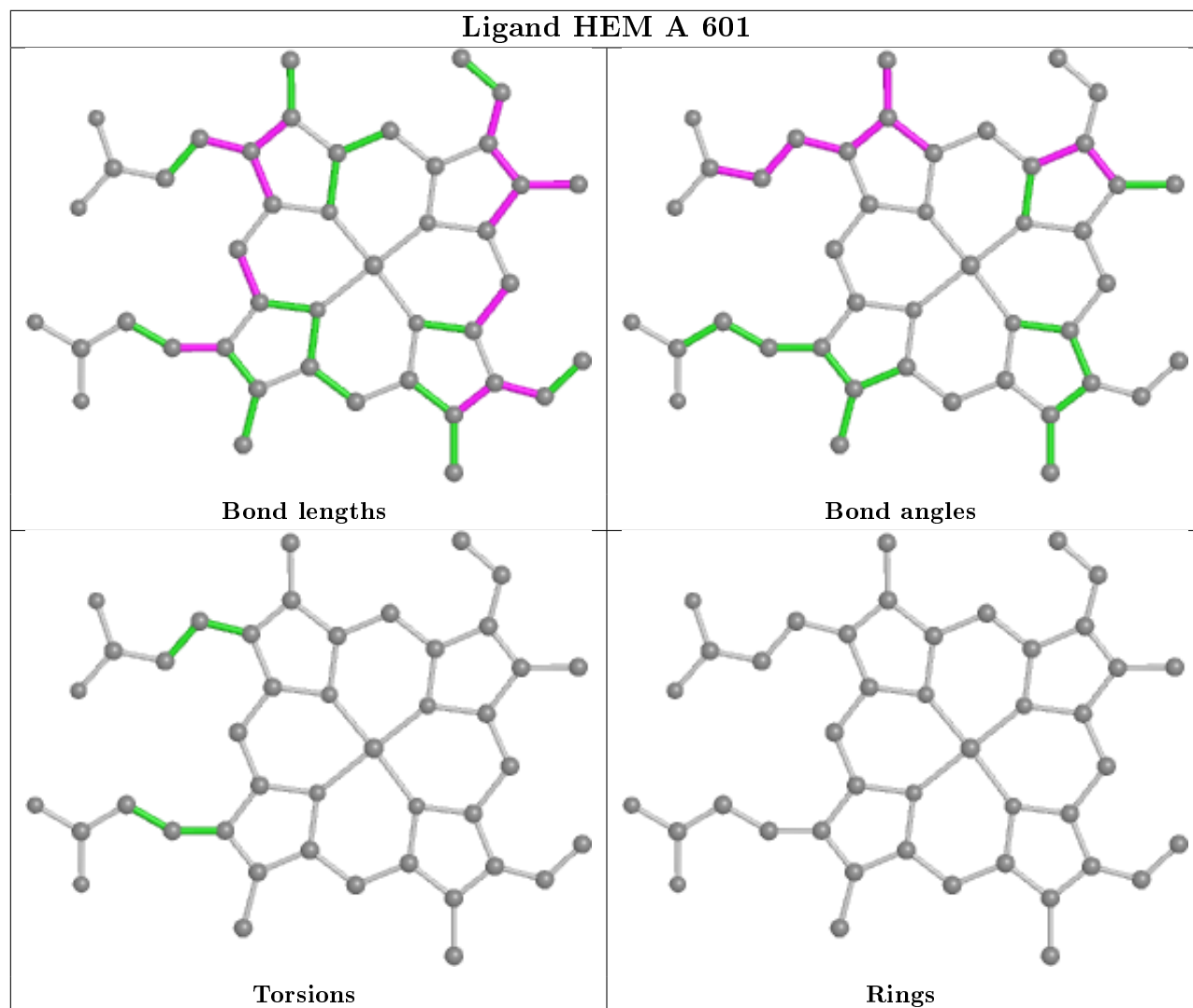
Mol	Chain	Res	Type	Atoms
11	A	622	HCI	C1-C2-C3-C1'
8	A	617	PEG	O2-C3-C4-O4
10	A	621	EDO	O1-C1-C2-O2
10	A	620	EDO	O1-C1-C2-O2
8	A	617	PEG	O1-C1-C2-O2
9	A	618	GOL	O2-C2-C3-O3
10	A	619	EDO	O1-C1-C2-O2
11	A	622	HCI	C6'-C1'-C3-C2
11	A	622	HCI	C2'-C1'-C3-C2

There are no ring outliers.

5 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	A	622	HCI	8	0
4	A	602	NAG	1	0
3	A	601	HEM	6	0
9	A	618	GOL	3	0
7	A	616	SCN	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

6.1 Protein, DNA and RNA chains

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%)	0.05	32 (5%) 25 20	28, 48, 93, 163	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	173	GLN	11.2
1	A	2	TRP	10.4
1	A	7	GLY	9.2
1	A	1	SER	8.4
1	A	121	SER	7.7
1	A	122	ASN	7.4
1	A	12	LEU	6.8
1	A	595	ASN	6.6
1	A	120	GLY	6.4
1	A	8	ALA	6.1
1	A	174	SER	6.1
1	A	13	VAL	4.4
1	A	4	VAL	4.4
1	A	594	GLU	4.1
1	A	172	TYR	3.2
1	A	119	LEU	3.1
1	A	587	LEU	3.0
1	A	64	ARG	2.9
1	A	17	GLU	2.9
1	A	279	GLU	2.9
1	A	124	HIS	2.7
1	A	123	GLU	2.6
1	A	592	SER	2.6
1	A	5	GLY	2.5
1	A	283	LEU	2.5
1	A	593	ARG	2.5
1	A	126	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	132	TYR	2.3
1	A	14	LYS	2.2
1	A	117	THR	2.2
1	A	128	GLN	2.1
1	A	9	PRO	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.

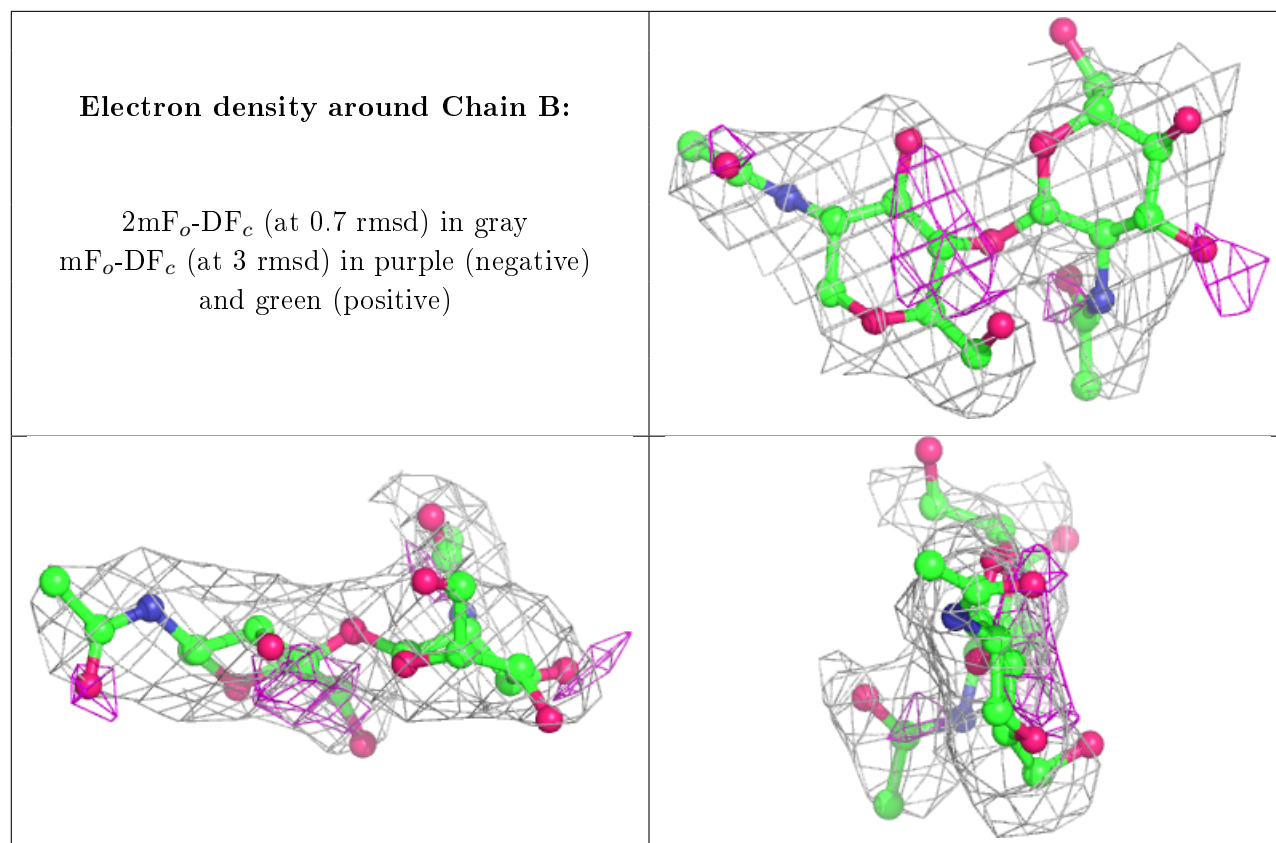
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	SEP	A	198	10/11	0.81	0.27	49,54,58,59	0

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	0.80	0.31	59,66,73,73	0
2	NAG	B	1	14/15	0.81	0.22	43,47,54,55	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.



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.

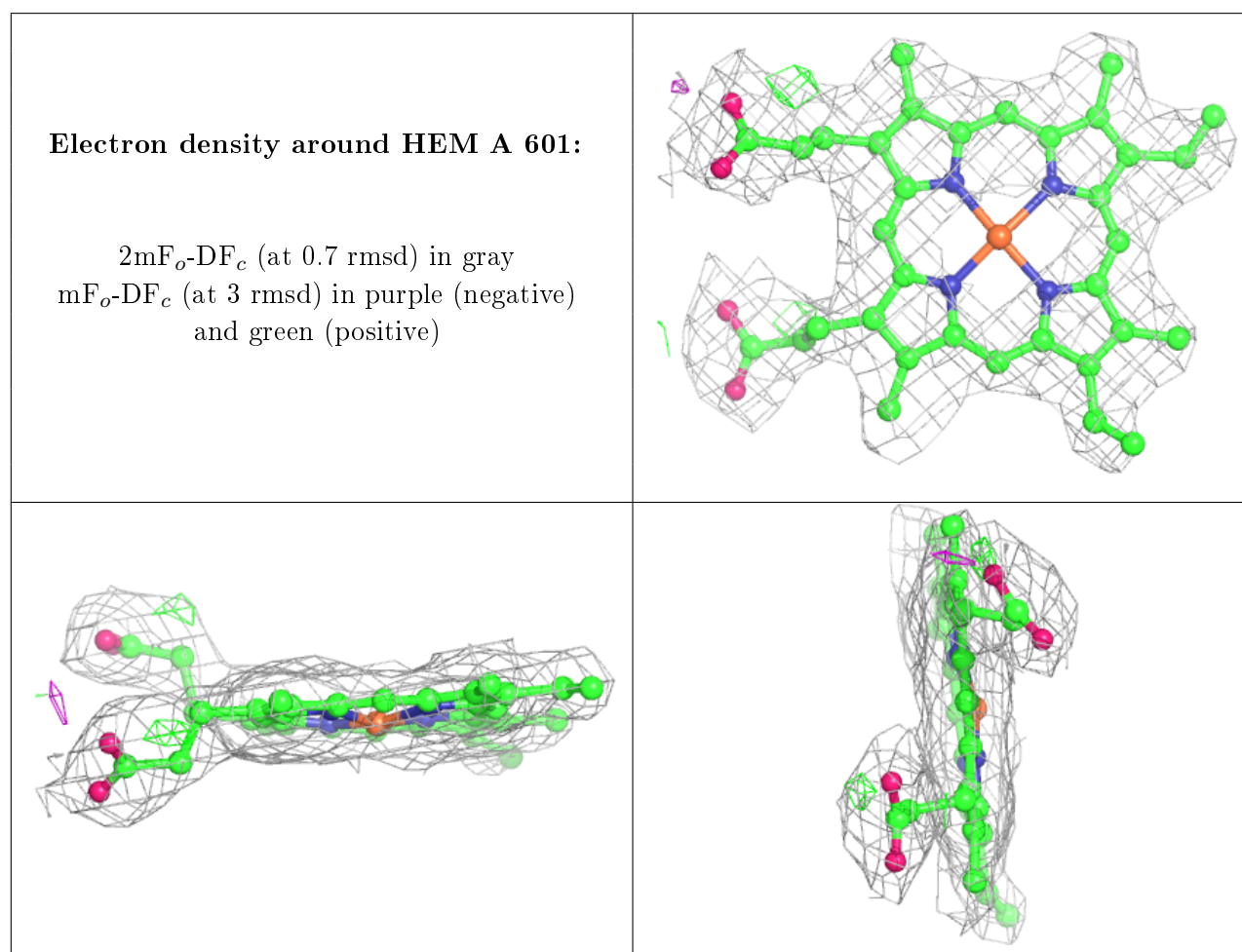
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	NAG	A	602	14/15	0.76	0.34	53,67,77,86	0
4	NAG	A	606	14/15	0.79	0.27	49,59,63,64	0
10	EDO	A	619	4/4	0.79	0.29	70,73,75,75	0
8	PEG	A	617	7/7	0.82	0.31	67,71,73,73	0
11	HCl	A	622	11/11	0.87	0.29	31,37,45,46	0
4	NAG	A	603	14/15	0.88	0.17	41,51,53,54	0
9	GOL	A	618	6/6	0.90	0.15	47,49,51,51	0
10	EDO	A	621	4/4	0.90	0.24	61,61,62,62	0
10	EDO	A	620	4/4	0.91	0.17	48,50,51,52	0
5	IOD	A	611	1/1	0.95	0.08	72,72,72,72	1
3	HEM	A	601	43/43	0.96	0.20	25,30,34,37	0
5	IOD	A	610	1/1	0.96	0.08	72,72,72,72	1
5	IOD	A	614	1/1	0.96	0.05	86,86,86,86	0
7	SCN	A	616	3/3	0.97	0.18	57,57,57,59	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	IOD	A	613	1/1	0.97	0.08	67,67,67,67	0
5	IOD	A	607	1/1	0.98	0.05	90,90,90,90	0
6	CA	A	615	1/1	0.98	0.17	40,40,40,40	0
5	IOD	A	612	1/1	0.99	0.11	49,49,49,49	0
5	IOD	A	609	1/1	0.99	0.06	73,73,73,73	0
5	IOD	A	608	1/1	1.00	0.05	79,79,79,79	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.