Full wwPDB X-ray Structure Validation Report (i)

May 23, 2024 – 02:41 PM EDT

PDB ID	:	4BQ5
Title	:	Structural analysis of an exo-beta-agarase
Authors	:	Pluvinage, B.; Hehemann, J.H.; Boraston, A.B.
Deposited on	:	2013-05-29
Resolution	:	2.30 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:

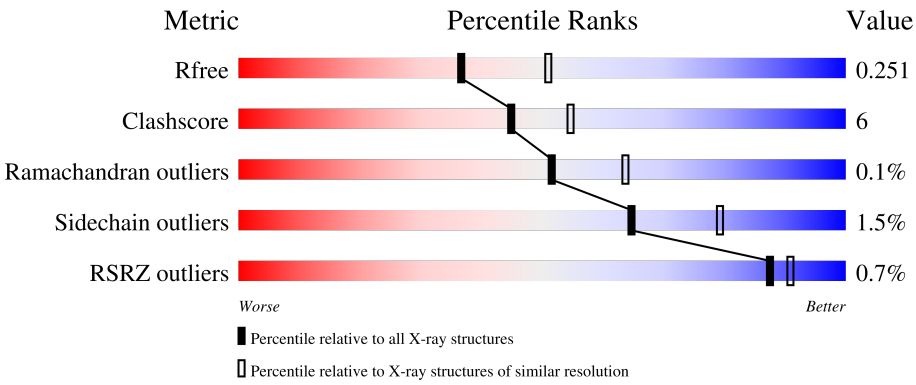
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	750	% 86% 13%
1	B	750	% 84% 16%
2	C	2	50% 50%
2	D	2	50% 50%
2	E	2	50% 50%

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Mol	Chain	Length	Quality of chain
2	G	2	100%
2	H	2	50% 50%
3	F	4	50% 50%

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
4	AAL	A	1795	-	-	X	-
5	GAL	A	1796	-	-	X	-

2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 12994 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.

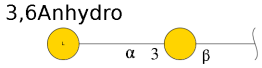
- Molecule 1 is a protein called B-AGARASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	748	5920	3776	991	1130	23	3	3	0
1	B	749	5912	3770	997	1122	23	14	2	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	44	GLY	-	expression tag	UNP Q21HC5
A	45	SER	-	expression tag	UNP Q21HC5
A	46	HIS	-	expression tag	UNP Q21HC5
A	534	GLN	GLU	engineered mutation	UNP Q21HC5
B	44	GLY	-	expression tag	UNP Q21HC5
B	45	SER	-	expression tag	UNP Q21HC5
B	46	HIS	-	expression tag	UNP Q21HC5
B	534	GLN	GLU	engineered mutation	UNP Q21HC5

- Molecule 2 is an oligosaccharide called 3,6-anhydro-alpha-L-galactopyranose-(1-3)-beta-D-galactopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
2	C	2	21	12	9	0	0	0
2	D	2	21	12	9	0	0	0
2	E	2	22	12	10	0	0	0

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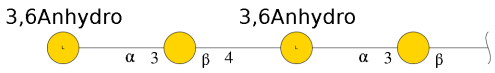
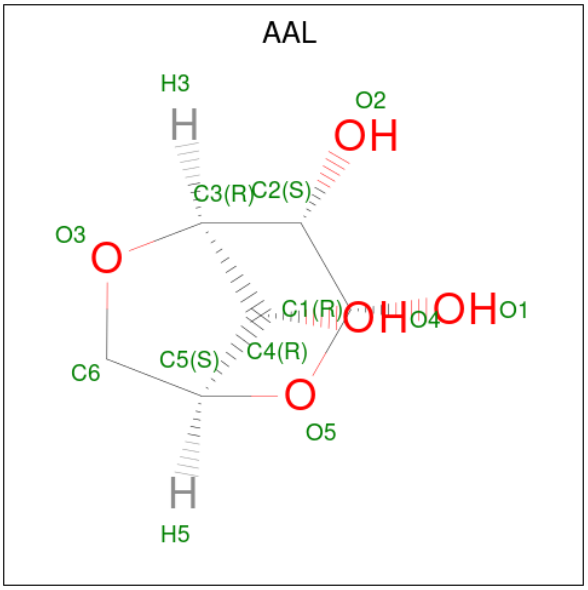
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
2	G	2	21	12	9	0	0	0
			Total	C	O			
2	H	2	22	12	10	0	0	0

- Molecule 3 is an oligosaccharide called 3,6-anhydro-alpha-L-galactopyranose-(1-3)-beta-D-galactopyranose-(1-4)-3,6-anhydro-alpha-L-galactopyranose-(1-3)-beta-D-galactopyranose.



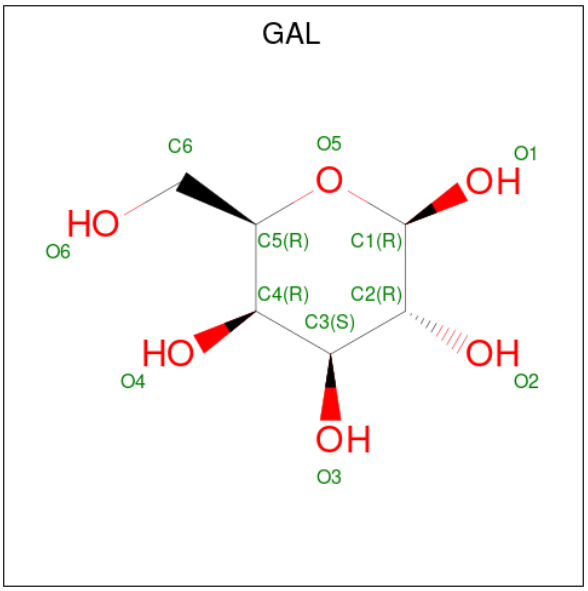
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
3	F	4	42	24	18	0	0	0

- Molecule 4 is 3,6-anhydro-alpha-L-galactopyranose (three-letter code: AAL) (formula: $C_6H_{10}O_5$).



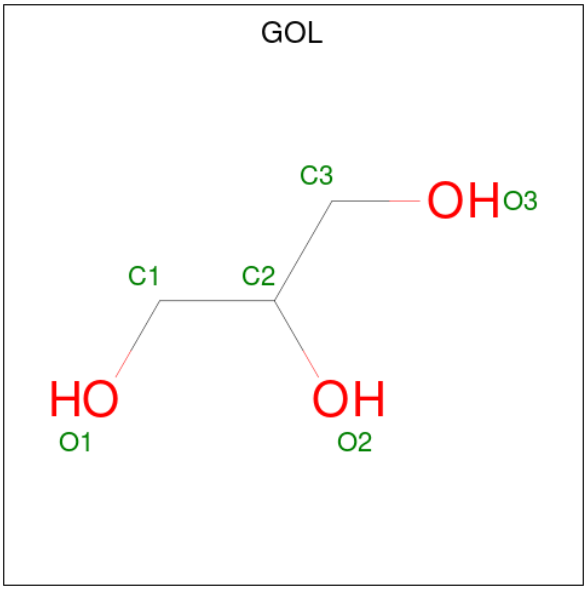
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
4	A	1	10	6	4	0	0

- Molecule 5 is beta-D-galactopyranose (three-letter code: GAL) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
5	A	1	11	6	5	0	0

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

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

- Molecule 7 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Ca		
7	A	3	3	3	0	0
7	B	3	Total	Ca	0	0
			3	3		

- Molecule 8 is water.

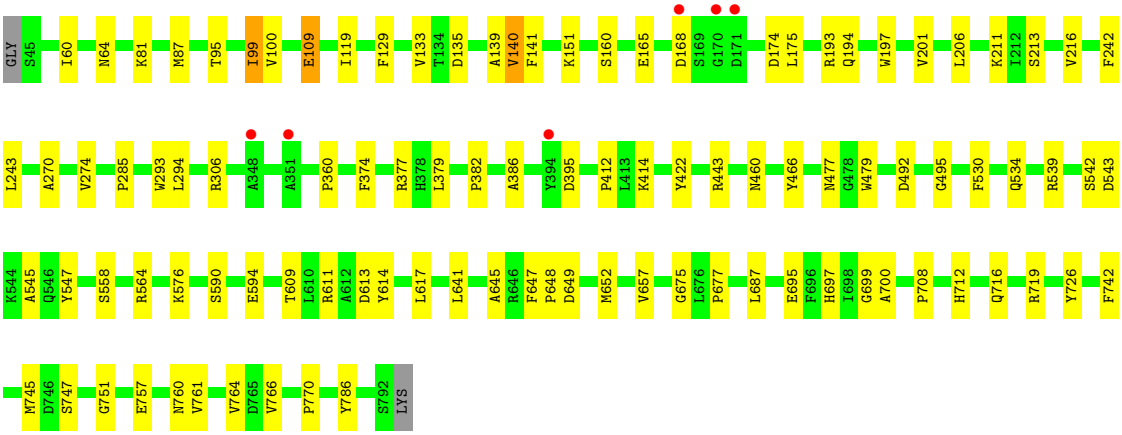
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
8	A	512	512	512	0	0
8	B	450	Total	O	0	0
			450	450		

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.

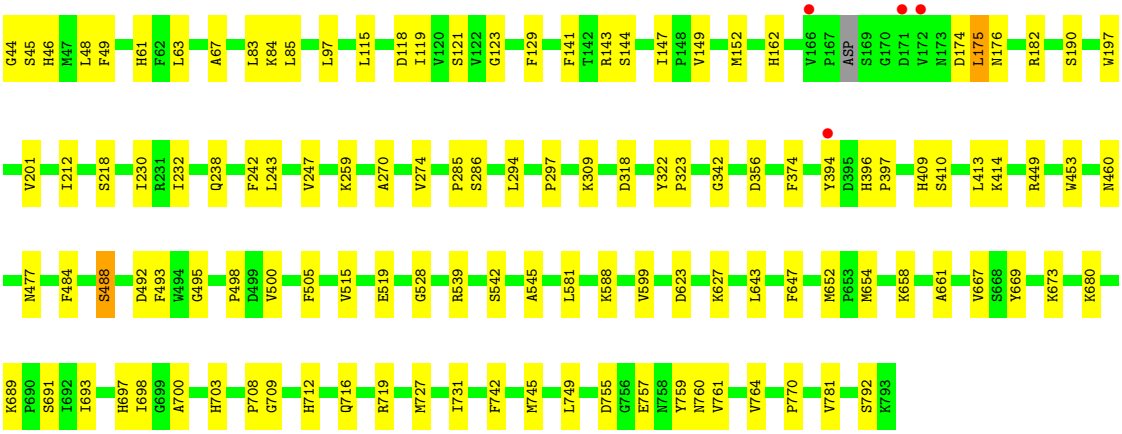
• Molecule 1: B-AGARASE

Chain A:



• Molecule 1: B-AGARASE

Chain B:



• Molecule 2: 3,6-anhydro-alpha-L-galactopyranose-(1-3)-beta-D-galactopyranose

Chain C:

• Molecule 2: 3,6-anhydro-alpha-L-galactopyranose-(1-3)-beta-D-galactopyranose

Chain D:

• Molecule 2: 3,6-anhydro-alpha-L-galactopyranose-(1-3)-beta-D-galactopyranose

Chain E:

• Molecule 2: 3,6-anhydro-alpha-L-galactopyranose-(1-3)-beta-D-galactopyranose

Chain G:

• Molecule 2: 3,6-anhydro-alpha-L-galactopyranose-(1-3)-beta-D-galactopyranose

Chain H:

• Molecule 3: 3,6-anhydro-alpha-L-galactopyranose-(1-3)-beta-D-galactopyranose-(1-4)-3,6-anhydro-alpha-L-galactopyranose-(1-3)-beta-D-galactopyranose

Chain F:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	69.03Å 116.17Å 207.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.19 – 2.30 39.16 – 2.30	Depositor EDS
% Data completeness (in resolution range)	94.1 (39.19-2.30) 94.1 (39.16-2.30)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.24 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.189 , 0.251 0.189 , 0.251	Depositor DCC
R_{free} test set	3582 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å^2)	17.5	Xtriage
Anisotropy	0.154	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 36.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	12994	wwPDB-VP
Average B, all atoms (Å^2)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.58% 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: GAL, AAL, GOL, CA

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	0.67	1/6094 (0.0%)	0.76	3/8293 (0.0%)
1	B	0.68	4/6085 (0.1%)	0.74	4/8277 (0.0%)
All	All	0.67	5/12179 (0.0%)	0.75	7/16570 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	165	GLU	CG-CD	-20.29	1.21	1.51
1	B	84	LYS	CE-NZ	-15.21	1.11	1.49
1	B	259	LYS	CG-CD	-15.01	1.01	1.52
1	B	414	LYS	CD-CE	-12.48	1.20	1.51
1	B	588	LYS	CG-CD	7.52	1.78	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	259	LYS	CB-CG-CD	11.21	140.74	111.60
1	B	588	LYS	CB-CG-CD	-7.41	92.32	111.60
1	B	581	LEU	CA-CB-CG	6.19	129.53	115.30
1	A	539	ARG	NE-CZ-NH2	-5.72	117.44	120.30
1	A	193	ARG	NE-CZ-NH2	-5.35	117.62	120.30
1	A	165	GLU	CB-CG-CD	5.24	128.34	114.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	414	LYS	CD-CE-NZ	5.10	123.42	111.70

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	285	PRO	Peptide

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	5920	0	5586	60	0
1	B	5912	0	5576	67	0
2	C	21	0	17	5	0
2	D	21	0	17	8	0
2	E	22	0	19	3	0
2	G	21	0	17	9	0
2	H	22	0	19	4	0
3	F	42	0	33	7	0
4	A	10	0	8	7	0
5	A	11	0	10	8	0
6	A	6	0	8	0	0
6	B	18	0	24	1	0
7	A	3	0	0	0	0
7	B	3	0	0	0	0
8	A	512	0	0	11	0
8	B	450	0	0	5	0
All	All	12994	0	11334	151	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 6.

All (151) 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)
3:F:1:GAL:C1	2:G:2:AAL:O4	1.69	1.40
4:A:1795:AAL:C1	5:A:1796:GAL:O3	1.72	1.38
5:A:1796:GAL:C1	2:D:2:AAL:O4	1.75	1.34
4:A:1795:AAL:O4	2:C:1:GAL:C1	1.90	1.20
3:F:1:GAL:C1	2:G:2:AAL:HO4	1.47	1.18
2:D:1:GAL:C1	2:E:2:AAL:O4	1.92	1.16
5:A:1796:GAL:C1	2:D:2:AAL:HO4	1.67	0.98
1:B:175:LEU:HD21	8:B:2095:HOH:O	1.67	0.93
1:A:700:ALA:H	1:A:760:ASN:HD22	1.13	0.93
4:A:1795:AAL:HO4	2:C:1:GAL:C1	1.79	0.92
1:A:109:GLU:HG2	8:A:2049:HOH:O	1.71	0.91
3:F:1:GAL:C1	2:G:2:AAL:C4	2.50	0.89
4:A:1795:AAL:C1	5:A:1796:GAL:HO3	1.78	0.88
5:A:1796:GAL:C1	2:D:2:AAL:C4	2.54	0.84
1:A:649:ASP:HB3	8:A:2452:HOH:O	1.81	0.81
4:A:1795:AAL:C1	5:A:1796:GAL:C3	2.60	0.80
1:A:700:ALA:H	1:A:760:ASN:ND2	1.82	0.76
1:B:764:VAL:HG12	1:B:770:PRO:HA	1.67	0.76
5:A:1796:GAL:C2	2:D:2:AAL:O4	2.34	0.75
1:B:742:PHE:CZ	3:F:3:GAL:H4	2.23	0.74
1:B:700:ALA:H	1:B:760:ASN:HD22	1.37	0.73
1:A:716:GLN:HE22	1:A:719:ARG:HH11	1.36	0.72
1:B:716:GLN:HE22	1:B:719:ARG:HH11	1.35	0.71
2:D:1:GAL:C1	2:E:2:AAL:C4	2.70	0.70
2:G:1:GAL:H2	2:H:2:AAL:O4	1.92	0.69
1:B:162:HIS:HA	1:B:238:GLN:OE1	1.93	0.67
1:A:151:LYS:O	8:A:2095:HOH:O	2.13	0.66
1:A:87:MET:SD	1:A:216:VAL:HG23	2.35	0.65
4:A:1795:AAL:O4	2:C:1:GAL:C2	2.43	0.65
1:B:492:ASP:HB3	1:B:495:GLY:O	1.96	0.65
2:G:1:GAL:C2	2:H:2:AAL:O4	2.47	0.63
1:B:703:HIS:O	6:B:1802:GOL:H12	1.99	0.63
1:A:492:ASP:HB3	1:A:495:GLY:O	1.99	0.61
1:B:48:LEU:HD11	1:B:232:ILE:CD1	2.31	0.61
1:B:141:PHE:HB3	1:B:201:VAL:O	2.01	0.60
1:A:422:TYR:CE1	1:A:745:MET:HG2	2.38	0.59
1:A:141:PHE:HB3	1:A:201:VAL:O	2.02	0.59
1:B:176:ASN:HD22	1:B:182:ARG:HA	1.65	0.58
1:A:697:HIS:CE1	1:A:761:VAL:HG11	2.38	0.58
1:B:176:ASN:ND2	1:B:182:ARG:HE	2.03	0.57
1:A:708:PRO:HG2	1:A:712:HIS:CD2	2.41	0.56
1:A:194:GLN:HB3	8:A:2140:HOH:O	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:294:LEU:HD23	1:B:322:TYR:CE1	2.41	0.56
1:B:764:VAL:HG12	1:B:770:PRO:CA	2.36	0.55
1:A:87:MET:SD	1:A:216:VAL:CG2	2.95	0.54
1:B:85:LEU:HD22	1:B:97:LEU:HD23	1.89	0.54
1:A:422:TYR:CZ	1:A:745:MET:HG2	2.42	0.54
3:F:1:GAL:C1	2:G:2:AAL:H4	2.35	0.54
1:A:534:GLN:HA	1:A:648:PRO:HD3	1.90	0.53
1:A:64:ASN:HB2	1:A:95:THR:HG22	1.90	0.53
1:A:645:ALA:HB2	8:A:2370:HOH:O	2.09	0.53
1:A:386:ALA:HB3	8:A:2174:HOH:O	2.08	0.52
1:B:742:PHE:CE1	3:F:3:GAL:H4	2.44	0.52
1:B:654:MET:HG3	1:B:658:LYS:HD2	1.90	0.52
1:B:680:LYS:HG3	8:B:2120:HOH:O	2.10	0.52
1:A:675:GLY:O	1:A:677:PRO:HD3	2.09	0.51
1:B:661:ALA:O	1:B:689:LYS:HE2	2.11	0.51
1:B:247:VAL:HG22	1:B:749:LEU:HD23	1.92	0.51
1:A:766:VAL:HB	8:A:2493:HOH:O	2.11	0.51
1:B:270:ALA:O	1:B:274:VAL:HG23	2.11	0.51
1:B:697:HIS:O	1:B:698:ILE:HG13	2.10	0.51
5:A:1796:GAL:C1	2:D:2:AAL:H4	2.40	0.50
1:B:61:HIS:O	1:B:97:LEU:HA	2.11	0.50
1:B:500:VAL:HG12	1:B:505:PHE:CE2	2.46	0.50
1:A:306:ARG:HA	1:A:641:LEU:HD21	1.93	0.50
1:B:528:GLY:HA3	1:B:643:LEU:HD11	1.93	0.50
1:A:700:ALA:N	1:A:760:ASN:HD22	1.96	0.50
1:B:48:LEU:HD11	1:B:232:ILE:HD12	1.93	0.50
1:B:409:HIS:HE1	1:B:493:PHE:O	1.95	0.50
1:B:488:SER:HB2	8:B:2301:HOH:O	2.12	0.50
1:B:727:MET:O	1:B:731:ILE:HG13	2.11	0.49
1:A:60:ILE:HG12	1:A:99:ILE:HG23	1.94	0.49
1:A:168:ASP:HB2	1:A:174:ASP:OD2	2.13	0.49
1:B:44:GLY:HA2	1:B:46:HIS:N	2.27	0.49
1:B:121:SER:O	1:B:149:VAL:HA	2.12	0.49
1:A:443:ARG:HG2	1:A:466:TYR:CZ	2.47	0.49
1:B:119:ILE:HG22	1:B:147:ILE:HG21	1.95	0.49
1:A:590:SER:OG	1:A:594:GLU:OE1	2.28	0.49
1:B:623:ASP:O	1:B:627:LYS:HB2	2.12	0.49
1:A:716:GLN:NE2	1:A:719:ARG:HH11	2.09	0.49
1:B:374:PHE:CD2	1:B:394:TYR:HB2	2.48	0.48
1:B:647:PHE:HB3	1:B:652:MET:HB3	1.95	0.48
1:B:449:ARG:HD3	1:B:453:TRP:CZ2	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:VAL:HG22	1:A:211:LYS:HG3	1.96	0.48
1:B:285:PRO:O	1:B:286:SER:HB2	2.13	0.48
1:A:460:ASN:OD1	1:A:477:ASN:HB3	2.14	0.47
1:A:675:GLY:HA2	1:A:726:TYR:CD2	2.49	0.47
1:A:764:VAL:HG12	1:A:770:PRO:HA	1.96	0.47
1:A:379:LEU:HD21	1:A:382:PRO:HG3	1.96	0.47
1:B:129:PHE:O	1:B:144:SER:HA	2.15	0.47
1:A:558:SER:HA	1:A:564:ARG:HG3	1.97	0.47
1:B:115:LEU:HD21	1:B:212:ILE:HD13	1.97	0.46
1:B:669:TYR:O	1:B:693:ILE:HA	2.15	0.46
1:B:44:GLY:HA2	1:B:45:SER:C	2.36	0.46
1:B:396:HIS:CG	1:B:397:PRO:HD2	2.50	0.46
1:B:697:HIS:HB3	1:B:742:PHE:O	2.16	0.46
1:B:118:ASP:HB3	1:B:152:MET:SD	2.56	0.46
1:A:613:ASP:O	1:A:617:LEU:HG	2.15	0.46
1:B:709:GLY:HA2	1:B:759:TYR:CG	2.51	0.45
1:A:293:TRP:CD2	1:A:786:TYR:HB3	2.51	0.45
1:B:218:SER:HB2	1:B:673:LYS:HD2	1.99	0.45
1:B:515:VAL:O	1:B:519:GLU:HG3	2.17	0.45
1:B:781:VAL:HG22	8:B:2425:HOH:O	2.16	0.45
1:A:360:PRO:HD2	1:A:479:TRP:CD1	2.51	0.45
1:A:160:SER:HB3	1:A:206:LEU:HD11	1.99	0.45
1:A:270:ALA:O	1:A:274:VAL:HG23	2.17	0.45
1:B:697:HIS:CE1	1:B:761:VAL:HG11	2.52	0.45
2:D:1:GAL:C2	2:E:2:AAL:O4	2.61	0.45
2:G:1:GAL:C1	2:H:2:AAL:O4	2.65	0.44
1:A:374:PHE:O	1:A:377:ARG:HB2	2.16	0.44
1:A:751:GLY:HA2	1:A:757:GLU:O	2.17	0.44
4:A:1795:AAL:O4	2:C:1:GAL:H2	2.14	0.44
1:B:174:ASP:OD1	1:B:410:SER:HB2	2.17	0.44
1:A:695:GLU:HG2	1:A:742:PHE:HB2	2.00	0.44
3:F:1:GAL:C2	2:G:2:AAL:O4	2.54	0.44
1:A:657:VAL:CG1	1:A:687:LEU:HD11	2.48	0.44
1:B:539:ARG:HD3	8:B:2335:HOH:O	2.16	0.44
1:A:412:PRO:HD2	8:A:2050:HOH:O	2.17	0.44
1:B:342:GLY:HA2	1:B:413:LEU:HD11	2.00	0.43
1:A:699:GLY:HA3	1:A:708:PRO:O	2.18	0.43
1:B:755:ASP:OD1	1:B:757:GLU:HG3	2.18	0.43
1:A:133:VAL:O	1:A:140:VAL:HA	2.18	0.43
1:A:543:ASP:CG	1:A:611:ARG:HH12	2.22	0.43
2:G:1:GAL:C1	2:H:2:AAL:C4	2.96	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:657:VAL:HG12	1:A:687:LEU:HD11	2.00	0.43
1:A:175:LEU:O	1:A:197:TRP:NE1	2.52	0.43
1:A:294:LEU:HD23	1:A:294:LEU:HA	1.76	0.42
1:B:708:PRO:HG2	1:B:712:HIS:CD2	2.54	0.42
1:B:542:SER:H	1:B:545:ALA:HB3	1.85	0.42
1:B:218:SER:CB	1:B:673:LYS:HD2	2.49	0.42
1:B:67:ALA:HB1	1:B:83:LEU:HD11	2.01	0.42
1:B:318:ASP:OD1	1:B:318:ASP:C	2.58	0.42
1:A:547:TYR:CD1	1:A:614:TYR:HB3	2.54	0.42
1:A:647:PHE:HB3	1:A:652:MET:HB3	2.01	0.42
1:B:667:VAL:O	1:B:691:SER:HA	2.19	0.42
1:A:609:THR:HG23	8:A:2398:HOH:O	2.19	0.42
1:B:484:PHE:CE1	1:B:498:PRO:HB3	2.55	0.42
1:A:695:GLU:OE1	2:C:1:GAL:C1	2.67	0.42
1:B:49:PHE:HB2	1:B:230:ILE:HD12	2.01	0.42
1:A:81:LYS:NZ	8:A:2004:HOH:O	2.52	0.41
1:A:477:ASN:HA	1:A:530:PHE:O	2.20	0.41
1:B:123:GLY:O	1:B:149:VAL:HG11	2.20	0.41
1:B:297:PRO:HD2	1:B:323:PRO:HG2	2.03	0.41
1:B:97:LEU:HD12	1:B:97:LEU:O	2.21	0.40
1:B:143:ARG:HG2	1:B:197:TRP:HA	2.03	0.40
1:A:135:ASP:OD2	1:A:139:ALA:HB3	2.21	0.40
1:A:242:PHE:CE2	1:A:243:LEU:HG	2.57	0.40
1:B:242:PHE:CZ	1:B:243:LEU:HG	2.56	0.40
1:A:109:GLU:CG	8:A:2049:HOH:O	2.47	0.40
1:A:542:SER:H	1:A:545:ALA:HB3	1.86	0.40
1:B:460:ASN:OD1	1:B:477:ASN:HB3	2.22	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	749/750 (100%)	722 (96%)	26 (4%)	1 (0%)	51	64
1	B	747/750 (100%)	721 (96%)	26 (4%)	0	100	100
All	All	1496/1500 (100%)	1443 (96%)	52 (4%)	1 (0%)	51	64

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	576	LYS

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	622/630 (99%)	613 (99%)	9 (1%)	67	81
1	B	617/630 (98%)	608 (98%)	9 (2%)	65	79
All	All	1239/1260 (98%)	1221 (98%)	18 (2%)	65	79

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	99	ILE
1	A	109	GLU
1	A	119	ILE
1	A	129	PHE
1	A	140	VAL
1	A	213	SER
1	A	395	ASP
1	A	414	LYS
1	A	747	SER
1	B	63	LEU
1	B	175	LEU
1	B	190	SER
1	B	309	LYS
1	B	356	ASP
1	B	488	SER

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Mol	Chain	Res	Type
1	B	599	VAL
1	B	745	MET
1	B	792	SER

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

Mol	Chain	Res	Type
1	A	194	GLN
1	A	409	HIS
1	A	716	GLN
1	A	760	ASN
1	B	176	ASN
1	B	194	GLN
1	B	409	HIS
1	B	546	GLN
1	B	716	GLN
1	B	760	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)

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

5.5 Carbohydrates (i)

14 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	GAL	C	1	2	11,11,12	0.82	0	15,15,17	2.45	5 (33%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	AAL	C	2	2	11,11,12	1.74	3 (27%)	15,16,18	1.30	2 (13%)
2	GAL	D	1	2	11,11,12	1.03	0	15,15,17	2.40	4 (26%)
2	AAL	D	2	2	11,11,12	0.47	0	15,16,18	0.65	0
2	GAL	E	1	2	12,12,12	0.65	0	17,17,17	1.17	2 (11%)
2	AAL	E	2	2	11,11,12	1.71	1 (9%)	15,16,18	2.12	3 (20%)
3	GAL	F	1	3	11,11,12	1.05	1 (9%)	15,15,17	2.17	4 (26%)
3	AAL	F	2	3	11,11,12	1.16	1 (9%)	15,16,18	1.82	2 (13%)
3	GAL	F	3	3	11,11,12	1.87	3 (27%)	15,15,17	3.26	7 (46%)
3	AAL	F	4	3	11,11,12	1.43	2 (18%)	15,16,18	1.87	4 (26%)
2	GAL	G	1	2	11,11,12	0.90	0	15,15,17	3.01	3 (20%)
2	AAL	G	2	2	11,11,12	1.26	2 (18%)	15,16,18	1.21	2 (13%)
2	GAL	H	1	2	12,12,12	0.58	0	17,17,17	1.79	4 (23%)
2	AAL	H	2	2	11,11,12	1.63	2 (18%)	15,16,18	1.90	4 (26%)

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	GAL	C	1	2	-	2/2/19/22	0/1/1/1
2	AAL	C	2	2	-	-	0/3/2/2
2	GAL	D	1	2	-	2/2/19/22	0/1/1/1
2	AAL	D	2	2	-	-	0/3/2/2
2	GAL	E	1	2	-	0/2/22/22	0/1/1/1
2	AAL	E	2	2	-	-	0/3/2/2
3	GAL	F	1	3	-	0/2/19/22	0/1/1/1
3	AAL	F	2	3	-	-	0/3/2/2
3	GAL	F	3	3	-	2/2/19/22	0/1/1/1
3	AAL	F	4	3	-	-	0/3/2/2
2	GAL	G	1	2	-	0/2/19/22	0/1/1/1
2	AAL	G	2	2	-	-	0/3/2/2
2	GAL	H	1	2	-	0/2/22/22	0/1/1/1
2	AAL	H	2	2	-	-	0/3/2/2

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	2	AAL	O5-C1	4.91	1.51	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	2	AAL	O5-C1	3.80	1.49	1.43
2	H	2	AAL	O5-C1	3.63	1.49	1.43
3	F	3	GAL	O5-C1	3.61	1.49	1.43
3	F	3	GAL	C2-C3	3.37	1.57	1.52
3	F	4	AAL	O5-C5	2.71	1.48	1.43
3	F	2	AAL	O5-C5	2.67	1.48	1.43
3	F	4	AAL	O5-C1	2.58	1.47	1.43
2	G	2	AAL	O5-C1	2.40	1.47	1.43
2	G	2	AAL	O5-C5	2.32	1.48	1.43
2	C	2	AAL	O3-C3	2.31	1.49	1.43
3	F	1	GAL	O5-C1	2.29	1.47	1.43
3	F	3	GAL	O5-C5	2.20	1.47	1.43
2	C	2	AAL	O5-C5	2.14	1.47	1.43
2	H	2	AAL	O3-C6	2.01	1.48	1.43

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	G	1	GAL	O5-C1-C2	-10.46	94.63	110.77
2	D	1	GAL	C1-O5-C5	-7.00	102.70	112.19
3	F	3	GAL	O5-C1-C2	6.74	121.18	110.77
3	F	3	GAL	C1-O5-C5	5.81	120.06	112.19
2	C	1	GAL	O5-C1-C2	5.69	119.55	110.77
3	F	3	GAL	O3-C3-C4	-5.37	97.93	110.35
2	E	2	AAL	C2-C3-C4	-5.34	100.63	113.85
3	F	1	GAL	O3-C3-C2	-5.28	99.89	109.99
2	C	1	GAL	C1-O5-C5	5.11	119.12	112.19
2	E	2	AAL	O5-C1-C2	4.89	118.32	110.77
3	F	1	GAL	C1-C2-C3	-4.80	103.77	109.67
3	F	3	GAL	C2-C3-C4	4.79	119.18	110.89
3	F	2	AAL	C1-C2-C3	-4.52	103.71	109.17
2	H	2	AAL	O3-C3-C2	4.10	117.44	108.07
3	F	4	AAL	O5-C5-C6	3.98	119.23	113.33
2	H	2	AAL	C1-O5-C5	3.84	117.40	112.19
3	F	4	AAL	O5-C1-C2	3.82	116.67	110.77
2	H	1	GAL	C1-C2-C3	3.82	118.23	110.31
2	H	1	GAL	O5-C1-C2	3.53	116.58	110.28
3	F	4	AAL	O4-C4-C5	-3.29	101.53	111.05
2	D	1	GAL	O3-C3-C2	-3.25	103.77	109.99
2	G	1	GAL	C1-O5-C5	-3.25	107.79	112.19
3	F	2	AAL	O5-C5-C6	-3.19	108.60	113.33
2	H	2	AAL	O5-C1-C2	3.10	115.56	110.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	H	1	GAL	C4-C3-C2	3.05	116.15	110.82
2	C	1	GAL	O3-C3-C4	-3.05	103.30	110.35
2	E	2	AAL	O3-C3-C2	2.96	114.83	108.07
2	D	1	GAL	C6-C5-C4	-2.86	106.30	113.00
2	E	1	GAL	C1-O5-C5	-2.79	108.40	113.66
3	F	4	AAL	C6-C5-C4	-2.78	97.42	102.28
2	C	1	GAL	O2-C2-C1	-2.70	103.62	109.15
3	F	1	GAL	O5-C5-C6	2.59	111.27	107.20
3	F	3	GAL	C3-C4-C5	2.55	114.78	110.24
2	D	1	GAL	O5-C5-C6	2.52	111.16	107.20
2	G	1	GAL	O3-C3-C2	-2.39	105.42	109.99
2	G	2	AAL	O5-C5-C6	2.35	116.81	113.33
2	E	1	GAL	O5-C5-C6	2.34	112.27	106.44
2	H	2	AAL	C2-C3-C4	-2.28	108.22	113.85
2	C	2	AAL	C1-O5-C5	2.24	115.23	112.19
3	F	3	GAL	O5-C5-C4	2.24	116.27	110.83
2	G	2	AAL	C6-C5-C4	-2.23	98.38	102.28
2	H	1	GAL	C1-O5-C5	2.23	117.87	113.66
3	F	3	GAL	O2-C2-C1	-2.21	104.64	109.15
2	C	2	AAL	C2-C3-C4	-2.19	108.43	113.85
2	C	1	GAL	C1-C2-C3	2.12	112.28	109.67
3	F	1	GAL	C2-C3-C4	-2.04	107.36	110.89

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	F	3	GAL	O5-C5-C6-O6
2	C	1	GAL	O5-C5-C6-O6
2	D	1	GAL	O5-C5-C6-O6
2	C	1	GAL	C4-C5-C6-O6
2	D	1	GAL	C4-C5-C6-O6
3	F	3	GAL	C4-C5-C6-O6

There are no ring outliers.

9 monomers are involved in 24 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1	GAL	3	0
2	H	2	AAL	4	0
3	F	1	GAL	5	0

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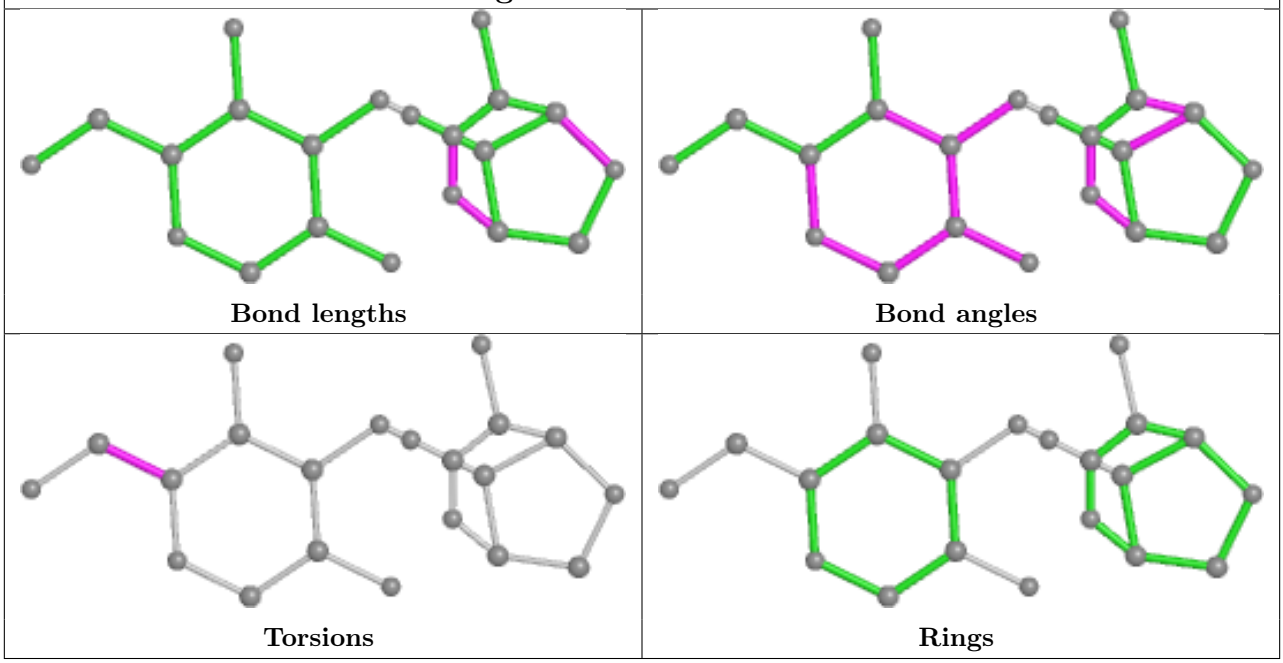
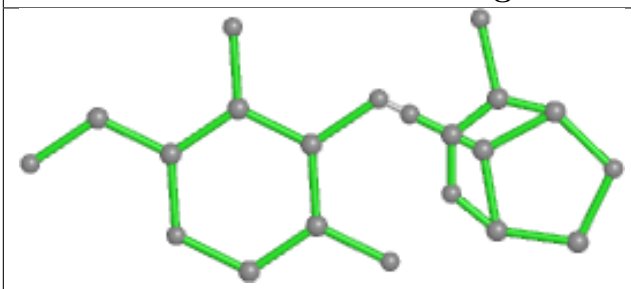
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	2	AAL	5	0
3	F	3	GAL	2	0
2	D	2	AAL	5	0
2	E	2	AAL	3	0
2	G	1	GAL	4	0
2	C	1	GAL	5	0

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

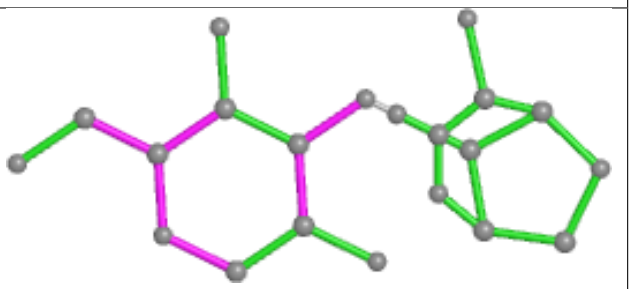
Oligosaccharide Chain C



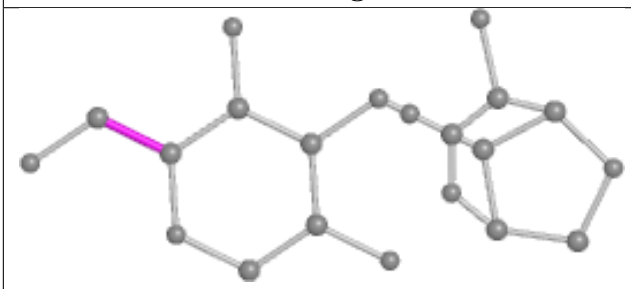
Oligosaccharide Chain D



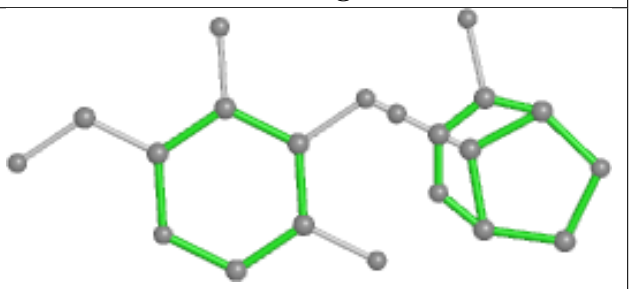
Bond lengths



Bond angles

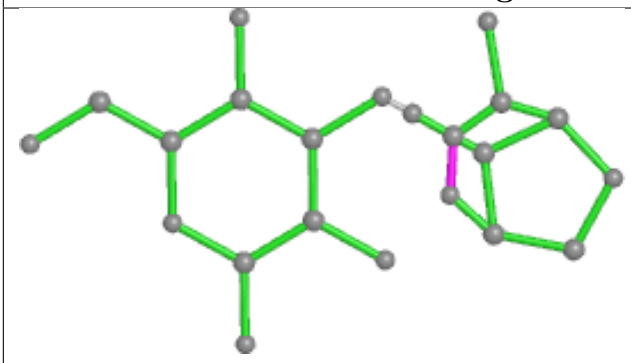


Torsions

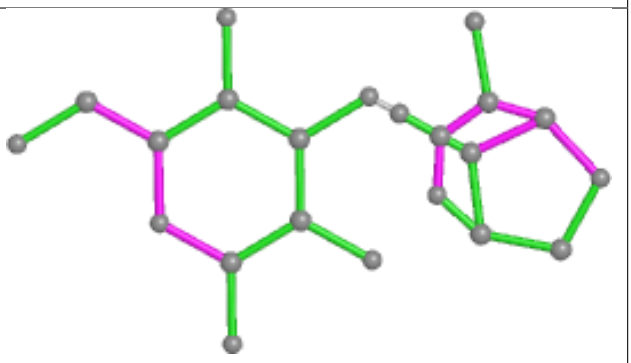


Rings

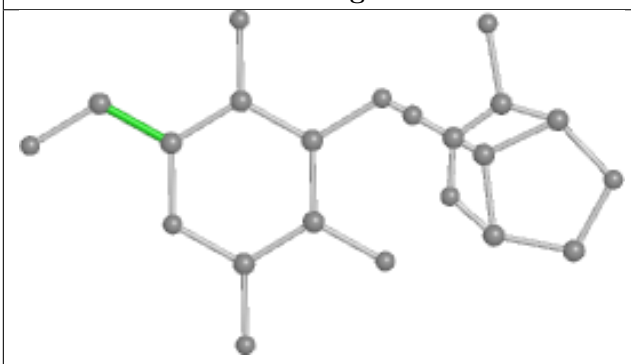
Oligosaccharide Chain E



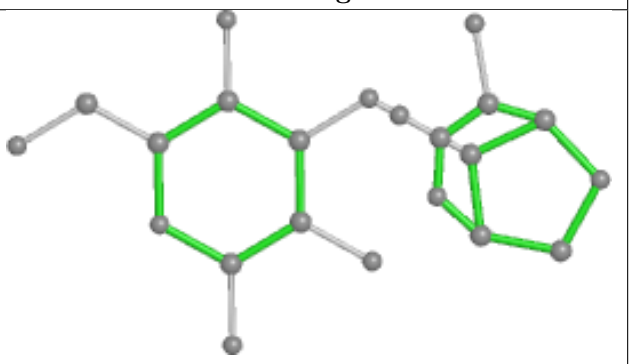
Bond lengths



Bond angles

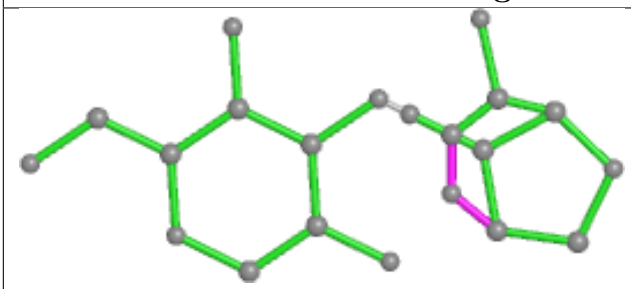


Torsions

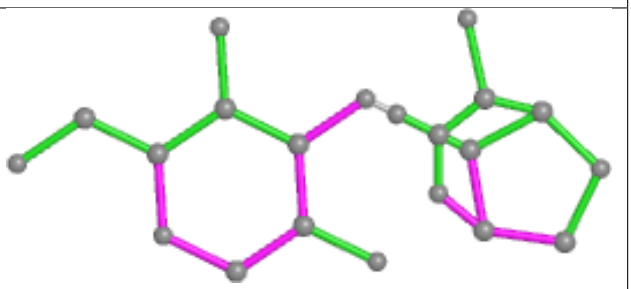


Rings

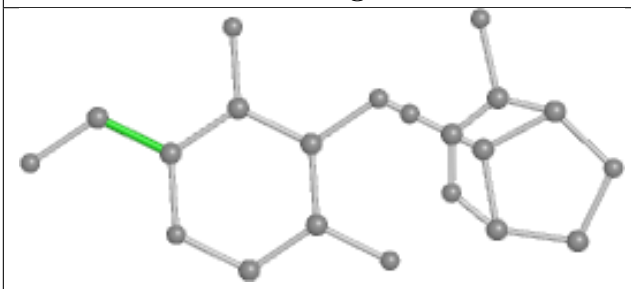
Oligosaccharide Chain G



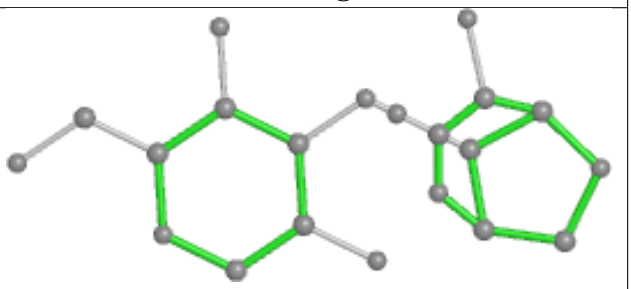
Bond lengths



Bond angles

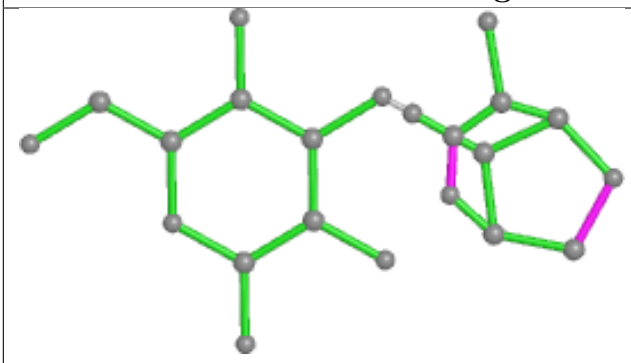


Torsions

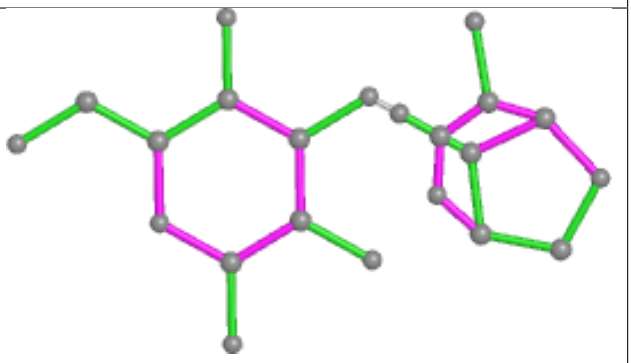


Rings

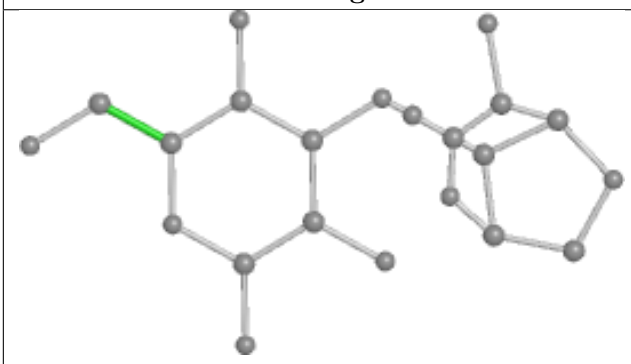
Oligosaccharide Chain H



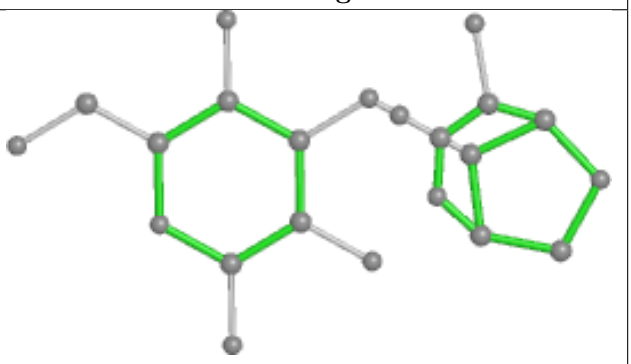
Bond lengths



Bond angles

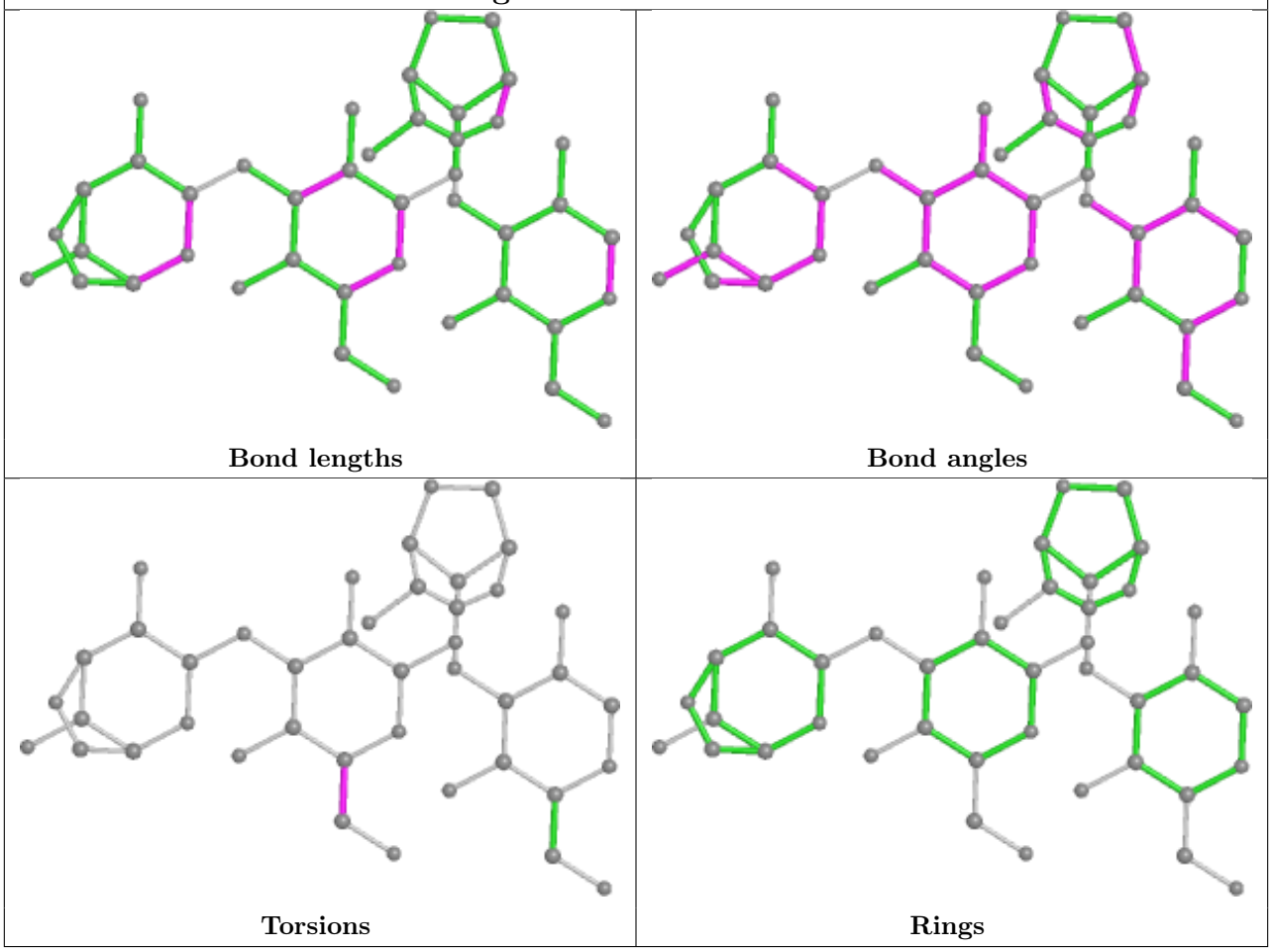


Torsions



Rings

Oligosaccharide Chain F



5.6 Ligand geometry (i)

Of 12 ligands modelled in this entry, 6 are monoatomic - leaving 6 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	$\#\|Z\| > 2$
6	GOL	B	1802	-	5,5,5	0.47	0	5,5,5	0.37	0
5	GAL	A	1796	-	11,11,12	0.99	1 (9%)	15,15,17	2.80	4 (26%)
6	GOL	A	1801	-	5,5,5	0.56	0	5,5,5	0.53	0
6	GOL	B	1804	-	5,5,5	0.31	0	5,5,5	0.38	0
4	AAL	A	1795	-	11,11,12	2.24	3 (27%)	15,16,18	1.57	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
6	GOL	B	1803	-	5,5,5	0.31	0	5,5,5	0.19	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
6	GOL	B	1802	-	-	2/4/4/4	-
5	GAL	A	1796	-	-	2/2/19/22	0/1/1/1
6	GOL	A	1801	-	-	4/4/4/4	-
6	GOL	B	1804	-	-	1/4/4/4	-
4	AAL	A	1795	-	-	-	0/3/2/2
6	GOL	B	1803	-	-	0/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1795	AAL	O5-C1	4.61	1.51	1.43
4	A	1795	AAL	O5-C5	3.82	1.51	1.43
4	A	1795	AAL	O3-C6	3.54	1.51	1.43
5	A	1796	GAL	O3-C3	-2.04	1.38	1.43

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	A	1796	GAL	O3-C3-C2	-7.45	95.72	109.99
5	A	1796	GAL	O5-C1-C2	-5.51	102.27	110.77
4	A	1795	AAL	C1-C2-C3	-3.76	104.64	109.17
5	A	1796	GAL	C1-O5-C5	-3.32	107.70	112.19
4	A	1795	AAL	O5-C1-C2	3.22	115.74	110.77
5	A	1796	GAL	C1-C2-C3	-2.63	106.43	109.67

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1796	GAL	C4-C5-C6-O6
6	A	1801	GOL	O1-C1-C2-C3
6	A	1801	GOL	C1-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
6	B	1802	GOL	O1-C1-C2-C3
6	B	1804	GOL	C1-C2-C3-O3
5	A	1796	GAL	O5-C5-C6-O6
6	A	1801	GOL	O2-C2-C3-O3
6	A	1801	GOL	O1-C1-C2-O2
6	B	1802	GOL	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	1802	GOL	1	0
5	A	1796	GAL	8	0
4	A	1795	AAL	7	0

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	748/750 (99%)	-0.11	6 (0%)	86	89	6, 15, 33, 60	1 (0%)
1	B	749/750 (99%)	-0.05	4 (0%)	91	94	7, 18, 35, 69	6 (0%)
All	All	1497/1500 (99%)	-0.08	10 (0%)	87	91	6, 16, 35, 69	7 (0%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	171	ASP	3.5
1	B	394	TYR	2.8
1	A	394	TYR	2.8
1	B	166	VAL	2.7
1	B	172	VAL	2.7
1	A	348	ALA	2.3
1	A	171	ASP	2.2
1	A	168	ASP	2.2
1	A	170	GLY	2.2
1	A	351	ALA	2.1

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

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

6.3 Carbohydrates (i)

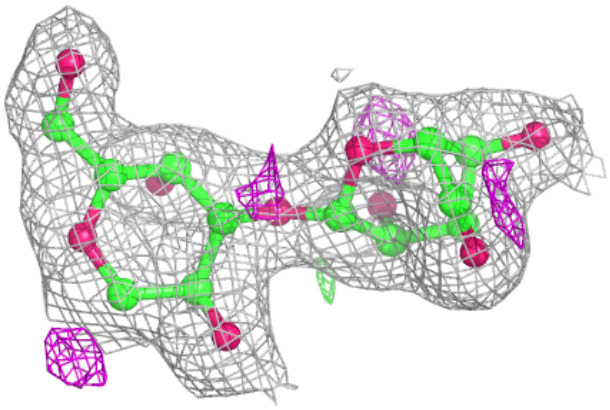
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

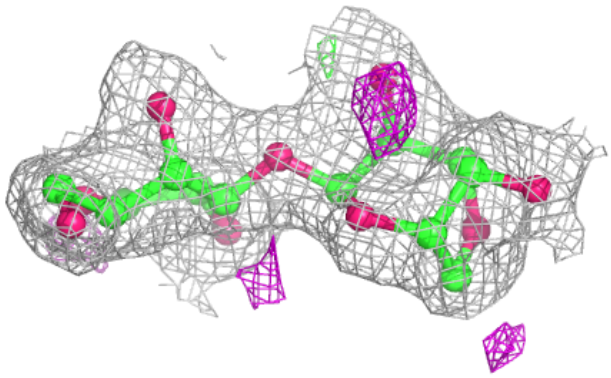
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	AAL	H	2	10/11	0.80	0.15	40,45,46,47	0
2	GAL	E	1	12/12	0.81	0.23	44,52,56,57	0
2	GAL	H	1	12/12	0.84	0.20	53,63,70,77	0
2	GAL	G	1	11/12	0.89	0.14	24,27,28,29	0
2	AAL	G	2	10/11	0.92	0.14	21,24,25,26	0
2	AAL	E	2	10/11	0.94	0.11	26,28,30,32	0
2	AAL	C	2	10/11	0.95	0.15	8,8,8,9	0
2	GAL	D	1	11/12	0.95	0.12	14,15,16,17	0
2	AAL	D	2	10/11	0.95	0.14	14,15,16,16	0
3	GAL	F	1	11/12	0.95	0.14	13,14,15,17	0
3	GAL	F	3	11/12	0.95	0.16	7,8,8,9	0
3	AAL	F	2	10/11	0.96	0.16	8,8,9,9	0
3	AAL	F	4	10/11	0.96	0.15	7,7,8,8	0
2	GAL	C	1	11/12	0.97	0.16	7,8,8,8	0

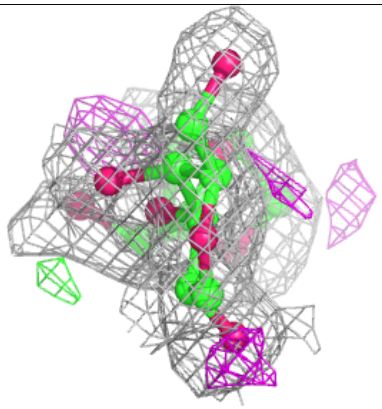
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain C:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative) and green (positive)

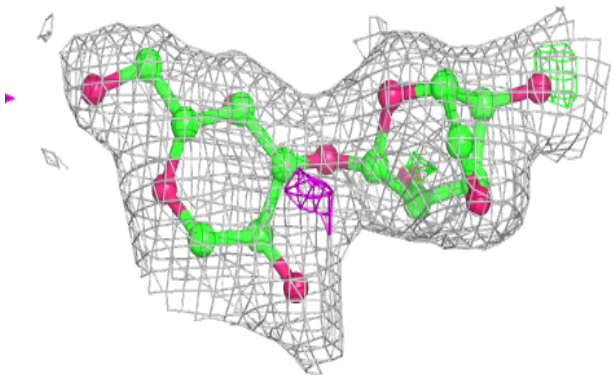


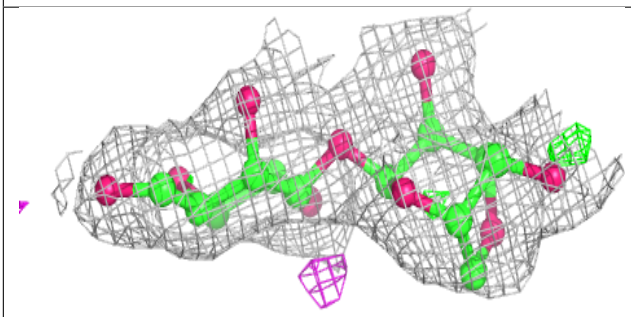


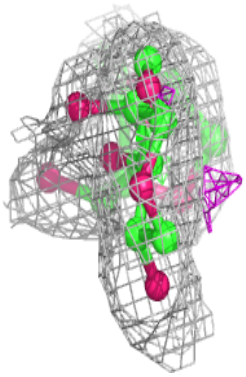


Electron density around Chain D:

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

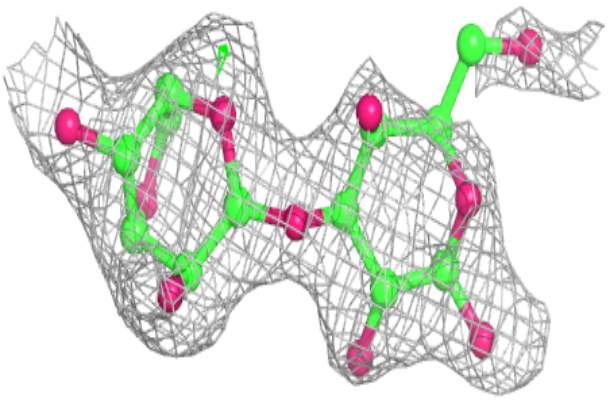


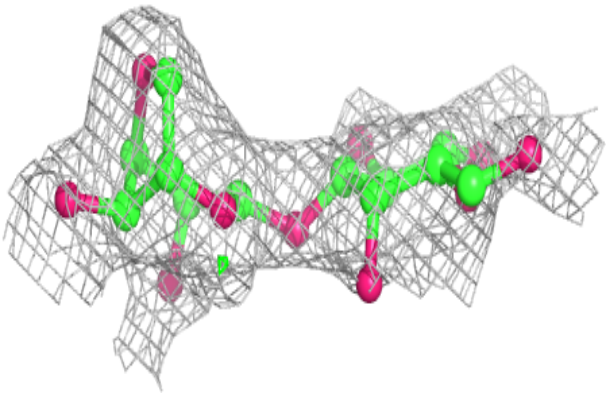


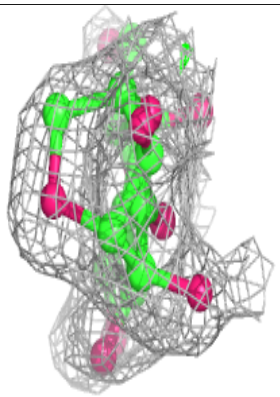


Electron density around Chain E:

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

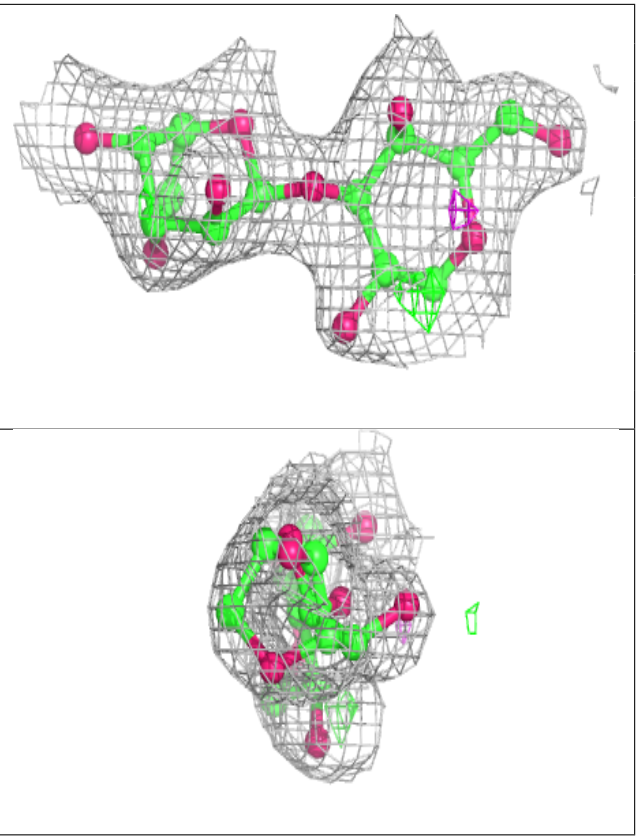






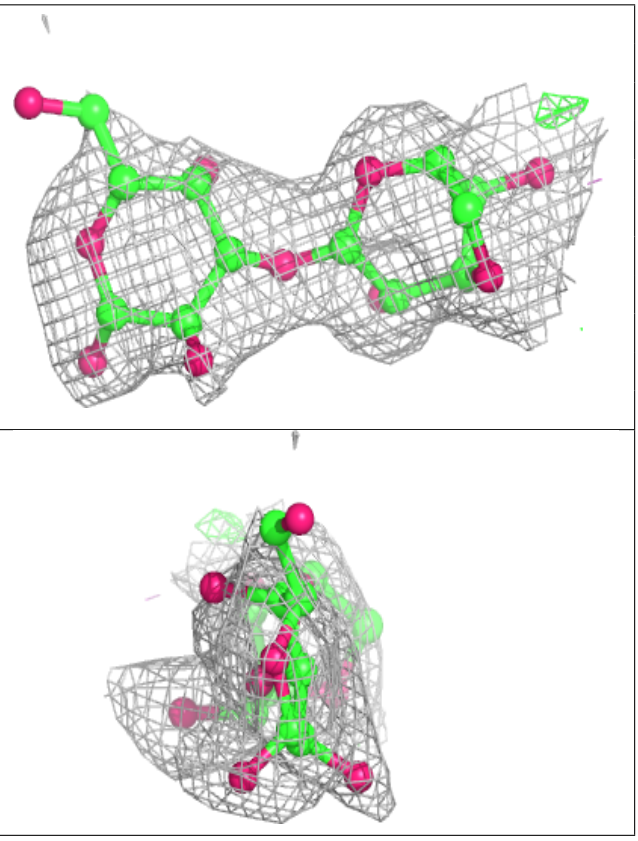
Electron density around Chain G:

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



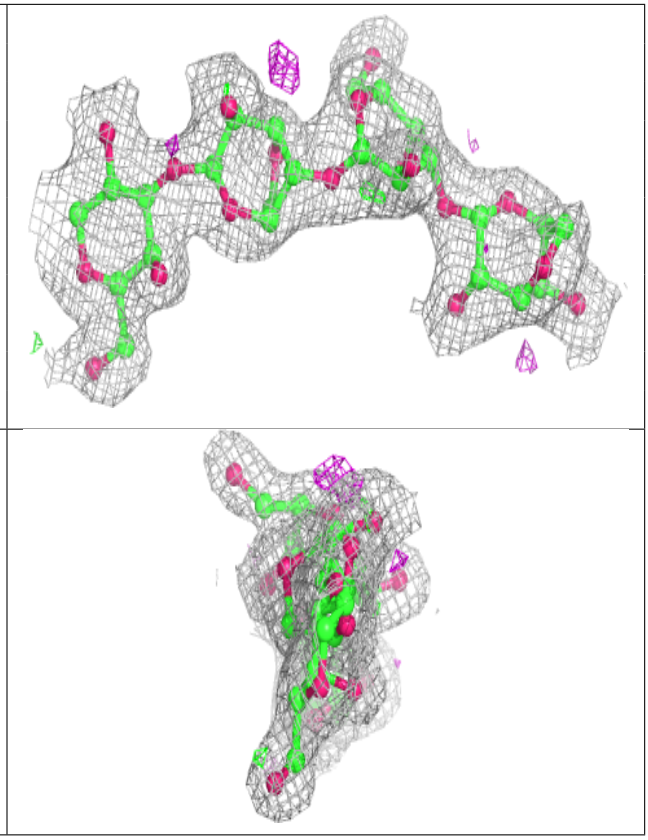
Electron density around Chain H:

$2mF_o$-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 F:

$2mF_o$-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
6	GOL	B	1804	6/6	0.84	0.26	52,57,58,58	0
6	GOL	A	1801	6/6	0.90	0.16	22,23,26,27	0
6	GOL	B	1803	6/6	0.93	0.16	32,36,39,47	0
6	GOL	B	1802	6/6	0.93	0.13	16,19,19,19	0
4	AAL	A	1795	10/11	0.96	0.17	7,7,7,7	0
7	CA	A	1804	1/1	0.97	0.07	35,35,35,35	0
5	GAL	A	1796	11/12	0.98	0.14	11,11,12,12	0
7	CA	B	1806	1/1	0.98	0.02	26,26,26,26	0
7	CA	B	1805	1/1	0.99	0.02	19,19,19,19	0
7	CA	A	1803	1/1	0.99	0.07	16,16,16,16	0
7	CA	B	1807	1/1	0.99	0.03	21,21,21,21	0
7	CA	A	1802	1/1	1.00	0.07	17,17,17,17	0

6.5 Other polymers (i)

There are no such residues in this entry.