Full wwPDB X-ray Structure Validation Report (i)

Oct 15, 2023 – 06:54 PM EDT

PDB ID	:	2FV0
Title	:	UGL_D88N/dGlcA-Glc-Rha-Glc
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Deposited on	:	2006-01-28
Resolution	:	1.91 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 (1)) were used in the production of this report:

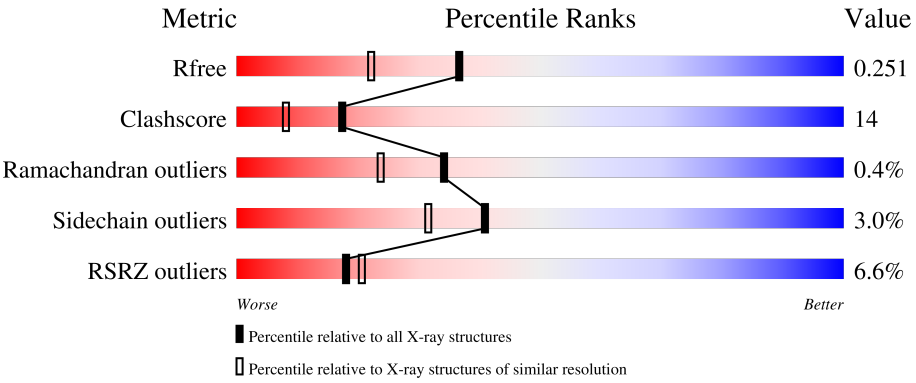
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, 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 1.91 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	7937 (1.94-1.90)
Clashscore	141614	8644 (1.94-1.90)
Ramachandran outliers	138981	8530 (1.94-1.90)
Sidechain outliers	138945	8530 (1.94-1.90)
RSRZ outliers	127900	7793 (1.94-1.90)

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	377	12% 69% 30%
1	B	377	2% 80% 18%
2	C	4	25% 75%

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
2	RAM	C	2	-	-	-	X
2	BGC	C	3	-	-	-	X

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 6928 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.

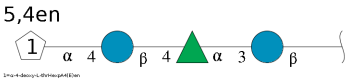
- Molecule 1 is a protein called Unsaturated glucuronyl hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	377	Total	C	N	O	S	0	10	0
			3091	1939	560	586	6			
1	B	377	Total	C	N	O	S	0	12	0
			3098	1949	563	579	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	88	ASN	ASP	engineered mutation	UNP Q9RC92
B	88	ASN	ASP	engineered mutation	UNP Q9RC92

- Molecule 2 is an oligosaccharide called 2,6-anhydro-3-deoxy-L-threo-hex-2-enonic acid-(1-4)-beta-D-glucopyranose-(1-4)-alpha-L-rhamnopyranose-(1-3)-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	4	Total	C	O	0	0	0
			44	24	20			

- Molecule 3 is water.

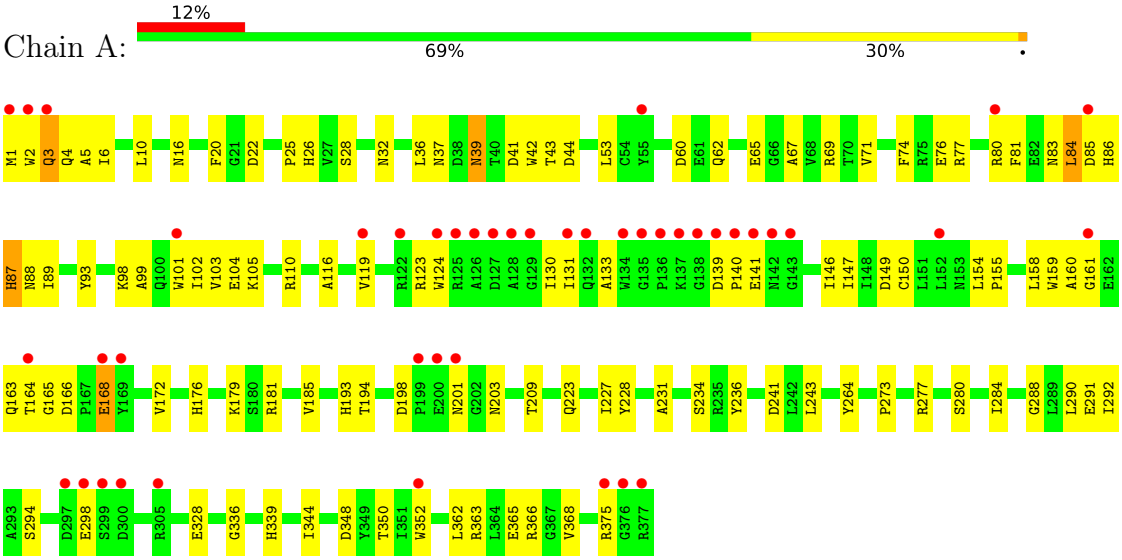
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	319	Total	O	0	0
			319	319		
3	B	376	Total	O	0	0
			376	376		

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.

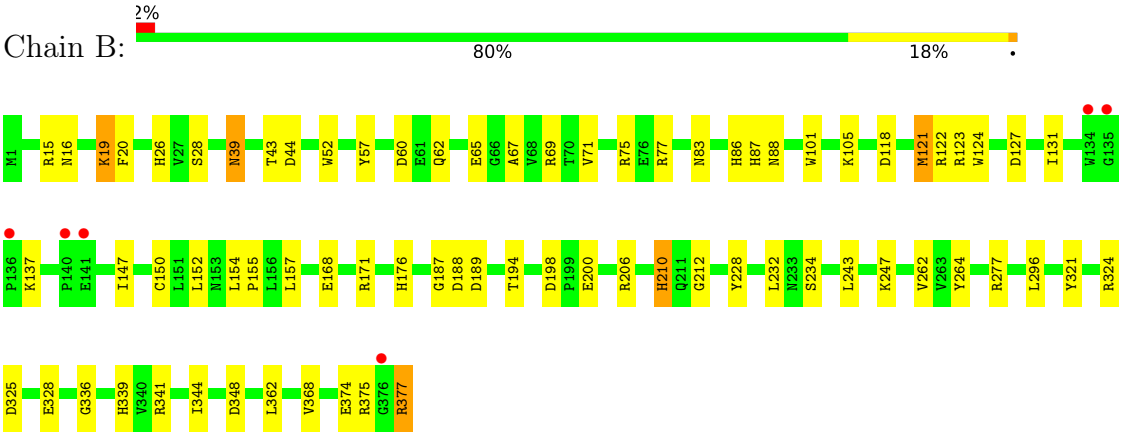
• Molecule 1: Unsaturated glucuronyl hydrolase

Chain A:



• Molecule 1: Unsaturated glucuronyl hydrolase

Chain B:



• Molecule 2: 2,6-anhydro-3-deoxy-L-threo-hex-2-enonic acid-(1-4)-beta-D-glucopyranose-(1-4)-alpha-L-rhamnopyranose-(1-3)-beta-D-glucopyranose

Chain C:


BGC1
RAM2
BGC3
GAD4

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	87.66Å 93.48Å 95.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	14.95 – 1.91 39.82 – 1.91	Depositor EDS
% Data completeness (in resolution range)	98.5 (14.95-1.91) 98.7 (39.82-1.91)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.90 (at 1.91Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.198 , 0.254 0.193 , 0.251	Depositor DCC
R_{free} test set	6109 reflections (10.10%)	wwPDB-VP
Wilson B-factor (Å^2)	18.6	Xtriage
Anisotropy	0.565	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 66.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.44$, $< L^2 > = 0.26$	Xtriage
Estimated twinning fraction	0.040 for -h,l,k	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	6928	wwPDB-VP
Average B, all atoms (Å^2)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.61% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BGC, RAM, GAD

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	0.31	0/3209	0.56	1/4343 (0.0%)
1	B	0.33	0/3227	0.59	1/4369 (0.0%)
All	All	0.32	0/6436	0.57	2/8712 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	336	GLY	N-CA-C	-5.39	99.62	113.10
1	A	336	GLY	N-CA-C	-5.25	99.97	113.10

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3091	0	2914	100	0
1	B	3098	0	2930	67	0
2	C	44	0	35	8	0
3	A	319	0	0	7	0
3	B	376	0	0	6	0
All	All	6928	0	5879	167	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 (167) 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:B:150[A]:CYS:SG	1:B:176:HIS:HE1	1.88	0.96
1:A:344:ILE:HG23	2:C:1:BGC:H1	1.46	0.95
1:A:103:VAL:HG12	1:A:104[B]:GLU:HG3	1.46	0.94
1:A:22:ASP:HB3	1:A:69[A]:ARG:HH22	1.36	0.89
1:A:344:ILE:CG2	2:C:1:BGC:H1	2.03	0.87
1:A:4:GLN:HG2	3:A:1159:HOH:O	1.77	0.84
1:B:131:ILE:H	1:B:176:HIS:HD2	1.26	0.83
1:A:131:ILE:H	1:A:176:HIS:HD2	1.26	0.82
1:A:60:ASP:OD1	1:A:62:GLN:HG2	1.79	0.81
1:B:171:ARG:HH11	1:B:171:ARG:HB3	1.45	0.81
2:C:3:BGC:H6C1	2:C:4:GAD:O5	1.80	0.81
1:A:65:GLU:HG3	1:A:69[B]:ARG:HH12	1.48	0.79
1:B:131:ILE:H	1:B:176:HIS:CD2	2.02	0.77
1:B:150[A]:CYS:SG	1:B:176:HIS:CE1	2.77	0.77
1:B:28:SER:OG	1:B:348:ASP:HB3	1.87	0.74
1:B:212:GLY:O	1:B:341:ARG:HD2	1.88	0.73
1:A:339:HIS:NE2	2:C:3:BGC:H6C2	2.05	0.72
1:B:101:TRP:O	1:B:105:LYS:HE2	1.90	0.72
1:A:363[B]:ARG:HH21	1:A:363[B]:ARG:HG2	1.55	0.71
1:A:65:GLU:HG3	1:A:69[B]:ARG:NH1	2.05	0.69
1:B:60:ASP:OD1	1:B:62:GLN:HG2	1.93	0.69
1:A:131:ILE:H	1:A:176:HIS:CD2	2.11	0.68
1:B:137:LYS:HE2	3:B:1299:HOH:O	1.93	0.68
1:A:124:TRP:CE3	1:A:172:VAL:HG13	2.29	0.67
1:B:39:ASN:ND2	1:B:77:ARG:HH11	1.91	0.67
1:B:86:HIS:CD2	1:B:88:ASN:HB2	2.30	0.66
1:A:39:ASN:HA	1:A:44:ASP:OD2	1.96	0.65
1:A:116:ALA:O	1:A:119:VAL:HG22	1.96	0.65
1:A:375:ARG:HH21	1:A:375:ARG:HG3	1.61	0.65
1:A:176:HIS:HE1	3:A:1084:HOH:O	1.80	0.64
1:A:193:HIS:HB3	1:A:209:THR:HB	1.81	0.62
1:B:171:ARG:HB3	1:B:171:ARG:NH1	2.13	0.62
1:A:159:TRP:O	1:A:163:GLN:HG2	2.00	0.62
1:B:86:HIS:HD2	1:B:88:ASN:H	1.48	0.61
1:B:123:ARG:HD3	3:B:1663:HOH:O	1.99	0.61
1:A:344:ILE:HG23	2:C:1:BGC:H5	1.83	0.61
1:A:28:SER:OG	1:A:348:ASP:HB3	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:HIS:ND1	1:A:150:CYS:SG	2.52	0.60
1:B:122[B]:ARG:HG3	1:B:122[B]:ARG:HH11	1.67	0.58
1:B:131:ILE:N	1:B:176:HIS:HD2	2.00	0.58
1:B:105:LYS:HE3	3:B:1681:HOH:O	2.03	0.57
1:B:198:ASP:OD1	1:B:200:GLU:HB2	2.04	0.57
1:B:121:MET:HG3	3:B:1453:HOH:O	2.05	0.56
1:A:130:ILE:CG2	1:A:179:LYS:HG2	2.36	0.56
1:A:39:ASN:ND2	1:A:77:ARG:HH11	2.03	0.56
1:A:339:HIS:HB2	1:A:344:ILE:HB	1.87	0.56
1:A:363[B]:ARG:HG2	1:A:363[B]:ARG:NH2	2.19	0.55
1:A:4:GLN:HG3	1:A:5:ALA:N	2.21	0.55
1:A:83:ASN:HD22	1:A:83:ASN:N	2.02	0.55
1:B:368:VAL:HG13	1:B:375:ARG:NH1	2.22	0.55
1:A:344:ILE:HG23	2:C:1:BGC:C1	2.31	0.54
1:A:164:THR:C	1:A:166:ASP:H	2.11	0.53
1:B:339:HIS:HB2	1:B:344:ILE:HB	1.90	0.52
1:A:87:HIS:HB3	1:A:133:ALA:HA	1.92	0.52
1:B:60:ASP:CG	1:B:62:GLN:HG2	2.29	0.52
1:B:154:LEU:N	1:B:155:PRO:CD	2.73	0.52
1:A:4:GLN:CG	1:A:5:ALA:N	2.72	0.51
1:A:84:LEU:HG	1:A:89:ILE:HD12	1.92	0.51
1:A:86:HIS:O	1:A:89:ILE:HG12	2.11	0.51
1:B:65:GLU:O	1:B:69:ARG:HG3	2.10	0.51
1:A:291:GLU:HA	1:A:363[A]:ARG:NH2	2.26	0.51
1:A:158:LEU:HD13	1:A:236:TYR:CB	2.41	0.51
1:A:160:ALA:O	1:A:164:THR:HG23	2.11	0.51
1:B:171:ARG:HH11	1:B:171:ARG:CB	2.18	0.51
1:A:130:ILE:HD12	1:A:130:ILE:O	2.09	0.50
1:B:121:MET:CE	1:B:168:GLU:HB3	2.41	0.50
1:B:324:ARG:NH2	1:B:324:ARG:HG2	2.26	0.50
1:A:42:TRP:CE3	2:C:4:GAD:H3	2.47	0.50
1:A:147:ILE:HG23	1:A:194:THR:HG22	1.92	0.50
1:A:166:ASP:OD1	1:A:168:GLU:HB2	2.12	0.50
1:B:86:HIS:HD2	1:B:88:ASN:HB2	1.76	0.50
1:A:203:ASN:HB3	3:A:1628:HOH:O	2.11	0.49
1:A:375:ARG:HG3	1:A:375:ARG:NH2	2.27	0.49
1:A:158:LEU:HB3	1:A:236:TYR:CD2	2.48	0.49
1:A:294:SER:HB3	1:A:363[B]:ARG:NH1	2.27	0.49
1:A:74:PHE:HB3	1:A:93:TYR:CE1	2.47	0.49
1:B:16:ASN:HB3	1:B:20:PHE:CE2	2.48	0.49
1:B:26:HIS:HD2	1:B:43:THR:OG1	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:39:ASN:HA	1:B:44:ASP:OD1	2.13	0.48
1:A:1:MET:HB3	3:A:1389:HOH:O	2.12	0.48
1:A:32:ASN:HD21	1:A:328[B]:GLU:HG2	1.78	0.48
1:B:324:ARG:HG2	1:B:324:ARG:HH21	1.78	0.48
1:A:198:ASP:HB3	1:A:201:ASN:OD1	2.14	0.47
1:B:247:LYS:HE2	1:B:296:LEU:HD21	1.96	0.47
1:B:147:ILE:O	1:B:150[A]:CYS:SG	2.72	0.47
1:A:139:ASP:OD1	1:A:140:PRO:HD2	2.15	0.47
1:A:87:HIS:H	1:A:87:HIS:CD2	2.32	0.47
1:B:168:GLU:CD	1:B:171:ARG:NH1	2.68	0.47
1:A:181:ARG:CZ	1:A:241:ASP:HB3	2.45	0.47
1:A:264:TYR:HA	1:A:277:ARG:HA	1.97	0.47
1:A:365:GLU:HG3	1:A:366[A]:ARG:NH2	2.30	0.47
1:A:366[B]:ARG:HB2	1:A:368:VAL:HG23	1.96	0.47
1:B:377:ARG:HD3	1:B:377:ARG:C	2.35	0.47
1:A:76:GLU:O	1:A:80:ARG:HG2	2.16	0.46
1:A:147:ILE:HG22	1:A:149:ASP:OD1	2.16	0.46
1:B:67:ALA:O	1:B:71:VAL:HG23	2.15	0.46
1:A:84:LEU:HG	1:A:89:ILE:CD1	2.46	0.46
1:B:15:ARG:NH2	1:B:325:ASP:OD1	2.49	0.45
1:A:3:GLN:HG3	1:A:4:GLN:N	2.31	0.45
1:B:206:ARG:HD2	3:B:1112:HOH:O	2.16	0.45
1:A:83:ASN:N	1:A:83:ASN:ND2	2.64	0.45
1:A:99:ALA:O	1:A:103:VAL:HG23	2.17	0.45
1:A:154:LEU:N	1:A:155:PRO:CD	2.79	0.45
1:B:168:GLU:OE1	1:B:171:ARG:NH1	2.50	0.45
1:B:375:ARG:HH21	1:B:375:ARG:HG3	1.82	0.45
1:A:291:GLU:OE2	1:A:363[A]:ARG:NH2	2.50	0.44
1:B:124:TRP:HA	1:B:131:ILE:HD13	2.00	0.44
1:A:298:GLU:CD	1:A:298:GLU:H	2.21	0.44
1:B:52:TRP:CE2	1:B:67:ALA:HB1	2.52	0.44
1:A:223:GLN:O	1:A:227:ILE:HG13	2.18	0.44
1:A:298:GLU:CD	1:A:298:GLU:N	2.71	0.44
1:A:53:LEU:HB3	1:A:362:LEU:CD2	2.48	0.44
1:B:122[B]:ARG:HG3	1:B:122[B]:ARG:NH1	2.32	0.44
1:B:234[A]:SER:OG	1:B:243:LEU:HB2	2.17	0.44
1:A:161:GLY:O	1:A:165:GLY:N	2.50	0.43
1:A:26:HIS:HD2	1:A:43:THR:OG1	2.00	0.43
1:A:98:LYS:O	1:A:102:ILE:HG13	2.19	0.43
1:B:188:ASP:O	1:B:189:ASP:HB2	2.18	0.43
1:A:65:GLU:O	1:A:69[A]:ARG:HG3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:290:LEU:HD13	1:A:363[B]:ARG:HD3	2.00	0.43
1:A:146:ILE:O	1:A:146:ILE:HG13	2.18	0.43
1:A:366[A]:ARG:HB2	1:A:368:VAL:HG23	1.99	0.43
1:B:121:MET:HE1	1:B:168:GLU:HB3	1.99	0.43
1:B:362:LEU:HB3	1:B:368:VAL:HB	2.01	0.43
1:A:16:ASN:HB3	1:A:20:PHE:CE2	2.54	0.43
1:A:181:ARG:HD3	3:A:1209:HOH:O	2.17	0.43
1:B:57:TYR:HB2	1:B:362:LEU:HD11	1.99	0.43
1:A:83:ASN:HD22	1:A:83:ASN:H	1.66	0.43
1:A:164:THR:C	1:A:166:ASP:N	2.71	0.43
1:A:130:ILE:HD12	1:A:130:ILE:C	2.39	0.43
1:A:362:LEU:HB3	1:A:368:VAL:HB	2.00	0.43
1:B:86:HIS:CD2	1:B:88:ASN:H	2.33	0.43
1:A:25:PRO:HA	1:A:36:LEU:HD23	1.99	0.43
1:B:247:LYS:CE	1:B:296:LEU:HD21	2.49	0.43
1:B:152:LEU:HD13	1:B:232:LEU:HD12	2.00	0.42
1:B:264:TYR:HA	1:B:277:ARG:HA	2.01	0.42
1:B:374:GLU:HG3	3:B:1068:HOH:O	2.19	0.42
1:A:234[A]:SER:OG	1:A:243:LEU:HB2	2.19	0.42
1:B:121:MET:HG2	1:B:157:LEU:HD22	2.00	0.42
1:B:262:VAL:HG21	1:B:321:TYR:CD2	2.54	0.42
1:A:344:ILE:HD13	2:C:2:RAM:O2	2.20	0.42
1:B:147:ILE:HG12	1:B:194:THR:HG22	2.01	0.42
1:A:2:TRP:O	1:A:6:ILE:HG13	2.19	0.42
1:A:84:LEU:O	1:A:89:ILE:HD11	2.19	0.42
1:B:28:SER:HG	1:B:348:ASP:HB3	1.83	0.41
1:A:365:GLU:HG3	1:A:366[A]:ARG:HH21	1.85	0.41
1:A:101:TRP:CE3	1:A:110:ARG:HB2	2.55	0.41
1:A:22:ASP:HA	1:A:69[A]:ARG:HH12	1.86	0.41
1:A:80:ARG:O	1:A:81:PHE:C	2.59	0.41
1:A:158:LEU:HD13	1:A:236:TYR:HB2	2.02	0.41
1:A:185:VAL:HB	1:B:187:GLY:HA3	2.02	0.41
1:A:350:THR:HB	1:A:352:TRP:CE2	2.56	0.41
1:B:15:ARG:NH1	1:B:328:GLU:N	2.69	0.41
1:A:288:GLY:O	1:A:292:ILE:HG13	2.20	0.41
1:B:15:ARG:O	1:B:19[A]:LYS:HE3	2.21	0.41
1:A:181:ARG:NH1	3:A:1415:HOH:O	2.42	0.41
1:A:231:ALA:O	1:A:234[B]:SER:OG	2.32	0.41
1:A:280:SER:O	1:A:284:ILE:HG13	2.20	0.41
1:B:83:ASN:O	1:B:86:HIS:HE1	2.04	0.40
1:A:37:ASN:OD1	1:A:44:ASP:OD2	2.40	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:67:ALA:O	1:A:71:VAL:HG23	2.20	0.40
1:A:85:ASP:OD2	1:A:123:ARG:NH1	2.55	0.40
1:A:105:LYS:NZ	3:A:1690:HOH:O	2.55	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	385/377 (102%)	356 (92%)	25 (6%)	4 (1%)	15 6
1	B	387/377 (103%)	372 (96%)	15 (4%)	0	100 100
All	All	772/754 (102%)	728 (94%)	40 (5%)	4 (0%)	34 18

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	88	ASN
1	A	141[A]	GLU
1	A	141[B]	GLU
1	A	273	PRO

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	314/304 (103%)	306 (98%)	8 (2%)	47 39
1	B	316/304 (104%)	303 (96%)	13 (4%)	30 20
All	All	630/608 (104%)	609 (97%)	21 (3%)	41 28

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	10	LEU
1	A	39	ASN
1	A	41	ASP
1	A	84	LEU
1	A	87	HIS
1	A	168	GLU
1	A	228	TYR
1	B	19[A]	LYS
1	B	19[B]	LYS
1	B	39	ASN
1	B	75[A]	ARG
1	B	75[B]	ARG
1	B	87	HIS
1	B	118	ASP
1	B	121	MET
1	B	127	ASP
1	B	210[A]	HIS
1	B	210[B]	HIS
1	B	228	TYR
1	B	377	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	4	GLN
1	A	26	HIS
1	A	39	ASN
1	A	62	GLN
1	A	83	ASN
1	A	176	HIS
1	A	223	GLN
1	B	26	HIS

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Mol	Chain	Res	Type
1	B	39	ASN
1	B	62	GLN
1	B	83	ASN
1	B	86	HIS
1	B	176	HIS
1	B	203	ASN
1	B	211	GLN
1	B	223	GLN
1	B	304	GLN

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)

4 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	BGC	C	1	2	12,12,12	0.51	0	17,17,17	0.70	0
2	RAM	C	2	2	10,10,11	0.45	0	14,14,16	0.90	1 (7%)
2	BGC	C	3	2	11,11,12	0.41	0	15,15,17	1.52	3 (20%)
2	GAD	C	4	1,2	10,11,11	1.78	2 (20%)	13,15,15	2.03	4 (30%)

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	BGC	C	1	2	-	0/2/22/22	0/1/1/1
2	RAM	C	2	2	-	-	1/1/1/1
2	BGC	C	3	2	-	0/2/19/22	0/1/1/1
2	GAD	C	4	1,2	-	1/4/17/17	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	4	GAD	C4-C5	4.30	1.39	1.33
2	C	4	GAD	C3-C4	2.18	1.53	1.50

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	4	GAD	O5-C5-C6	5.20	119.33	111.52
2	C	3	BGC	C3-C4-C5	-2.89	105.08	110.24
2	C	3	BGC	O4-C4-C5	2.70	115.99	109.30
2	C	4	GAD	C4-C5-C6	-2.63	117.98	123.65
2	C	4	GAD	O5-C5-C4	-2.52	122.69	124.81
2	C	4	GAD	C2-C3-C4	-2.36	109.09	112.32
2	C	3	BGC	C1-O5-C5	2.12	115.07	112.19
2	C	2	RAM	C6-C5-C4	-2.01	109.35	113.07

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	4	GAD	O5-C5-C6-O6A

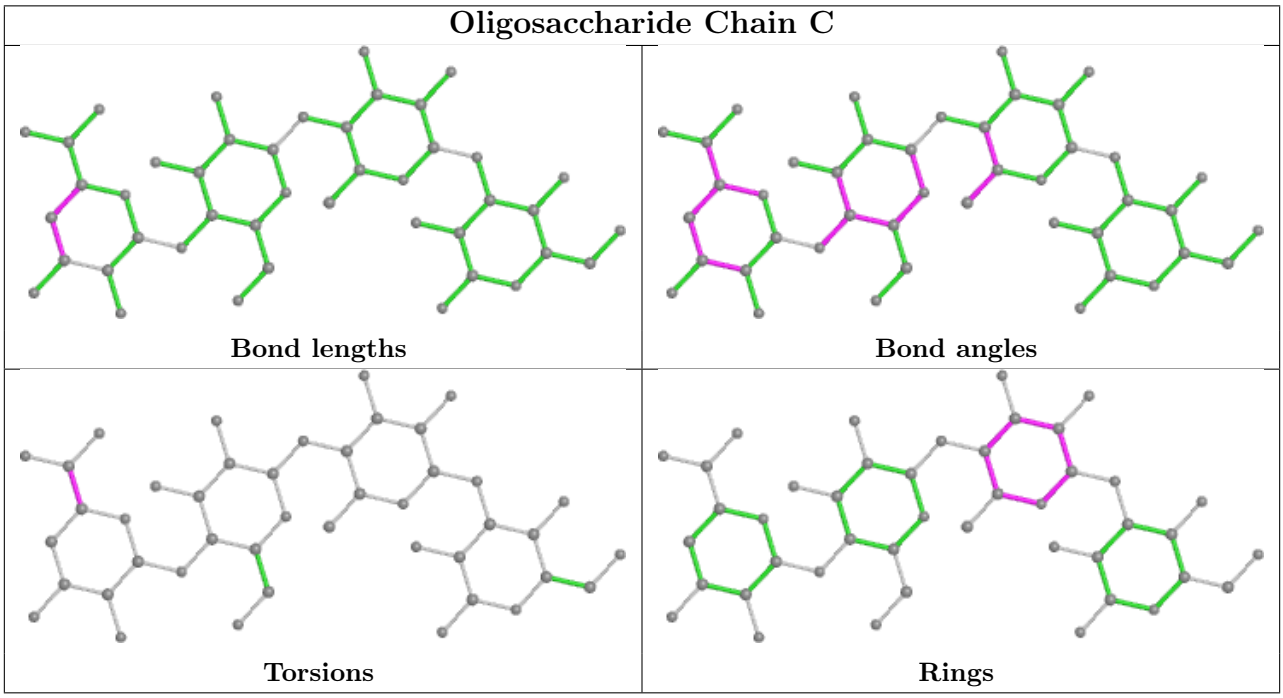
All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	2	RAM	C1-C2-C3-C4-C5-O5

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	3	BGC	2	0
2	C	2	RAM	1	0
2	C	1	BGC	4	0
2	C	4	GAD	2	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

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 95^{th} 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	377/377 (100%)	0.75	44 (11%) 4 5	8, 22, 49, 62	0
1	B	377/377 (100%)	0.12	6 (1%) 72 74	9, 17, 36, 54	0
All	All	754/754 (100%)	0.44	50 (6%) 18 20	8, 19, 44, 62	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	140	PRO	9.9
1	A	134	TRP	7.3
1	A	136	PRO	7.2
1	A	376	GLY	6.8
1	A	138	GLY	6.1
1	A	124	TRP	6.0
1	B	134[A]	TRP	5.9
1	A	135	GLY	5.6
1	A	164	THR	4.9
1	A	139	ASP	4.8
1	A	199	PRO	4.6
1	A	141[A]	GLU	4.2
1	A	125	ARG	4.1
1	A	137	LYS	4.0
1	A	3	GLN	3.8
1	B	140	PRO	3.8
1	A	122	ARG	3.6
1	A	126	ALA	3.5
1	A	297	ASP	3.5
1	A	101	TRP	3.5
1	A	128	ALA	3.4
1	B	141	GLU	3.3
1	A	377	ARG	3.3
1	B	136	PRO	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	85	ASP	3.3
1	A	142	ASN	3.1
1	A	298	GLU	3.0
1	A	299	SER	3.0
1	B	376	GLY	2.8
1	A	119	VAL	2.7
1	A	169	TYR	2.7
1	A	80	ARG	2.7
1	B	135	GLY	2.7
1	A	2	TRP	2.6
1	A	201	ASN	2.5
1	A	161	GLY	2.5
1	A	152	LEU	2.4
1	A	55	TYR	2.4
1	A	127	ASP	2.4
1	A	132	GLN	2.4
1	A	129	GLY	2.3
1	A	143	GLY	2.3
1	A	200	GLU	2.3
1	A	1	MET	2.3
1	A	375	ARG	2.2
1	A	305	ARG	2.2
1	A	168	GLU	2.2
1	A	300	ASP	2.2
1	A	352	TRP	2.1
1	A	131	ILE	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, 95^{th} 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	BGC	C	3	11/12	0.48	0.42	53,58,63,66	0
2	BGC	C	1	12/12	0.62	0.39	24,53,58,61	0

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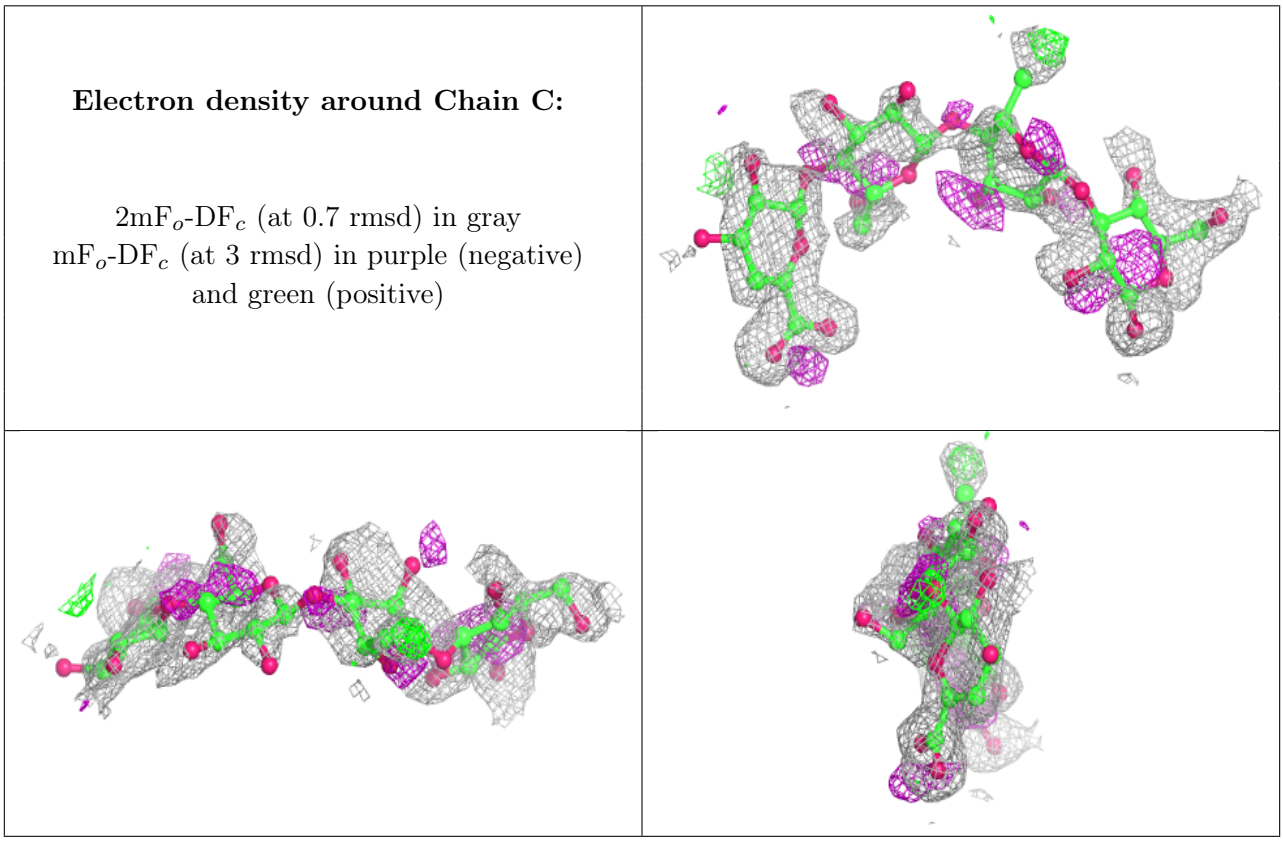
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	RAM	C	2	10/11	0.63	0.46	52,56,65,67	0
2	GAD	C	4	11/11	0.85	0.23	15,40,52,59	0

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)



6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.