



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:12 PM GMT

PDB ID : 4BPZ
Title : Crystal structure of lamA_E269S from Zobellia galactanivorans in complex with a trisaccharide of 1,3-1,4-beta-D-glucan.
Authors : Labourel, A.; Jam, M.; Jeudy, A.; Czjzek, M.; Michel, G.
Deposited on : 2013-05-29
Resolution : 1.13 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

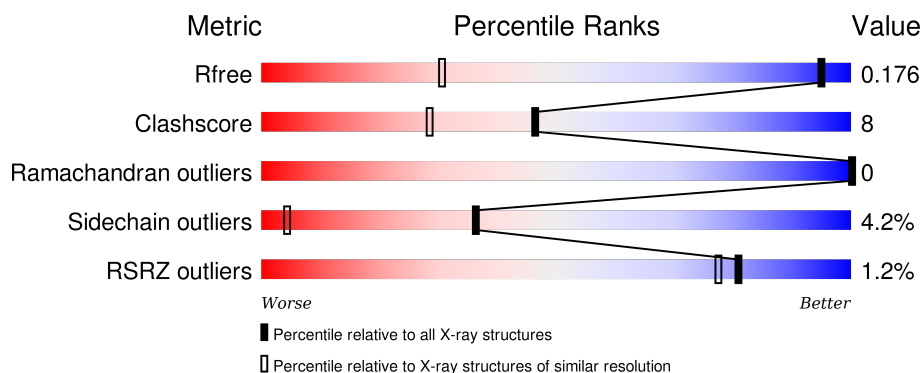
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

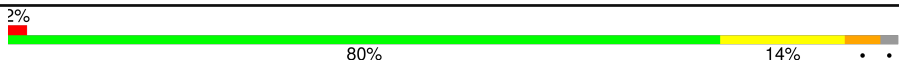

The reported resolution of this entry is 1.13 Å.

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}	91344	1158 (1.18-1.06)
Clashscore	102246	1215 (1.18-1.06)
Ramachandran outliers	100387	1162 (1.18-1.06)
Sidechain outliers	100360	1160 (1.18-1.06)
RSRZ outliers	91569	1161 (1.18-1.06)

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. 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	256	
1	B	256	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4855 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 ENDO-1,3-BETA-GLUCANASE, FAMILY GH16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	251	Total	C	N	O	S	0	18	0
			2133	1364	345	420	4			
1	B	252	Total	C	N	O	S	0	8	0
			2068	1318	337	409	4			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	128	HIS	-	EXPRESSION TAG	UNP G0L5X4
A	129	HIS	-	EXPRESSION TAG	UNP G0L5X4
A	130	HIS	-	EXPRESSION TAG	UNP G0L5X4
A	131	HIS	-	EXPRESSION TAG	UNP G0L5X4
A	132	HIS	-	EXPRESSION TAG	UNP G0L5X4
A	133	HIS	-	EXPRESSION TAG	UNP G0L5X4
A	134	GLY	-	EXPRESSION TAG	UNP G0L5X4
A	135	SER	-	EXPRESSION TAG	UNP G0L5X4
A	269	SER	GLU	ENGINEERED MUTATION	UNP G0L5X4
B	128	HIS	-	EXPRESSION TAG	UNP G0L5X4
B	129	HIS	-	EXPRESSION TAG	UNP G0L5X4
B	130	HIS	-	EXPRESSION TAG	UNP G0L5X4
B	131	HIS	-	EXPRESSION TAG	UNP G0L5X4
B	132	HIS	-	EXPRESSION TAG	UNP G0L5X4
B	133	HIS	-	EXPRESSION TAG	UNP G0L5X4
B	134	GLY	-	EXPRESSION TAG	UNP G0L5X4
B	135	SER	-	EXPRESSION TAG	UNP G0L5X4
B	269	SER	GLU	ENGINEERED MUTATION	UNP G0L5X4

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Ca	0	0
			1	1		

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

- Molecule 3 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	3	Total	C	O	0	0
			34	18	16		
3	B	3	Total	C	O	0	0
			34	18	16		

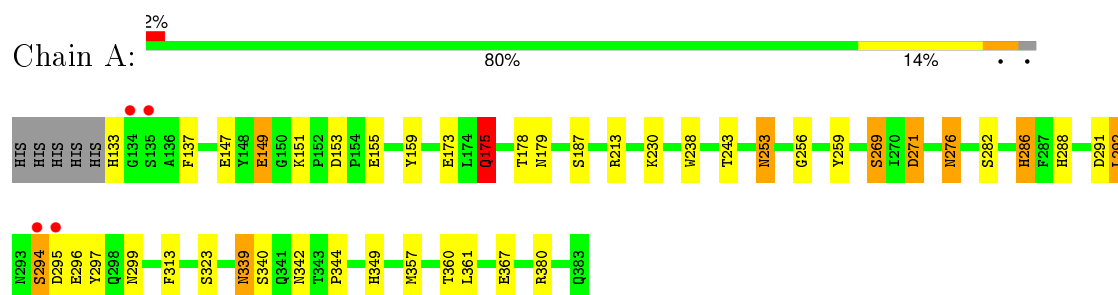
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	333	Total	O	0	0
			333	333		
4	B	251	Total	O	0	0
			251	251		

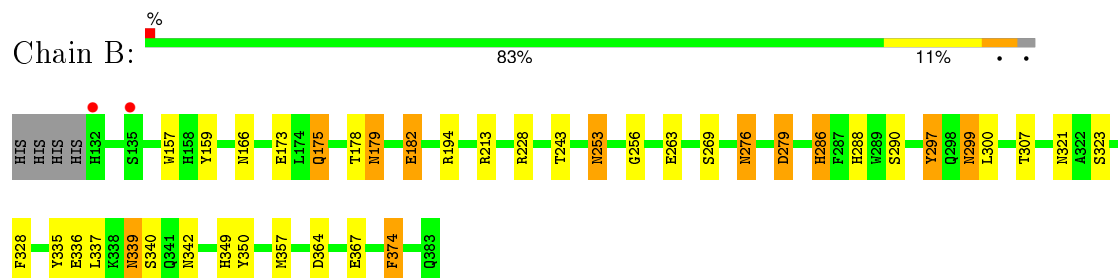
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: ENDO-1,3-BETA-GLUCANASE, FAMILY GH16



- Molecule 1: ENDO-1,3-BETA-GLUCANASE, FAMILY GH16



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	44.53 Å 76.48 Å 142.67 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	71.34 – 1.13 42.51 – 1.13	Depositor EDS
% Data completeness (in resolution range)	99.1 (71.34-1.13) 99.1 (42.51-1.13)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.53 (at 1.13 Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.142 , 0.175 0.144 , 0.176	Depositor DCC
R_{free} test set	9104 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	13.7	Xtriage
Anisotropy	0.450	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 181280 reflections	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	4855	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.90% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	1.29	12/2240 (0.5%)	1.23	17/3044 (0.6%)
1	B	1.27	6/2155 (0.3%)	1.24	16/2932 (0.5%)
All	All	1.28	18/4395 (0.4%)	1.24	33/5976 (0.6%)

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	2

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	147	GLU	CD-OE2	11.32	1.38	1.25
1	A	271[A]	ASP	CG-OD1	7.55	1.42	1.25
1	A	271[B]	ASP	CG-OD1	7.55	1.42	1.25
1	A	149[A]	GLU	CD-OE2	-7.41	1.17	1.25
1	A	149[B]	GLU	CD-OE2	-7.41	1.17	1.25
1	A	367	GLU	CG-CD	6.67	1.61	1.51
1	B	297	TYR	CG-CD2	6.42	1.47	1.39
1	A	344	PRO	C-O	6.08	1.35	1.23
1	B	269[A]	SER	CB-OG	-5.99	1.34	1.42
1	B	269[B]	SER	CB-OG	-5.99	1.34	1.42
1	A	367	GLU	CD-OE2	5.60	1.31	1.25
1	A	367	GLU	CD-OE1	5.58	1.31	1.25
1	B	290	SER	CA-CB	5.53	1.61	1.52
1	A	259	TYR	CE2-CZ	-5.43	1.31	1.38
1	A	137	PHE	CG-CD2	5.42	1.46	1.38
1	B	336	GLU	CD-OE1	-5.29	1.19	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	159	TYR	CD1-CE1	5.07	1.47	1.39
1	B	157	TRP	CD2-CE2	5.03	1.47	1.41

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	291	ASP	CB-CG-OD2	-11.95	107.55	118.30
1	A	291	ASP	CB-CG-OD1	9.99	127.29	118.30
1	A	292[A]	LEU	CB-CG-CD1	-8.87	95.93	111.00
1	A	292[B]	LEU	CB-CG-CD1	-8.87	95.93	111.00
1	B	374	PHE	CB-CG-CD2	-7.62	115.47	120.80
1	A	271[A]	ASP	CB-CG-OD1	-7.23	111.79	118.30
1	A	271[B]	ASP	CB-CG-OD1	-7.23	111.79	118.30
1	B	279	ASP	CB-CG-OD2	-7.15	111.86	118.30
1	B	228	ARG	NE-CZ-NH1	-7.05	116.77	120.30
1	B	182	GLU	CG-CD-OE1	6.91	132.12	118.30
1	A	269[A]	SER	N-CA-CB	-6.80	100.30	110.50
1	A	269[B]	SER	N-CA-CB	-6.80	100.30	110.50
1	B	350	TYR	CZ-CE2-CD2	-6.73	113.74	119.80
1	A	380	ARG	NE-CZ-NH2	-6.52	117.04	120.30
1	B	350	TYR	CG-CD2-CE2	6.44	126.45	121.30
1	A	259	TYR	CB-CG-CD1	6.11	124.67	121.00
1	B	159	TYR	CB-CG-CD2	-6.07	117.36	121.00
1	A	213	ARG	NE-CZ-NH2	-5.92	117.34	120.30
1	B	364[A]	ASP	CB-CG-OD1	5.87	123.58	118.30
1	B	364[B]	ASP	CB-CG-OD1	5.87	123.58	118.30
1	B	299	ASN	N-CA-CB	-5.83	100.10	110.60
1	B	328	PHE	CB-CG-CD1	-5.81	116.73	120.80
1	A	271[A]	ASP	CB-CG-OD2	5.78	123.50	118.30
1	A	271[B]	ASP	CB-CG-OD2	5.78	123.50	118.30
1	B	182	GLU	OE1-CD-OE2	-5.62	116.56	123.30
1	B	213	ARG	NE-CZ-NH1	5.46	123.03	120.30
1	B	159	TYR	CB-CG-CD1	5.45	124.27	121.00
1	A	159	TYR	CB-CG-CD2	-5.27	117.84	121.00
1	A	213	ARG	NE-CZ-NH1	5.27	122.93	120.30
1	B	335	TYR	CB-CG-CD2	-5.24	117.86	121.00
1	A	153	ASP	CB-CG-OD1	5.15	122.94	118.30
1	B	178	THR	CA-CB-CG2	-5.05	105.33	112.40
1	A	151	LYS	CD-CE-NZ	5.04	123.30	111.70

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	175	GLN	Peptide
1	A	271[A]	ASP	Sidechain

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	2133	0	2001	37	1
1	B	2068	0	1924	25	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	34	0	30	0	0
3	B	34	0	30	2	0
4	A	333	0	0	14	1
4	B	251	0	0	3	0
All	All	4855	0	3985	63	2

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 8.

All (63) 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:294[A]:SER:HB3	4:A:2244:HOH:O	1.32	1.26
1:A:360[B]:THR:HG22	4:A:2070:HOH:O	1.41	1.21
1:A:286[A]:HIS:CE1	1:A:299:ASN:OD1	2.17	0.98
1:A:286[B]:HIS:CE1	1:A:299:ASN:OD1	2.17	0.98
1:A:286[B]:HIS:HE1	1:A:299:ASN:OD1	1.46	0.98
1:A:286[A]:HIS:HE1	1:A:299:ASN:OD1	1.47	0.97
1:A:296:GLU:O	4:A:2241:HOH:O	1.88	0.90
1:A:294[B]:SER:HB3	4:A:2244:HOH:O	1.81	0.80
1:A:155:GLU:HG2	4:A:2038:HOH:O	1.81	0.79
1:B:276:ASN:HD21	1:B:279:ASP:H	1.29	0.79
1:A:286[C]:HIS:CE1	1:A:299:ASN:OD1	2.37	0.77
1:B:276:ASN:ND2	1:B:279:ASP:H	1.85	0.74
1:A:253:ASN:HD21	1:A:256:GLY:H	1.41	0.67
1:A:286[C]:HIS:HD2	4:A:2236:HOH:O	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:HIS:N	4:A:2002:HOH:O	2.30	0.65
1:A:173:GLU:HB3	1:A:175:GLN:HE22	1.62	0.64
1:B:173:GLU:HB3	1:B:175:GLN:HE22	1.63	0.64
1:B:253:ASN:HD21	1:B:256:GLY:H	1.45	0.62
1:A:294[A]:SER:CB	4:A:2244:HOH:O	2.13	0.62
1:B:175:GLN:NE2	1:B:357:MET:H	1.99	0.60
1:A:360[B]:THR:CG2	4:A:2070:HOH:O	2.20	0.57
1:B:179:ASN:C	1:B:179:ASN:HD22	2.06	0.56
1:A:360[A]:THR:HG21	4:A:2067:HOH:O	2.06	0.56
1:B:288:HIS:HD2	1:B:297:TYR:OH	1.89	0.55
1:A:175:GLN:NE2	1:A:357:MET:H	2.04	0.55
1:B:300[A]:LEU:HD11	1:B:337:LEU:HD13	1.88	0.55
1:B:276:ASN:C	1:B:276:ASN:HD22	2.09	0.54
1:B:263[A]:GLU:HG3	3:B:403:BGC:O3	2.07	0.54
1:B:175:GLN:HE21	1:B:357:MET:H	1.57	0.53
1:B:286:HIS:HD2	4:B:2173:HOH:O	1.92	0.52
1:A:253:ASN:ND2	1:A:256:GLY:H	2.08	0.52
1:A:286[C]:HIS:HE1	1:A:299:ASN:OD1	1.81	0.52
1:A:294[B]:SER:OG	1:A:295:ASP:N	2.44	0.51
1:B:253:ASN:ND2	1:B:256:GLY:H	2.09	0.51
1:A:339:ASN:HD22	1:A:340:SER:H	1.56	0.51
1:A:286[B]:HIS:HD2	4:A:2236:HOH:O	1.94	0.51
1:A:286[A]:HIS:HD2	4:A:2236:HOH:O	1.95	0.50
1:A:175:GLN:HE21	1:A:357:MET:H	1.59	0.50
1:B:339:ASN:HD22	1:B:340:SER:H	1.58	0.50
1:B:286:HIS:CE1	1:B:299:ASN:OD1	2.65	0.49
1:A:276:ASN:HD22	1:A:276:ASN:C	2.15	0.49
1:A:243:THR:OG1	1:A:349:HIS:HD2	1.96	0.49
1:B:286:HIS:HE1	1:B:299:ASN:OD1	1.96	0.48
1:A:269[B]:SER:OG	4:A:2217:HOH:O	1.65	0.48
1:B:175:GLN:NE2	1:B:175:GLN:H	2.13	0.47
1:A:253:ASN:C	1:A:253:ASN:HD22	2.19	0.45
3:B:401:GLC:H2	4:B:2161:HOH:O	2.15	0.45
1:B:243:THR:OG1	1:B:349:HIS:HD2	1.99	0.45
1:A:238:TRP:CE2	1:A:361:LEU:HB2	2.52	0.44
1:A:253:ASN:C	1:A:253:ASN:ND2	2.71	0.43
1:A:149[B]:GLU:HG3	1:A:187:SER:HA	2.01	0.43
1:B:321:ASN:OD1	1:B:323[A]:SER:HB2	2.19	0.42
1:A:286[C]:HIS:CD2	4:A:2236:HOH:O	2.61	0.42
1:B:340:SER:OG	1:B:342[B]:ASN:ND2	2.52	0.42
1:B:179:ASN:C	1:B:179:ASN:ND2	2.73	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:288:HIS:HE1	4:B:2158:HOH:O	2.02	0.41
1:B:253:ASN:HD22	1:B:253:ASN:C	2.24	0.41
1:A:288:HIS:HD2	1:A:297:TYR:OH	2.03	0.41
1:B:182:GLU:O	1:B:194:ARG:HD2	2.20	0.41
1:A:230:LYS:HE3	1:A:313:PHE:CE2	2.56	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:323[B]:SER:OG	4:A:2317:HOH:O[3_544]	2.01	0.19
1:B:166:ASN:OD1	1:B:307:THR:OG1[1_655]	2.08	0.12

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	268/256 (105%)	259 (97%)	9 (3%)	0	100	100
1	B	258/256 (101%)	248 (96%)	10 (4%)	0	100	100
All	All	526/512 (103%)	507 (96%)	19 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	235/221 (106%)	220 (94%)	15 (6%)	22	1
1	B	225/221 (102%)	217 (96%)	8 (4%)	42	6
All	All	460/442 (104%)	437 (95%)	23 (5%)	36	3

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	175	GLN
1	A	178	THR
1	A	179	ASN
1	A	253	ASN
1	A	276	ASN
1	A	282[A]	SER
1	A	282[B]	SER
1	A	286[A]	HIS
1	A	286[B]	HIS
1	A	286[C]	HIS
1	A	294[A]	SER
1	A	294[B]	SER
1	A	339	ASN
1	A	342[A]	ASN
1	A	342[B]	ASN
1	B	175	GLN
1	B	179	ASN
1	B	253	ASN
1	B	276	ASN
1	B	286	HIS
1	B	339	ASN
1	B	367	GLU
1	B	374	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	175	GLN
1	A	179	ASN
1	A	215	ASN
1	A	253	ASN
1	A	276	ASN
1	A	288	HIS
1	A	339	ASN

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Mol	Chain	Res	Type
1	A	349	HIS
1	A	383	GLN
1	B	175	GLN
1	B	179	ASN
1	B	215	ASN
1	B	253	ASN
1	B	276	ASN
1	B	286	HIS
1	B	288	HIS
1	B	298	GLN
1	B	339	ASN
1	B	349	HIS
1	B	383	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

6 carbohydrates 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	$\# Z > 2$
3	GLC	A	401	3	12,12,12	2.34	4 (33%)	17,17,17	1.46	4 (23%)
3	BGC	A	402	3	11,11,12	1.21	1 (9%)	14,15,17	1.02	0
3	BGC	A	403	3	11,11,12	0.76	0	14,15,17	1.26	2 (14%)
3	GLC	B	401	3	12,12,12	1.23	1 (8%)	17,17,17	1.59	5 (29%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	BGC	B	402	3	11,11,12	1.49	2 (18%)	14,15,17	0.80	0
3	BGC	B	403	3	11,11,12	1.51	1 (9%)	14,15,17	1.70	4 (28%)

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
3	GLC	A	401	3	-	0/2/22/22	0/1/1/1
3	BGC	A	402	3	-	0/2/19/22	0/1/1/1
3	BGC	A	403	3	-	0/2/19/22	0/1/1/1
3	GLC	B	401	3	-	0/2/22/22	0/1/1/1
3	BGC	B	402	3	-	0/2/19/22	0/1/1/1
3	BGC	B	403	3	-	0/2/19/22	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	401	GLC	O5-C1	-6.87	1.30	1.43
3	B	403	BGC	O5-C1	-3.96	1.37	1.43
3	B	402	BGC	C2-C3	-3.34	1.47	1.52
3	A	401	GLC	C4-C3	-2.03	1.47	1.52
3	A	402	BGC	C4-C3	2.10	1.57	1.52
3	A	401	GLC	C1-C2	2.17	1.57	1.52
3	B	402	BGC	C4-C3	2.18	1.58	1.52
3	B	401	GLC	C1-C2	2.33	1.57	1.52
3	A	401	GLC	O3-C3	2.71	1.49	1.43

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	GLC	C3-C4-C5	-3.20	104.61	110.20
3	B	401	GLC	O2-C2-C1	-2.96	103.30	109.82
3	B	401	GLC	C3-C4-C5	-2.55	105.74	110.20
3	A	401	GLC	C1-C2-C3	-2.42	106.83	110.43
3	B	403	BGC	C1-O5-C5	-2.28	109.36	112.25
3	B	401	GLC	O5-C1-C2	-2.21	106.27	109.80
3	A	403	BGC	O3-C3-C4	-2.01	105.81	110.34
3	A	401	GLC	O5-C1-C2	2.17	113.26	109.80
3	B	403	BGC	O2-C2-C3	2.33	114.81	110.12
3	A	401	GLC	C1-O5-C5	2.53	118.14	113.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	401	GLC	O1-C1-O5	2.54	117.20	110.25
3	B	403	BGC	O2-C2-C1	2.68	114.58	109.21
3	B	401	GLC	C1-O5-C5	2.76	118.58	113.47
3	A	403	BGC	O5-C5-C6	2.90	113.63	107.35
3	B	403	BGC	O5-C1-C2	2.91	115.57	110.86

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	401	GLC	1	0
3	B	403	BGC	1	0

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	251/256 (98%)	-0.30	4 (1%) 74 71	10, 15, 30, 60	1 (0%)
1	B	252/256 (98%)	-0.17	2 (0%) 87 84	13, 19, 32, 48	1 (0%)
All	All	503/512 (98%)	-0.24	6 (1%) 81 77	10, 17, 31, 60	2 (0%)

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	132	HIS	3.7
1	A	134	GLY	3.5
1	A	295	ASP	3.2
1	A	135	SER	2.7
1	B	135	SER	2.4
1	A	294[A]	SER	2.0

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	GLC	B	401	12/12	0.95	0.09	2.00	13,15,20,29	0
3	BGC	A	403	11/12	0.97	0.09	0.58	16,20,27,29	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	BGC	B	402	11/12	0.97	0.06	-0.20	17,18,20,20	0
3	GLC	A	401	12/12	0.97	0.06	-0.37	12,14,21,23	0
3	BGC	A	402	11/12	0.98	0.04	-1.35	14,15,17,18	0
3	BGC	B	403	11/12	0.96	0.11	-	23,26,35,36	0

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	CA	B	400	1/1	0.99	0.06	-0.87	23,23,23,23	1
2	CA	A	400	1/1	1.00	0.04	-0.94	16,16,16,16	1

6.5 Other polymers [i](#)

There are no such residues in this entry.