



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:20 PM UTC

PDB ID : 1N7R / pdb_00001n7r
Title : Streptococcus pneumoniae Hyaluronate Lyase W291A/W292A/F343V Mutant complex with hexasaccharide hyaluronan
Authors : Nukui, M.; Taylor, K.B.; McPherson, D.T.; Shigenaga, M.; Jedrzejas, M.J.
Deposited on : 2002-11-16
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

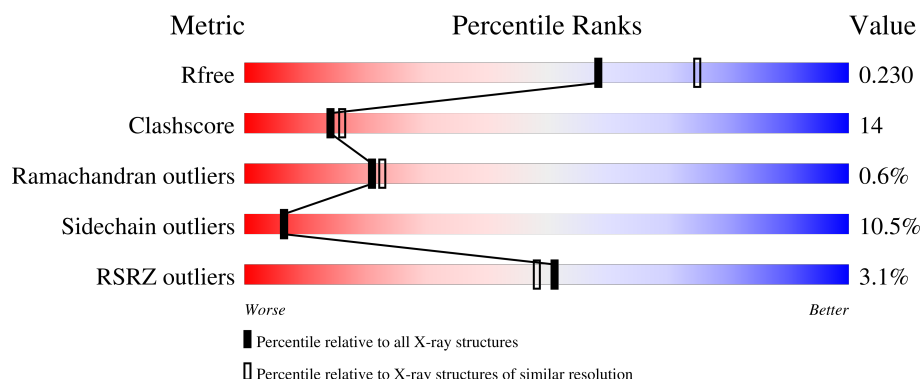
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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 $\leq 5\%$. 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	721	
2	B	6	

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	NAG	B	3	-	-	X	-
2	NAG	B	5	-	-	X	-
2	BDP	B	6	X	-	X	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6332 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 HYALURONIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	721	Total	C	N	O	S	0	0	0
			5762	3618	966	1156	22			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	291	ALA	TRP	engineered mutation	GB 437705
A	292	ALA	TRP	engineered mutation	GB 437705
A	343	VAL	PHE	engineered mutation	GB 437705
A	731	VAL	GLY	SEE REMARK 999	GB 437705

- Molecule 2 is an oligosaccharide called beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	6	Total	C	N	O	0	0	0
			79	42	3	34			

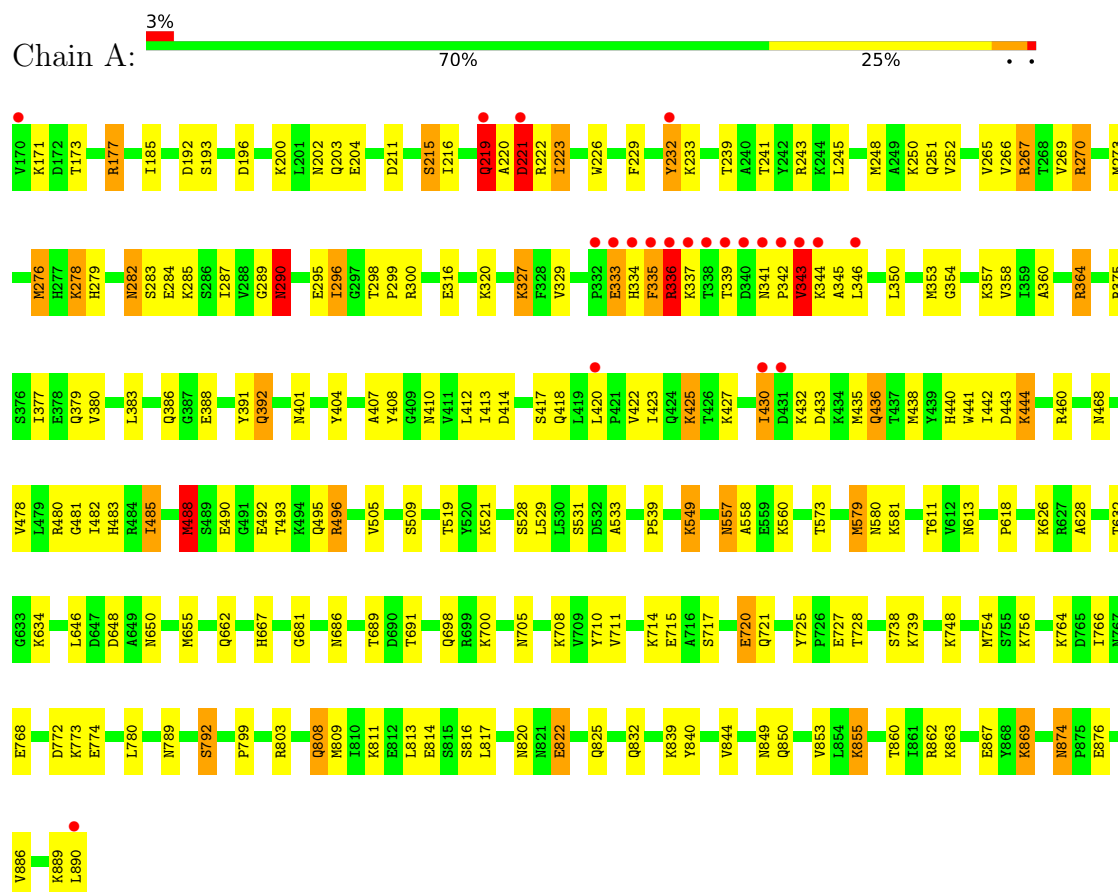
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	491	Total	O	0	0
			491	491		

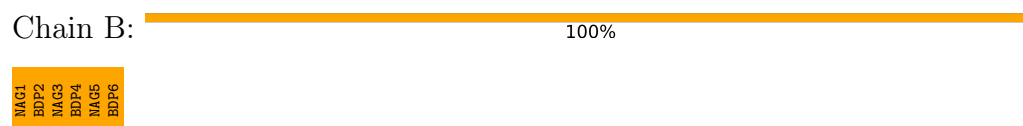
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: HYALURONIDASE



• Molecule 2: beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-beta-D-glucopyranuronic acid-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	84.49Å 102.49Å 103.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.20 20.00 – 2.21	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.20) 99.5 (20.00-2.21)	Depositor EDS
R_{merge}	0.21	Depositor
R_{sym}	0.21	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.43 (at 2.21Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.177 , 0.222 0.189 , 0.230	Depositor DCC
R_{free} test set	937 reflections (2.05%)	wwPDB-VP
Wilson B-factor (Å ²)	29.7	Xtriage
Anisotropy	0.042	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 43.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.012 for -h,l,k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6332	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.81% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BDP, NAG

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	1.63	55/5876 (0.9%)	1.34	21/7935 (0.3%)

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	4

All (55) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	276	MET	SD-CE	-15.90	1.39	1.79
1	A	632	THR	CA-CB	10.66	1.69	1.53
1	A	579	MET	SD-CE	8.99	2.02	1.79
1	A	267	ARG	NE-CZ	8.95	1.42	1.33
1	A	844	VAL	CA-CB	6.99	1.61	1.53
1	A	267	ARG	CG-CD	6.98	1.73	1.52
1	A	267	ARG	CD-NE	6.77	1.55	1.46
1	A	799	PRO	CA-C	6.72	1.59	1.52
1	A	229	PHE	CA-C	6.53	1.61	1.52
1	A	418	GLN	C-O	-6.50	1.16	1.24
1	A	648	ASP	CG-OD1	6.48	1.37	1.25
1	A	539	PRO	C-O	-6.48	1.14	1.23
1	A	792	SER	CB-OG	-6.35	1.29	1.42
1	A	655	MET	SD-CE	6.31	1.95	1.79
1	A	485	ILE	CA-CB	6.29	1.61	1.54
1	A	219	GLN	CA-C	6.23	1.61	1.52
1	A	221	ASP	CA-C	6.22	1.60	1.52
1	A	816	SER	C-O	-6.04	1.16	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	270	ARG	CD-NE	-6.04	1.37	1.46
1	A	766	ILE	CA-CB	6.03	1.61	1.54
1	A	444	LYS	C-O	-5.87	1.16	1.24
1	A	573	THR	CA-CB	5.87	1.64	1.53
1	A	558	ALA	CA-CB	-5.87	1.43	1.53
1	A	341	ASN	CA-C	5.85	1.60	1.52
1	A	488	MET	SD-CE	5.83	1.94	1.79
1	A	478	VAL	CA-CB	-5.79	1.47	1.54
1	A	273	MET	SD-CE	-5.68	1.65	1.79
1	A	809	MET	C-O	-5.61	1.17	1.24
1	A	442	ILE	C-O	-5.53	1.17	1.24
1	A	691	THR	CA-CB	5.48	1.62	1.53
1	A	267	ARG	CB-CG	5.45	1.68	1.52
1	A	628	ALA	CA-CB	5.40	1.61	1.53
1	A	840	TYR	CA-C	5.38	1.60	1.52
1	A	860	THR	C-O	5.37	1.30	1.24
1	A	388	GLU	C-O	5.30	1.30	1.23
1	A	233	LYS	CA-C	5.27	1.59	1.52
1	A	290	ASN	CA-C	5.26	1.58	1.52
1	A	216	ILE	C-O	-5.25	1.18	1.24
1	A	611	THR	CA-CB	5.25	1.61	1.53
1	A	531	SER	C-O	-5.21	1.16	1.24
1	A	193	SER	N-CA	5.20	1.52	1.46
1	A	533	ALA	CA-CB	5.19	1.62	1.53
1	A	245	LEU	C-O	-5.18	1.18	1.24
1	A	727	GLU	CA-C	5.18	1.59	1.53
1	A	343	VAL	CA-C	5.15	1.59	1.52
1	A	223	ILE	C-O	-5.14	1.18	1.24
1	A	509	SER	C-O	-5.13	1.17	1.24
1	A	204	GLU	CA-C	5.10	1.59	1.52
1	A	267	ARG	CA-CB	5.06	1.61	1.53
1	A	417	SER	CA-C	5.06	1.59	1.52
1	A	505	VAL	CA-CB	5.04	1.60	1.54
1	A	700	LYS	CA-C	5.04	1.59	1.52
1	A	460	ARG	N-CA	5.01	1.52	1.46
1	A	248	MET	SD-CE	5.01	1.92	1.79
1	A	725	TYR	CA-C	5.00	1.58	1.52

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	270	ARG	NE-CZ-NH2	-9.75	110.43	119.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	401	ASN	N-CA-C	7.60	120.11	110.61
1	A	221	ASP	CA-C-O	7.35	129.00	121.07
1	A	221	ASP	N-CA-C	6.03	118.79	110.35
1	A	482	ILE	N-CA-C	-5.97	104.92	110.53
1	A	792	SER	N-CA-CB	-5.92	100.51	110.99
1	A	270	ARG	NE-CZ-NH1	5.88	127.38	121.50
1	A	267	ARG	CA-CB-CG	5.85	125.79	114.10
1	A	289	GLY	N-CA-C	-5.84	104.27	111.45
1	A	221	ASP	CB-CA-C	5.81	118.79	109.61
1	A	270	ARG	CG-CD-NE	-5.81	99.22	112.00
1	A	283	SER	N-CA-C	5.78	119.59	112.54
1	A	267	ARG	NE-CZ-NH2	5.74	124.36	119.20
1	A	270	ARG	CD-NE-CZ	5.57	132.19	124.40
1	A	886	VAL	N-CA-C	5.50	119.03	113.47
1	A	728	THR	N-CA-C	5.36	117.45	109.25
1	A	748	LYS	N-CA-C	-5.14	101.03	109.40
1	A	267	ARG	N-CA-C	-5.08	105.82	111.36
1	A	192	ASP	CB-CA-C	5.05	118.00	110.62
1	A	266	VAL	CA-C-N	5.02	127.41	120.29
1	A	266	VAL	C-N-CA	5.02	127.41	120.29

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	219	GLN	Peptide
1	A	220	ALA	Peptide
1	A	221	ASP	Peptide
1	A	270	ARG	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	5762	0	5585	157	0
2	B	79	0	51	50	0
3	A	491	0	0	58	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	6332	0	5636	163	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 (163) 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:408:TYR:CE1	2:B:3:NAG:H82	1.44	1.52
1:A:579:MET:SD	1:A:579:MET:CE	2.02	1.48
1:A:580:ASN:ND2	2:B:2:BDP:O3	1.76	1.16
1:A:408:TYR:CZ	2:B:3:NAG:H82	1.82	1.14
1:A:243:ARG:NH2	2:B:5:NAG:H83	1.64	1.12
1:A:241:THR:HG21	1:A:276:MET:HE3	1.36	1.03
1:A:358:VAL:HG22	3:A:995:HOH:O	1.58	1.03
1:A:408:TYR:CE1	2:B:3:NAG:C8	2.41	1.03
3:A:1318:HOH:O	2:B:3:NAG:H62	1.55	1.02
1:A:613:ASN:H	1:A:698:GLN:HE22	1.07	1.02
1:A:243:ARG:CZ	2:B:5:NAG:H83	1.91	0.99
1:A:243:ARG:HH12	2:B:4:BDP:C6	1.77	0.97
1:A:483:HIS:HE1	3:A:1189:HOH:O	1.50	0.93
1:A:408:TYR:CD1	2:B:3:NAG:H82	2.04	0.93
1:A:822:GLU:HG3	3:A:927:HOH:O	1.70	0.90
1:A:488:MET:HB3	3:A:1384:HOH:O	1.73	0.88
1:A:243:ARG:NH1	2:B:4:BDP:O6A	2.08	0.86
1:A:250:LYS:HZ1	2:B:6:BDP:C3	1.88	0.86
1:A:773:LYS:HE3	3:A:1352:HOH:O	1.75	0.86
1:A:243:ARG:HD3	2:B:5:NAG:C8	2.07	0.84
1:A:267:ARG:HD2	3:A:1203:HOH:O	1.76	0.84
1:A:278:LYS:HG3	1:A:279:HIS:HD2	1.44	0.82
1:A:408:TYR:CZ	2:B:3:NAG:C8	2.62	0.81
1:A:267:ARG:CD	3:A:1203:HOH:O	2.27	0.81
1:A:243:ARG:HD3	2:B:5:NAG:H82	1.65	0.79
1:A:890:LEU:C	3:A:1054:HOH:O	2.26	0.77
1:A:243:ARG:CZ	2:B:5:NAG:C8	2.64	0.76
1:A:890:LEU:CB	3:A:1349:HOH:O	2.33	0.75
1:A:278:LYS:HG3	1:A:279:HIS:CD2	2.21	0.74
1:A:336:ARG:HH12	2:B:1:NAG:H4	1.52	0.73
1:A:185:ILE:O	2:B:6:BDP:O4	2.06	0.72
1:A:414:ASP:OD2	2:B:5:NAG:C6	2.38	0.72
1:A:557:ASN:HD22	1:A:557:ASN:C	1.98	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:243:ARG:CD	2:B:5:NAG:H82	2.23	0.68
1:A:221:ASP:HB3	3:A:1201:HOH:O	1.93	0.68
1:A:440:HIS:HE1	3:A:1068:HOH:O	1.77	0.67
1:A:414:ASP:OD2	2:B:5:NAG:H61	1.95	0.67
1:A:720:GLU:HG3	3:A:1308:HOH:O	1.94	0.67
1:A:243:ARG:HH21	2:B:5:NAG:H83	1.60	0.66
1:A:650:ASN:HD21	1:A:832:GLN:HE22	1.43	0.66
1:A:360:ALA:O	1:A:364:ARG:HG3	1.96	0.65
1:A:300:ARG:HH22	2:B:5:NAG:H4	1.61	0.65
1:A:354:GLY:HA3	1:A:377:ILE:HD11	1.79	0.65
1:A:243:ARG:NH1	2:B:4:BDP:C6	2.56	0.65
1:A:756:LYS:HE3	3:A:1229:HOH:O	1.97	0.64
1:A:708:LYS:HE3	1:A:715:GLU:OE2	1.98	0.64
3:A:1281:HOH:O	2:B:5:NAG:O4	2.12	0.64
1:A:241:THR:HG21	1:A:276:MET:CE	2.20	0.63
1:A:839:LYS:HD2	1:A:853:VAL:HG23	1.79	0.63
1:A:808:GLN:HG3	3:A:1383:HOH:O	1.99	0.63
1:A:560:LYS:HD3	3:A:1221:HOH:O	1.98	0.63
1:A:203:GLN:HG2	3:A:1390:HOH:O	1.99	0.62
1:A:243:ARG:CD	2:B:5:NAG:C8	2.75	0.62
1:A:219:GLN:C	3:A:903:HOH:O	2.42	0.62
1:A:232:TYR:HB3	3:A:1376:HOH:O	1.99	0.62
1:A:408:TYR:CD1	2:B:3:NAG:C8	2.79	0.61
1:A:241:THR:CG2	1:A:276:MET:HE3	2.21	0.60
1:A:519:THR:CG2	2:B:6:BDP:C6	2.80	0.59
3:A:1318:HOH:O	2:B:3:NAG:H4	2.02	0.58
1:A:243:ARG:NE	2:B:5:NAG:C8	2.67	0.58
1:A:444:LYS:NZ	3:A:914:HOH:O	2.36	0.57
1:A:392:GLN:HG2	3:A:1290:HOH:O	2.05	0.57
1:A:221:ASP:HB3	3:A:1255:HOH:O	2.05	0.57
1:A:243:ARG:NH1	2:B:5:NAG:HN2	2.03	0.57
1:A:850:GLN:HE21	1:A:863:LYS:HE3	1.70	0.56
1:A:278:LYS:HB3	3:A:1087:HOH:O	2.06	0.56
1:A:557:ASN:ND2	1:A:560:LYS:H	2.03	0.56
1:A:290:ASN:ND2	3:A:920:HOH:O	2.39	0.55
1:A:215:SER:HB3	1:A:226:TRP:CD1	2.42	0.55
1:A:436:GLN:HE21	1:A:436:GLN:N	2.03	0.55
1:A:414:ASP:OD2	2:B:5:NAG:O6	2.25	0.55
1:A:423:ILE:HG21	1:A:430:ILE:HD12	1.88	0.55
1:A:391:TYR:CD2	1:A:549:LYS:HG3	2.42	0.54
3:A:1318:HOH:O	2:B:3:NAG:C5	2.56	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:336:ARG:HH22	2:B:1:NAG:H62	1.73	0.54
1:A:519:THR:HG23	2:B:6:BDP:O6B	2.07	0.54
1:A:874:ASN:C	1:A:874:ASN:HD22	2.15	0.54
1:A:483:HIS:CE1	3:A:934:HOH:O	2.60	0.54
1:A:243:ARG:CZ	2:B:5:NAG:HN2	2.21	0.54
1:A:336:ARG:NH1	2:B:1:NAG:H4	2.22	0.53
1:A:250:LYS:NZ	3:A:905:HOH:O	2.25	0.52
1:A:343:VAL:HG13	1:A:344:LYS:N	2.24	0.52
1:A:443:ASP:OD2	1:A:496:ARG:NH2	2.43	0.52
1:A:855:LYS:CG	3:A:1333:HOH:O	2.57	0.52
1:A:285:LYS:O	1:A:327:LYS:NZ	2.43	0.51
1:A:521:LYS:HD2	3:A:913:HOH:O	2.11	0.51
1:A:689:THR:HA	3:A:1335:HOH:O	2.10	0.51
1:A:890:LEU:CB	3:A:1334:HOH:O	2.59	0.51
3:A:1318:HOH:O	2:B:3:NAG:C6	2.32	0.51
1:A:720:GLU:CG	3:A:1308:HOH:O	2.55	0.51
1:A:412:LEU:C	1:A:412:LEU:HD23	2.36	0.51
1:A:282:ASN:C	1:A:282:ASN:HD22	2.19	0.50
1:A:200:LYS:CE	3:A:969:HOH:O	2.59	0.50
1:A:820:ASN:HD22	1:A:825:GLN:HG2	1.75	0.50
1:A:295:GLU:C	1:A:296:ILE:HD13	2.36	0.50
1:A:296:ILE:HD11	1:A:335:PHE:CZ	2.47	0.50
1:A:557:ASN:C	1:A:557:ASN:ND2	2.64	0.50
1:A:579:MET:CE	1:A:579:MET:CG	2.89	0.50
1:A:350:LEU:HD21	1:A:379:GLN:HG2	1.93	0.49
1:A:689:THR:HG22	3:A:1335:HOH:O	2.13	0.49
1:A:354:GLY:CA	1:A:377:ILE:HD11	2.42	0.49
1:A:581:LYS:HB3	1:A:768:GLU:HB2	1.95	0.49
1:A:686:ASN:H	1:A:789:ASN:ND2	2.11	0.49
1:A:346:LEU:HD21	3:A:1158:HOH:O	2.12	0.48
1:A:822:GLU:CG	3:A:927:HOH:O	2.45	0.48
1:A:334:HIS:O	3:A:902:HOH:O	2.20	0.48
1:A:295:GLU:OE2	1:A:336:ARG:HA	2.13	0.48
1:A:764:LYS:HD2	1:A:772:ASP:HB3	1.95	0.48
1:A:874:ASN:HD22	1:A:876:GLU:H	1.63	0.47
1:A:250:LYS:NZ	2:B:6:BDP:C3	2.70	0.47
1:A:329:VAL:HG12	1:A:357:LYS:HD2	1.95	0.47
1:A:483:HIS:CE1	3:A:1189:HOH:O	2.39	0.47
1:A:335:PHE:CZ	1:A:353:MET:HB2	2.49	0.47
1:A:689:THR:CB	3:A:1335:HOH:O	2.62	0.47
1:A:243:ARG:NE	2:B:5:NAG:H83	2.26	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:890:LEU:CA	3:A:1054:HOH:O	2.62	0.47
1:A:667:HIS:O	1:A:681:GLY:HA2	2.15	0.46
3:A:1318:HOH:O	2:B:3:NAG:C4	2.63	0.46
1:A:855:LYS:CD	3:A:1333:HOH:O	2.63	0.46
1:A:243:ARG:NE	2:B:5:NAG:H82	2.29	0.46
1:A:667:HIS:HE1	3:A:1048:HOH:O	1.99	0.45
1:A:250:LYS:NZ	2:B:6:BDP:C4	2.80	0.45
1:A:869:LYS:HE2	3:A:1059:HOH:O	2.16	0.45
1:A:412:LEU:HD21	1:A:441:TRP:CZ3	2.52	0.45
1:A:346:LEU:HB2	3:A:943:HOH:O	2.17	0.45
1:A:267:ARG:HD3	3:A:1203:HOH:O	2.06	0.45
1:A:646:LEU:HD21	1:A:862:ARG:HB2	1.99	0.44
1:A:221:ASP:OD1	1:A:222:ARG:NH1	2.51	0.44
1:A:430:ILE:HG22	1:A:435:MET:HE3	1.99	0.44
1:A:634:LYS:HD3	3:A:1034:HOH:O	2.17	0.44
1:A:739:LYS:HE2	3:A:1050:HOH:O	2.18	0.44
1:A:343:VAL:HG21	2:B:1:NAG:H61	2.00	0.44
1:A:480:ARG:NH2	2:B:5:NAG:H61	2.33	0.44
1:A:425:LYS:HE2	3:A:1232:HOH:O	2.18	0.43
1:A:316:GLU:H	1:A:316:GLU:CD	2.25	0.43
1:A:440:HIS:CE1	3:A:1068:HOH:O	2.60	0.43
1:A:265:VAL:O	1:A:269:VAL:HG23	2.19	0.43
1:A:407:ALA:O	1:A:410:ASN:HB2	2.18	0.43
1:A:408:TYR:CZ	2:B:3:NAG:C7	3.01	0.43
1:A:211:ASP:O	1:A:215:SER:HB2	2.19	0.43
1:A:618:PRO:HG3	3:A:1163:HOH:O	2.19	0.42
1:A:173:THR:HG22	3:A:1035:HOH:O	2.18	0.42
1:A:202:ASN:ND2	1:A:251:GLN:HE22	2.16	0.42
1:A:711:VAL:CG2	1:A:754:MET:HE1	2.49	0.42
1:A:711:VAL:HG22	1:A:754:MET:HE1	2.02	0.42
1:A:298:THR:HB	1:A:299:PRO:HD3	2.02	0.42
1:A:386:GLN:NE2	3:A:960:HOH:O	2.52	0.42
1:A:333:GLU:O	1:A:345:ALA:HB3	2.20	0.42
1:A:480:ARG:HH21	2:B:5:NAG:H61	1.83	0.42
1:A:710:TYR:CE2	1:A:803:ARG:NH1	2.87	0.42
1:A:284:GLU:CD	1:A:284:GLU:N	2.78	0.42
1:A:177:ARG:HG3	1:A:422:VAL:HG13	2.02	0.42
1:A:243:ARG:NH1	2:B:5:NAG:N2	2.68	0.41
1:A:243:ARG:CD	2:B:5:NAG:H83	2.49	0.41
1:A:287:ILE:HD12	1:A:339:THR:CG2	2.51	0.41
1:A:822:GLU:CD	3:A:927:HOH:O	2.64	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:481:GLY:O	1:A:485:ILE:HG12	2.21	0.41
1:A:250:LYS:HE3	2:B:6:BDP:O3	2.21	0.41
1:A:549:LYS:HE2	3:A:955:HOH:O	2.21	0.41
1:A:343:VAL:HG11	2:B:1:NAG:H62	2.02	0.41
1:A:420:LEU:HD23	1:A:420:LEU:HA	1.93	0.41
1:A:874:ASN:ND2	1:A:876:GLU:H	2.18	0.40
1:A:662:GLN:HE21	1:A:662:GLN:HB2	1.63	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	719/721 (100%)	679 (94%)	36 (5%)	4 (1%)	21 23

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	219	GLN
1	A	342	PRO
1	A	336	ARG
1	A	343	VAL

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	636/638 (100%)	569 (90%)	67 (10%)	6 6

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	171	LYS
1	A	177	ARG
1	A	196	ASP
1	A	215	SER
1	A	219	GLN
1	A	223	ILE
1	A	232	TYR
1	A	239	THR
1	A	252	VAL
1	A	278	LYS
1	A	282	ASN
1	A	290	ASN
1	A	296	ILE
1	A	320	LYS
1	A	327	LYS
1	A	333	GLU
1	A	335	PHE
1	A	336	ARG
1	A	337	LYS
1	A	343	VAL
1	A	364	ARG
1	A	375	ARG
1	A	380	VAL
1	A	383	LEU
1	A	392	GLN
1	A	404	TYR
1	A	413	ILE
1	A	425	LYS
1	A	427	LYS
1	A	430	ILE
1	A	432	LYS
1	A	433	ASP
1	A	436	GLN
1	A	438	MET
1	A	468	ASN
1	A	488	MET
1	A	490	GLU
1	A	492	GLU

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Mol	Chain	Res	Type
1	A	493	THR
1	A	495	GLN
1	A	496	ARG
1	A	528	SER
1	A	529	LEU
1	A	549	LYS
1	A	557	ASN
1	A	626	LYS
1	A	705	ASN
1	A	714	LYS
1	A	717	SER
1	A	720	GLU
1	A	721	GLN
1	A	738	SER
1	A	774	GLU
1	A	780	LEU
1	A	792	SER
1	A	808	GLN
1	A	811	LYS
1	A	813	LEU
1	A	814	GLU
1	A	817	LEU
1	A	822	GLU
1	A	849	ASN
1	A	855	LYS
1	A	867	GLU
1	A	869	LYS
1	A	874	ASN
1	A	889	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	202	ASN
1	A	261	GLN
1	A	282	ASN
1	A	341	ASN
1	A	349	ASN
1	A	379	GLN
1	A	436	GLN
1	A	498	GLN
1	A	557	ASN

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Mol	Chain	Res	Type
1	A	580	ASN
1	A	662	GLN
1	A	667	HIS
1	A	685	GLN
1	A	698	GLN
1	A	759	GLN
1	A	789	ASN
1	A	820	ASN
1	A	825	GLN
1	A	832	GLN
1	A	850	GLN
1	A	874	ASN

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 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	B	1	2	15,15,15	2.20	7 (46%)	21,21,21	1.06	2 (9%)
2	BDP	B	2	2	12,12,13	3.77	4 (33%)	14,17,19	1.78	5 (35%)
2	NAG	B	3	2	14,14,15	2.39	7 (50%)	17,19,21	1.06	1 (5%)
2	BDP	B	4	2	12,12,13	3.67	6 (50%)	14,17,19	2.04	5 (35%)
2	NAG	B	5	2	14,14,15	2.54	7 (50%)	17,19,21	1.29	3 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	BDP	B	6	2	12,12,13	5.56	5 (41%)	14,17,19	4.31	8 (57%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	B	1	2	-	1/6/26/26	0/1/1/1
2	BDP	B	2	2	-	2/4/21/24	0/1/1/1
2	NAG	B	3	2	-	1/6/23/26	0/1/1/1
2	BDP	B	4	2	-	0/4/21/24	0/1/1/1
2	NAG	B	5	2	-	4/6/23/26	0/1/1/1
2	BDP	B	6	2	1/1/5/6	1/4/21/24	0/1/1/1

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	6	BDP	O4-C4	-14.49	1.07	1.43
2	B	6	BDP	C4-C5	-10.21	1.36	1.53
2	B	2	BDP	C4-C5	8.27	1.66	1.53
2	B	4	BDP	C5-C6	-7.54	1.37	1.53
2	B	2	BDP	C5-C6	-6.95	1.38	1.53
2	B	4	BDP	C4-C5	6.35	1.63	1.53
2	B	6	BDP	O5-C5	5.77	1.54	1.43
2	B	2	BDP	O5-C5	5.50	1.54	1.43
2	B	4	BDP	O5-C5	5.44	1.54	1.43
2	B	3	NAG	O5-C1	5.30	1.52	1.43
2	B	5	NAG	O5-C1	4.56	1.51	1.43
2	B	1	NAG	C2-N2	4.27	1.52	1.45
2	B	5	NAG	C2-N2	3.78	1.52	1.46
2	B	1	NAG	O5-C1	3.62	1.51	1.42
2	B	5	NAG	C1-C2	3.56	1.57	1.52
2	B	6	BDP	C5-C6	-3.49	1.45	1.53
2	B	5	NAG	O5-C5	3.36	1.50	1.43
2	B	5	NAG	C4-C5	3.31	1.60	1.53
2	B	3	NAG	O5-C5	3.30	1.49	1.43
2	B	3	NAG	C1-C2	3.24	1.56	1.52
2	B	1	NAG	C1-C2	3.10	1.56	1.52
2	B	4	BDP	C2-C3	3.07	1.57	1.52
2	B	4	BDP	O5-C1	2.83	1.48	1.43
2	B	2	BDP	C2-C3	2.79	1.56	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	NAG	C4-C5	2.79	1.59	1.53
2	B	3	NAG	C2-N2	2.74	1.50	1.46
2	B	1	NAG	C3-C2	2.54	1.57	1.53
2	B	3	NAG	C4-C5	2.51	1.58	1.53
2	B	5	NAG	C7-N2	2.34	1.41	1.34
2	B	1	NAG	O5-C5	2.34	1.50	1.44
2	B	3	NAG	C3-C2	2.31	1.57	1.52
2	B	4	BDP	O4-C4	2.27	1.48	1.43
2	B	1	NAG	C7-N2	2.27	1.41	1.34
2	B	5	NAG	C4-C3	2.13	1.57	1.52
2	B	3	NAG	C7-N2	2.11	1.41	1.34
2	B	6	BDP	O6A-C6	2.04	1.28	1.22

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	6	BDP	C2-C3-C4	12.17	132.27	110.86
2	B	6	BDP	O4-C4-C3	7.25	127.46	110.38
2	B	4	BDP	O4-C4-C5	-4.25	100.06	109.76
2	B	6	BDP	O4-C4-C5	4.17	119.28	109.76
2	B	6	BDP	C1-C2-C3	-3.60	104.40	109.64
2	B	4	BDP	O4-C4-C3	3.59	118.83	110.38
2	B	6	BDP	O5-C1-C2	-3.20	103.16	110.79
2	B	2	BDP	O4-C4-C5	-3.08	102.72	109.76
2	B	2	BDP	O4-C4-C3	2.80	116.99	110.38
2	B	4	BDP	C3-C4-C5	2.66	113.87	109.30
2	B	4	BDP	O6B-C6-O6A	2.65	130.09	124.08
2	B	2	BDP	C2-C3-C4	-2.57	106.35	110.86
2	B	2	BDP	C3-C4-C5	2.55	113.68	109.30
2	B	5	NAG	C4-C3-C2	-2.52	107.32	111.02
2	B	2	BDP	O6B-C6-O6A	2.50	129.76	124.08
2	B	5	NAG	C8-C7-N2	-2.49	111.99	116.12
2	B	5	NAG	O7-C7-N2	2.45	126.31	121.98
2	B	6	BDP	O6B-C6-O6A	2.44	129.61	124.08
2	B	6	BDP	O3-C3-C2	-2.29	105.39	110.05
2	B	1	NAG	C8-C7-N2	-2.24	112.39	116.12
2	B	1	NAG	O7-C7-N2	2.21	125.89	121.98
2	B	4	BDP	O5-C5-C6	2.20	114.43	106.59
2	B	6	BDP	C3-C4-C5	-2.08	105.73	109.30
2	B	3	NAG	C8-C7-N2	-2.07	112.68	116.12

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	6	BDP	C4

All (9) torsion outliers are listed below:

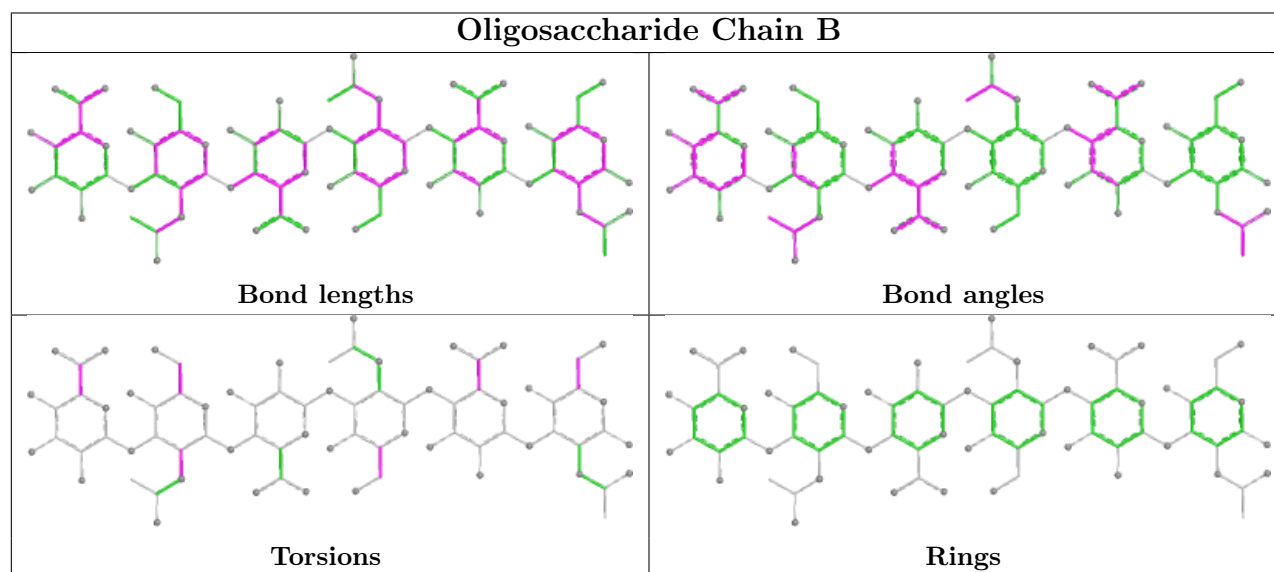
Mol	Chain	Res	Type	Atoms
2	B	2	BDP	C4-C5-C6-O6A
2	B	2	BDP	C4-C5-C6-O6B
2	B	5	NAG	C4-C5-C6-O6
2	B	5	NAG	O5-C5-C6-O6
2	B	1	NAG	O5-C5-C6-O6
2	B	3	NAG	O5-C5-C6-O6
2	B	5	NAG	C3-C2-N2-C7
2	B	5	NAG	C1-C2-N2-C7
2	B	6	BDP	O5-C5-C6-O6B

There are no ring outliers.

6 monomers are involved in 50 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	4	BDP	3	0
2	B	2	BDP	1	0
2	B	1	NAG	5	0
2	B	5	NAG	22	0
2	B	3	NAG	12	0
2	B	6	BDP	7	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 [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	721/721 (100%)	-0.11	22 (3%)	51	48	20, 31, 56, 110	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	342	PRO	5.0
1	A	219	GLN	4.6
1	A	334	HIS	4.6
1	A	335	PHE	4.4
1	A	343	VAL	4.3
1	A	337	LYS	3.6
1	A	336	ARG	3.5
1	A	232	TYR	3.4
1	A	344	LYS	3.4
1	A	338	THR	3.3
1	A	346	LEU	3.2
1	A	341	ASN	3.1
1	A	332	PRO	3.0
1	A	431	ASP	3.0
1	A	339	THR	3.0
1	A	890	LEU	2.6
1	A	420	LEU	2.5
1	A	221	ASP	2.5
1	A	430	ILE	2.2
1	A	333	GLU	2.1
1	A	340	ASP	2.1
1	A	170	VAL	2.0

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

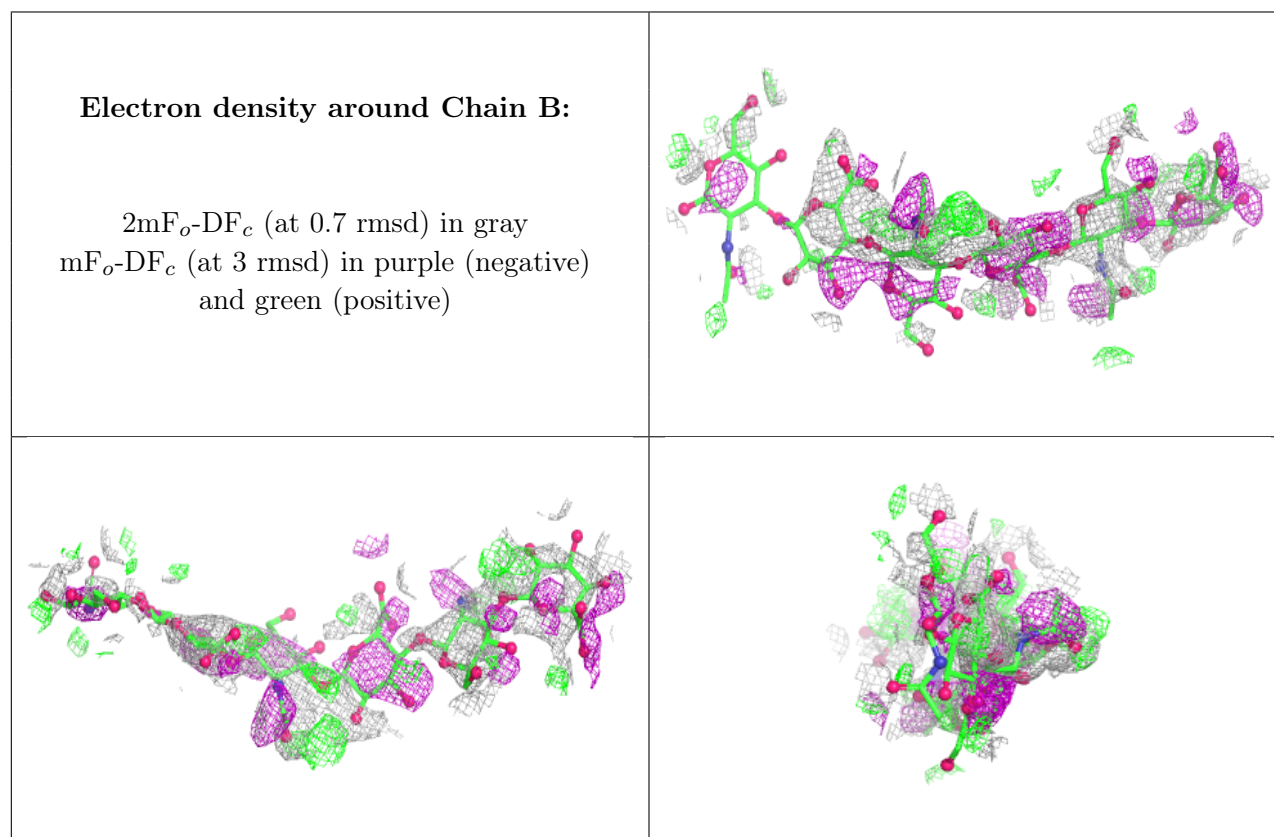
There are no non-standard protein/DNA/RNA residues in this entry.

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. 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($\text{\AA}^2$)	Q<0.9
2	NAG	B	1	15/15	0.04	0.23	94,96,97,97	0
2	BDP	B	4	12/13	0.11	0.29	88,91,94,95	0
2	NAG	B	5	14/15	0.30	0.24	86,89,91,92	0
2	NAG	B	3	14/15	0.40	0.28	81,87,88,90	0
2	BDP	B	2	12/13	0.45	0.23	81,87,90,91	0
2	BDP	B	6	12/13	0.63	0.23	92,94,94,95	0

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



6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.