



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 02:01 AM UTC

PDB ID : 4MTV / pdb_00004mtv
Title : Crystal structure of the complex of Buffalo Signalling Glycoprotein with pentasaccharide at 2.8Å resolution
Authors : Shukla, P.K.; Chaudhary, A.; Singh, A.K.; Sinha, M.; Kaur, P.; Sharma, S.; Singh, T.P.
Deposited on : 2013-09-20
Resolution : 2.80 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

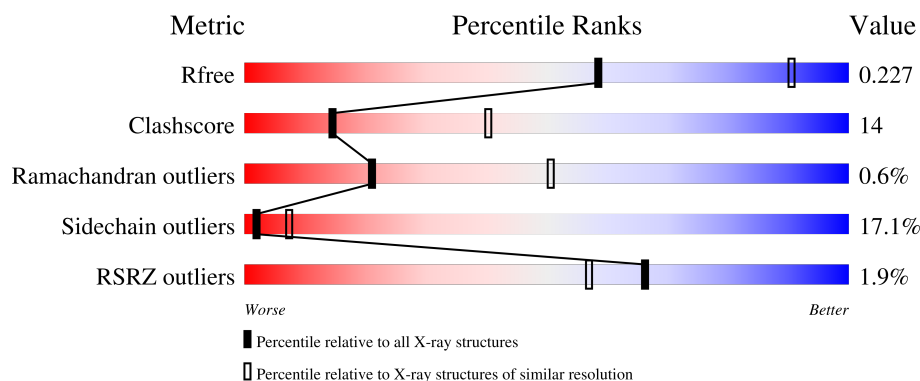
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	361	2% 64% 28% 7%
2	B	5	100%

2 Entry composition [i](#)

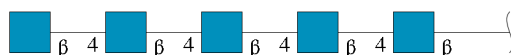
There are 4 unique types of molecules in this entry. The entry contains 3073 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Chitinase-3-like protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	361	Total	C	N	O	S	0	0	0
			2893	1851	507	526	9			

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	5	Total	C	N	O	0	0	0
			71	40	5	26			

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		

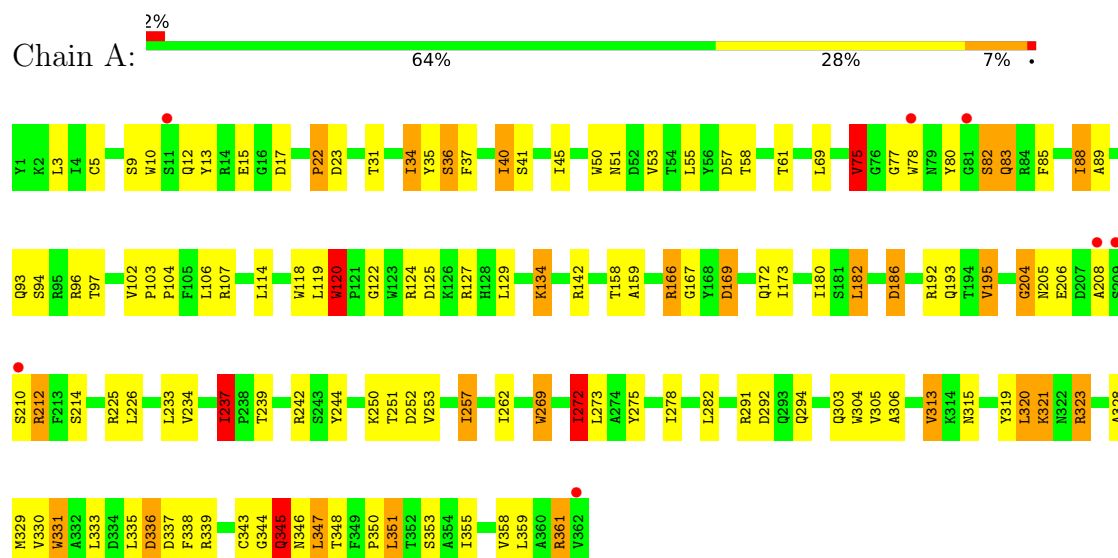
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	95	Total	O	0	0
			95	95		

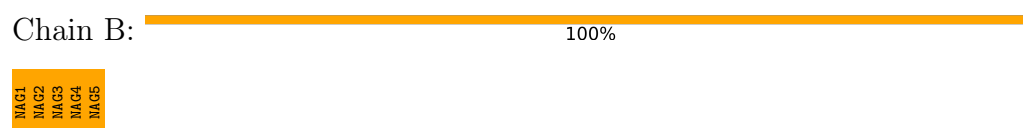
3 Residue-property plots [i](#)

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: Chitinase-3-like protein 1



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	62.87Å 66.50Å 107.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	56.50 – 2.80 56.50 – 2.80	Depositor EDS
% Data completeness (in resolution range)	92.4 (56.50-2.80) 92.4 (56.50-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.181 , 0.229 0.178 , 0.227	Depositor DCC
R_{free} test set	508 reflections (4.39%)	wwPDB-VP
Wilson B-factor (Å ²)	36.1	Xtriage
Anisotropy	0.283	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 66.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3073	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.69% 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: 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	0.93	19/2973 (0.6%)	1.11	26/4037 (0.6%)

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	34	ILE	C-O	-7.52	1.15	1.24
1	A	40	ILE	C-O	-7.50	1.16	1.24
1	A	35	TYR	C-O	-6.87	1.15	1.24
1	A	331	TRP	CA-C	-6.84	1.44	1.52
1	A	237	ILE	C-O	-6.72	1.16	1.23
1	A	331	TRP	NE1-CE2	-6.72	1.30	1.37
1	A	330	VAL	C-O	-6.53	1.17	1.24
1	A	180	ILE	C-O	-6.46	1.17	1.24
1	A	169	ASP	C-O	-6.41	1.15	1.23
1	A	122	GLY	C-O	-6.41	1.15	1.23
1	A	331	TRP	C-O	-6.08	1.17	1.24
1	A	214	SER	C-O	-6.07	1.16	1.24
1	A	272	ILE	C-O	-6.04	1.18	1.24
1	A	36	SER	C-O	-6.00	1.16	1.24
1	A	166	ARG	C-O	-5.60	1.17	1.24
1	A	320	LEU	C-O	-5.46	1.17	1.24
1	A	118	TRP	NE1-CE2	-5.44	1.31	1.37
1	A	120	TRP	NE1-CE2	-5.24	1.31	1.37
1	A	45	ILE	C-O	-5.05	1.18	1.24

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	89	ALA	N-CA-C	7.41	121.56	112.23
1	A	78	TRP	N-CA-C	7.26	120.11	111.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	212	ARG	N-CA-C	-7.22	104.21	113.16
1	A	82	SER	N-CA-C	-7.05	103.64	111.82
1	A	257	ILE	N-CA-C	6.56	117.75	108.17
1	A	167	GLY	N-CA-C	6.52	128.64	113.18
1	A	206	GLU	N-CA-C	-6.12	106.35	113.88
1	A	343	CYS	N-CA-C	6.06	120.97	112.72
1	A	125	ASP	N-CA-C	6.05	120.40	112.89
1	A	173	ILE	N-CA-C	5.99	118.58	111.09
1	A	51	ASN	N-CA-C	5.89	119.52	112.93
1	A	192	ARG	N-CA-C	-5.72	101.22	110.32
1	A	102	VAL	CA-C-N	-5.61	114.60	120.38
1	A	102	VAL	C-N-CA	-5.61	114.60	120.38
1	A	159	ALA	N-CA-C	5.60	121.00	113.72
1	A	336	ASP	N-CA-C	-5.57	102.51	110.59
1	A	361	ARG	N-CA-C	5.52	118.69	109.96
1	A	208	ALA	N-CA-C	5.51	119.18	111.30
1	A	335	LEU	N-CA-C	-5.42	106.03	113.30
1	A	294	GLN	N-CA-C	5.36	118.98	111.74
1	A	22	PRO	CA-C-N	5.34	128.83	120.82
1	A	22	PRO	C-N-CA	5.34	128.83	120.82
1	A	15	GLU	N-CA-C	5.32	117.73	110.55
1	A	338	PHE	N-CA-C	5.24	117.90	111.82
1	A	75	VAL	CB-CA-C	-5.21	103.45	111.31
1	A	31	THR	N-CA-C	-5.09	107.20	113.41

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2893	0	2817	79	0
2	B	71	0	63	14	0
3	A	14	0	13	0	0
4	A	95	0	0	4	0
All	All	3073	0	2893	82	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 (82) 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:321:LYS:NZ	1:A:358:VAL:O	2.00	0.94
1:A:269:TRP:CD1	2:B:3:NAG:HO6	1.87	0.91
1:A:269:TRP:CZ2	2:B:2:NAG:H3	2.07	0.90
2:B:4:NAG:O4	2:B:5:NAG:H83	1.75	0.85
1:A:83:GLN:HE21	1:A:83:GLN:H	1.21	0.84
1:A:57:ASP:O	1:A:61:THR:HG23	1.79	0.82
1:A:262:ILE:H	1:A:303:GLN:HE22	1.29	0.81
1:A:204:GLY:HA2	1:A:292:ASP:HB3	1.66	0.78
1:A:262:ILE:H	1:A:303:GLN:NE2	1.80	0.78
2:B:4:NAG:H62	2:B:5:NAG:C7	2.14	0.78
1:A:269:TRP:HD1	2:B:3:NAG:HO6	1.30	0.74
1:A:269:TRP:CZ3	1:A:272:ILE:HD13	2.23	0.73
1:A:269:TRP:HZ3	1:A:272:ILE:HD13	1.54	0.73
1:A:323:ARG:HH11	1:A:323:ARG:HG2	1.54	0.71
1:A:22:PRO:HB2	1:A:58:THR:HG22	1.71	0.71
1:A:292:ASP:HB2	4:A:585:HOH:O	1.91	0.70
1:A:244:TYR:HB3	1:A:257:ILE:HD12	1.76	0.68
1:A:313:VAL:HG13	1:A:355:ILE:HG13	1.76	0.68
1:A:195:VAL:HG23	1:A:257:ILE:HD11	1.77	0.67
1:A:319:TYR:CZ	1:A:323:ARG:HD2	2.30	0.67
1:A:83:GLN:H	1:A:83:GLN:NE2	1.92	0.66
1:A:278:ILE:CD1	1:A:305:VAL:HG11	2.25	0.66
1:A:278:ILE:HD11	1:A:305:VAL:HG11	1.79	0.65
1:A:186:ASP:OD1	1:A:242:ARG:NH1	2.31	0.64
1:A:169:ASP:OD2	1:A:172:GLN:HG3	1.99	0.61
1:A:93:GLN:HE21	1:A:96:ARG:NH2	1.99	0.61
1:A:323:ARG:HG2	1:A:323:ARG:NH1	2.13	0.61
1:A:331:TRP:CE2	2:B:1:NAG:H5	2.38	0.59
1:A:323:ARG:HH11	1:A:323:ARG:CG	2.17	0.57
1:A:119:LEU:HA	1:A:120:TRP:HB3	1.86	0.57
1:A:75:VAL:HG22	1:A:114:LEU:HD11	1.87	0.56
1:A:40:ILE:HG21	1:A:88:ILE:CD1	2.37	0.55
1:A:124:ARG:HG3	1:A:124:ARG:NH1	2.22	0.55
1:A:237:ILE:HG21	1:A:313:VAL:HG22	1.89	0.55
1:A:40:ILE:HG21	1:A:88:ILE:HD13	1.88	0.55
1:A:337:ASP:HB3	1:A:350:PRO:CD	2.37	0.55
1:A:348:THR:HG22	4:A:564:HOH:O	2.06	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:PRO:HB2	1:A:58:THR:CG2	2.37	0.54
1:A:272:ILE:HG12	2:B:2:NAG:H83	1.89	0.53
1:A:313:VAL:HG13	1:A:355:ILE:CG1	2.38	0.53
1:A:124:ARG:HG3	1:A:124:ARG:HH11	1.75	0.52
1:A:272:ILE:HG12	2:B:2:NAG:C8	2.40	0.52
1:A:119:LEU:HA	1:A:120:TRP:CB	2.39	0.52
1:A:262:ILE:N	1:A:303:GLN:HE22	2.02	0.51
1:A:93:GLN:NE2	1:A:96:ARG:NH2	2.58	0.51
1:A:331:TRP:NE1	2:B:1:NAG:H5	2.25	0.51
1:A:103:PRO:HB2	1:A:104:PRO:HD3	1.94	0.50
1:A:82:SER:N	1:A:83:GLN:HE21	2.10	0.50
1:A:304:TRP:CZ3	1:A:306:ALA:HB2	2.48	0.49
1:A:269:TRP:HE3	1:A:272:ILE:HB	1.77	0.49
1:A:313:VAL:CG1	1:A:355:ILE:HG13	2.43	0.48
1:A:273:LEU:HD12	1:A:278:ILE:HD11	1.94	0.48
1:A:339:ARG:HH11	1:A:339:ARG:HG2	1.79	0.48
1:A:320:LEU:HD21	1:A:328:ALA:HB2	1.94	0.47
1:A:93:GLN:HE21	1:A:96:ARG:HH22	1.61	0.47
1:A:304:TRP:HZ3	1:A:306:ALA:HB2	1.79	0.47
1:A:93:GLN:NE2	1:A:96:ARG:HH22	2.12	0.47
1:A:13:TYR:CD1	2:B:5:NAG:H5	2.50	0.46
1:A:344:GLY:O	1:A:345:GLN:C	2.58	0.46
2:B:4:NAG:H62	2:B:5:NAG:C8	2.46	0.46
1:A:134:LYS:HB2	1:A:134:LYS:HE2	1.56	0.46
1:A:127:ARG:NH1	4:A:517:HOH:O	2.50	0.45
1:A:10:TRP:CG	2:B:3:NAG:H3	2.52	0.45
1:A:77:GLY:O	1:A:80:TYR:HB3	2.16	0.44
1:A:12:GLN:NE2	1:A:55:LEU:HD11	2.32	0.44
1:A:237:ILE:CG2	1:A:313:VAL:HG22	2.48	0.43
1:A:275:TYR:CD2	1:A:351:LEU:HD22	2.53	0.43
1:A:344:GLY:HA3	4:A:556:HOH:O	2.17	0.43
1:A:336:ASP:OD1	1:A:337:ASP:N	2.51	0.42
1:A:234:VAL:CG1	1:A:329:MET:HB2	2.49	0.42
1:A:5:CYS:HB3	1:A:333:LEU:HG	2.01	0.42
1:A:351:LEU:HD12	1:A:351:LEU:HA	1.74	0.42
1:A:36:SER:HA	1:A:37:PHE:HA	1.78	0.42
1:A:50:TRP:CD1	1:A:50:TRP:H	2.37	0.42
1:A:10:TRP:CZ2	2:B:2:NAG:H4	2.54	0.41
1:A:347:LEU:HD23	1:A:348:THR:H	1.85	0.41
1:A:182:LEU:HD23	1:A:182:LEU:HA	1.90	0.41
1:A:269:TRP:CE3	1:A:272:ILE:HD13	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:278:ILE:HD12	1:A:305:VAL:HG11	1.99	0.41
1:A:269:TRP:CH2	2:B:2:NAG:H3	2.54	0.41
1:A:17:ASP:CA	1:A:339:ARG:HE	2.34	0.40
1:A:75:VAL:CG2	1:A:114:LEU:HD11	2.50	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	359/361 (99%)	338 (94%)	19 (5%)	2 (1%)	21	51

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	345	GLN
1	A	204	GLY

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	304/304 (100%)	252 (83%)	52 (17%)	2	7

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	9	SER
1	A	23	ASP
1	A	34	ILE
1	A	41	SER
1	A	53	VAL
1	A	69	LEU
1	A	75	VAL
1	A	83	GLN
1	A	85	PHE
1	A	88	ILE
1	A	94	SER
1	A	97	THR
1	A	106	LEU
1	A	107	ARG
1	A	120	TRP
1	A	129	LEU
1	A	134	LYS
1	A	142	ARG
1	A	158	THR
1	A	166	ARG
1	A	182	LEU
1	A	186	ASP
1	A	193	GLN
1	A	195	VAL
1	A	205	ASN
1	A	210	SER
1	A	212	ARG
1	A	225	ARG
1	A	226	LEU
1	A	233	LEU
1	A	237	ILE
1	A	239	THR
1	A	250	LYS
1	A	251	THR
1	A	252	ASP
1	A	253	VAL
1	A	269	TRP
1	A	272	ILE
1	A	282	LEU
1	A	291	ARG
1	A	313	VAL
1	A	315	ASN

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Mol	Chain	Res	Type
1	A	321	LYS
1	A	323	ARG
1	A	345	GLN
1	A	346	ASN
1	A	347	LEU
1	A	351	LEU
1	A	353	SER
1	A	359	LEU
1	A	361	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	ASN
1	A	79	ASN
1	A	83	GLN
1	A	93	GLN
1	A	128	HIS
1	A	193	GLN
1	A	231	ASN
1	A	294	GLN
1	A	303	GLN
1	A	310	GLN
1	A	345	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](#)

5 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	0.69	0	21,21,21	3.35	11 (52%)
2	NAG	B	2	2	14,14,15	0.60	0	17,19,21	1.32	2 (11%)
2	NAG	B	3	2	14,14,15	0.60	0	17,19,21	1.49	3 (17%)
2	NAG	B	4	2	14,14,15	0.56	0	17,19,21	1.88	3 (17%)
2	NAG	B	5	2	14,14,15	0.56	0	17,19,21	1.89	6 (35%)

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	NAG	B	2	2	-	0/6/23/26	0/1/1/1
2	NAG	B	3	2	-	2/6/23/26	0/1/1/1
2	NAG	B	4	2	-	5/6/23/26	0/1/1/1
2	NAG	B	5	2	-	3/6/23/26	0/1/1/1

There are no bond length outliers.

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	NAG	C1-C2-N2	-11.07	97.90	110.73
2	B	4	NAG	C2-N2-C7	5.74	130.60	122.90
2	B	1	NAG	O4-C4-C5	-5.17	96.58	109.32
2	B	1	NAG	C3-C2-N2	4.72	119.31	110.62
2	B	5	NAG	C2-N2-C7	4.70	129.20	122.90
2	B	4	NAG	C1-C2-N2	4.12	116.92	110.43
2	B	2	NAG	C1-O5-C5	3.56	116.95	112.19
2	B	3	NAG	O5-C5-C6	3.55	114.57	107.66
2	B	1	NAG	O5-C5-C4	3.46	115.93	109.70
2	B	3	NAG	C4-C3-C2	3.34	115.91	111.02
2	B	5	NAG	C1-C2-N2	3.26	115.57	110.43
2	B	1	NAG	O5-C1-C2	3.11	112.64	109.52
2	B	5	NAG	O4-C4-C3	-2.94	103.44	110.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	NAG	O5-C5-C6	2.82	113.44	106.44
2	B	1	NAG	C4-C3-C2	2.77	114.43	110.40
2	B	1	NAG	O7-C7-N2	2.41	126.24	121.98
2	B	5	NAG	O3-C3-C2	2.40	114.38	109.40
2	B	5	NAG	O3-C3-C4	-2.37	104.80	110.38
2	B	3	NAG	O4-C4-C5	-2.23	103.84	109.32
2	B	1	NAG	O7-C7-C8	-2.22	118.10	122.05
2	B	5	NAG	O4-C4-C5	2.14	114.59	109.32
2	B	1	NAG	C1-O5-C5	2.12	117.75	113.65
2	B	2	NAG	O3-C3-C4	2.11	115.36	110.38
2	B	1	NAG	C1-C2-C3	2.09	113.39	110.54
2	B	4	NAG	O5-C1-C2	-2.09	108.06	111.29

There are no chirality outliers.

All (11) torsion outliers are listed below:

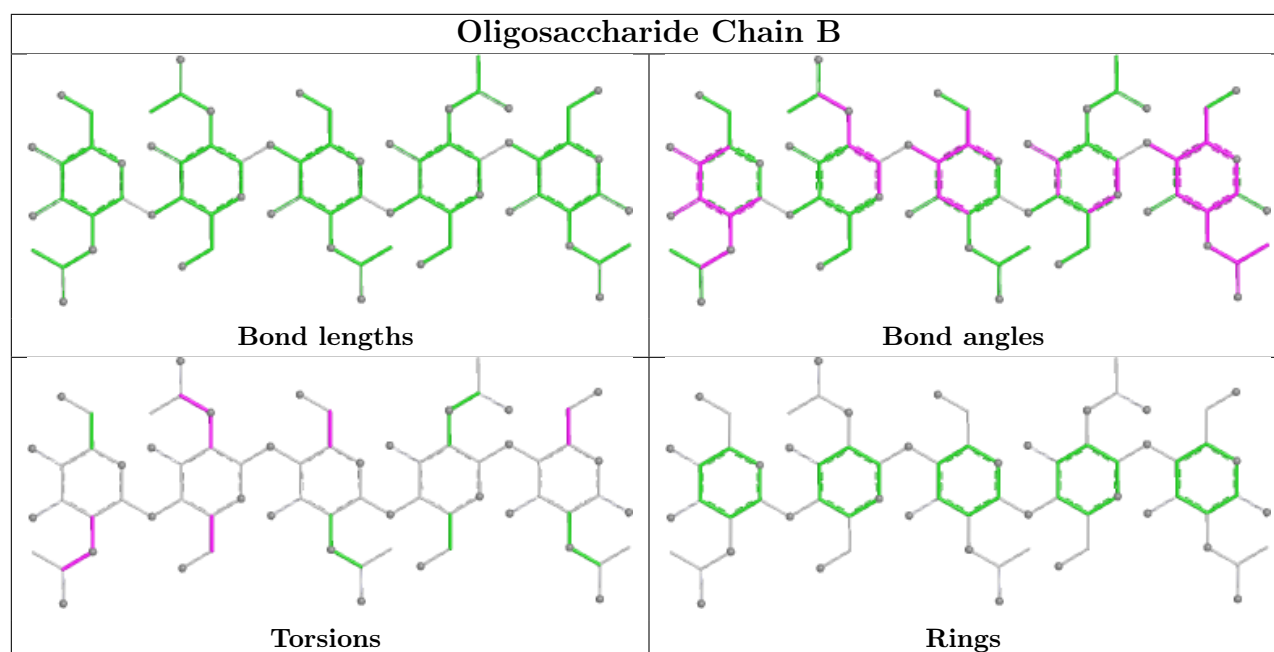
Mol	Chain	Res	Type	Atoms
2	B	4	NAG	C1-C2-N2-C7
2	B	4	NAG	C8-C7-N2-C2
2	B	4	NAG	O7-C7-N2-C2
2	B	5	NAG	C1-C2-N2-C7
2	B	5	NAG	C8-C7-N2-C2
2	B	5	NAG	O7-C7-N2-C2
2	B	3	NAG	C4-C5-C6-O6
2	B	1	NAG	C4-C5-C6-O6
2	B	4	NAG	C4-C5-C6-O6
2	B	3	NAG	O5-C5-C6-O6
2	B	4	NAG	O5-C5-C6-O6

There are no ring outliers.

5 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	3	NAG	3	0
2	B	5	NAG	4	0
2	B	1	NAG	2	0
2	B	4	NAG	3	0
2	B	2	NAG	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.



5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NAG	A	401	1	14,14,15	0.79	0	17,19,21	1.76	5 (29%)

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	NAG	A	401	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	NAG	C4-C3-C2	3.06	115.50	111.02
3	A	401	NAG	C2-N2-C7	2.99	126.90	122.90
3	A	401	NAG	O7-C7-N2	2.78	126.89	121.98
3	A	401	NAG	O7-C7-C8	-2.47	117.65	122.05
3	A	401	NAG	C3-C4-C5	-2.18	106.28	110.23

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	401	NAG	C4-C5-C6-O6
3	A	401	NAG	O5-C5-C6-O6

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	361/361 (100%)	-0.21	7 (1%) 66 57	11, 27, 60, 100	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	362	VAL	3.8
1	A	208	ALA	3.6
1	A	210	SER	3.6
1	A	78	TRP	2.8
1	A	11	SER	2.3
1	A	209	SER	2.3
1	A	81	GLY	2.2

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

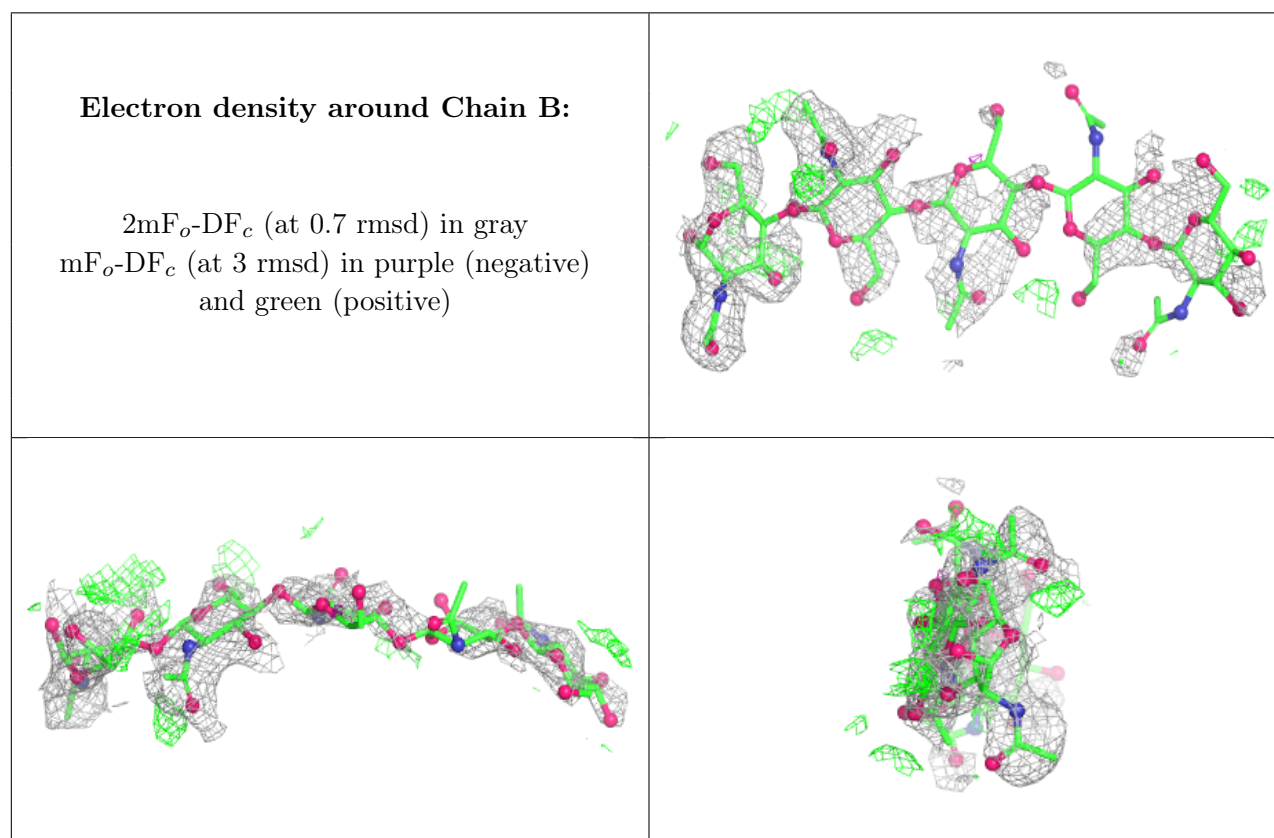
There are no non-standard protein/DNA/RNA residues in this entry.

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	NAG	B	5	14/15	0.63	0.29	45,46,50,52	14
2	NAG	B	4	14/15	0.65	0.20	38,42,47,48	14
2	NAG	B	2	14/15	0.85	0.16	33,38,42,43	14
2	NAG	B	1	15/15	0.88	0.15	21,30,35,41	15
2	NAG	B	3	14/15	0.88	0.18	40,42,45,46	14

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



6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	A	401	14/15	0.87	0.15	28,31,33,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.