



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:28 AM UTC

PDB ID : 4BPA / pdb_00004bpa
Title : Crystal structure of AmpDh2 from Pseudomonas aeruginosa in complex with NAG-NAM-NAG-NAM tetrasaccharide
Authors : Artola-Recolons, C.; Martinez-Caballero, S.; Lee, M.; Carrasco-Lopez, C.; Hesek, D.; Spink, E.; Lastochkin, E.; Zhang, W.; Hellman, L.; Boggess, B.; Mobashery, S.; Hermoso, J.A.
Deposited on : 2013-05-23
Resolution : 2.70 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

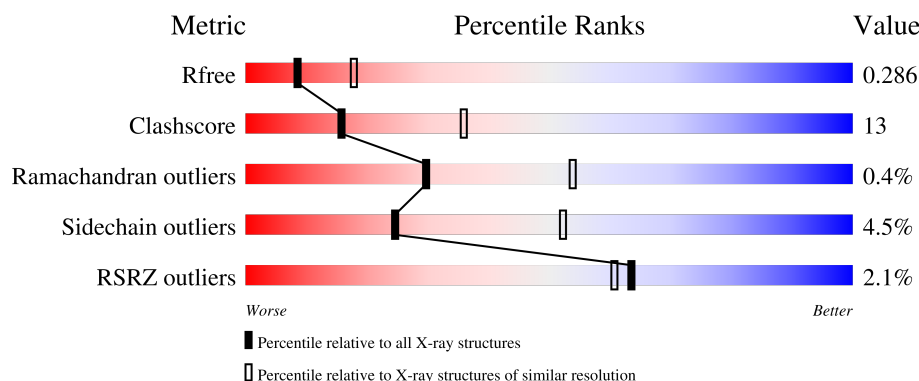
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

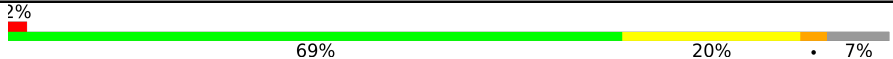
The reported resolution of this entry is 2.70 Å.

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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	259	
1	B	259	
2	C	4	

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	AMU	C	3	-	-	X	-

2 Entry composition [i](#)

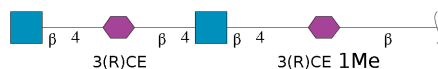
There are 4 unique types of molecules in this entry. The entry contains 3930 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 AMPDH2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	241	Total	C	N	O	S	0	0	0
			1912	1214	347	350	1			
1	B	241	Total	C	N	O	S	0	0	0
			1906	1211	344	350	1			

- Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-N-acetyl-beta-muramic acid-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-methyl 2-acetamido-3-O-[(1R)-1-carboxyethyl]-2-deoxy-beta-D-glucopyranoside.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	4	Total	C	N	O	0	0	0
			68	39	4	25			

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		
3	B	1	Total	Zn	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	23	Total	O	0	0
			23	23		
4	B	19	Total	O	0	0
			19	19		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	45.23Å 93.56Å 104.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	69.66 – 2.70 69.66 – 2.70	Depositor EDS
% Data completeness (in resolution range)	76.0 (69.66-2.70) 76.0 (69.66-2.70)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.35 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.223 , 0.290 0.226 , 0.286	Depositor DCC
R_{free} test set	623 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	42.0	Xtriage
Anisotropy	0.366	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 59.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3930	wwPDB-VP
Average B, all atoms (Å ²)	50.0	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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.53	0/1966	0.81	2/2682 (0.1%)
1	B	0.49	0/1960	0.82	1/2675 (0.0%)
All	All	0.51	0/3926	0.82	3/5357 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	44	THR	N-CA-C	-10.15	95.50	110.52
1	A	110	GLN	N-CA-C	-7.22	99.59	110.28
1	A	44	THR	N-CA-C	5.76	119.49	112.23

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	1912	0	1876	53	0
1	B	1906	0	1865	47	0
2	C	68	0	59	24	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	23	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	19	0	0	1	0
All	All	3930	0	3800	102	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 13.

All (102) 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:61:SER:OG	1:B:31:ASN:HB2	1.58	1.00
1:A:237:ARG:NH1	1:A:246:ASP:OD2	2.00	0.94
1:A:201:VAL:HG22	1:A:221:LEU:HD23	1.52	0.91
1:B:162:VAL:HG12	2:C:2:NAG:C6	2.12	0.79
1:B:25:THR:HG22	1:B:25:THR:O	1.82	0.77
1:B:119:ARG:NE	2:C:3:AMU:H82	1.99	0.77
1:A:201:VAL:CG2	1:A:221:LEU:HD23	2.16	0.75
1:A:249:THR:O	1:A:253:LEU:HD12	1.88	0.74
1:A:162:VAL:HG23	1:A:234:MET:HE1	1.71	0.72
1:B:162:VAL:HG12	2:C:2:NAG:H62	1.71	0.72
1:A:242:ASP:OD1	4:A:2021:HOH:O	2.09	0.71
1:B:160:ARG:O	2:C:2:NAG:H3	1.92	0.69
1:B:161:LYS:HE2	2:C:2:NAG:C8	2.22	0.69
1:A:187:LEU:O	1:A:190:ARG:N	2.27	0.68
2:C:1:AMV:H112	2:C:1:AMV:H2	1.76	0.67
1:A:246:ASP:HB2	1:A:249:THR:OG1	1.94	0.66
1:B:119:ARG:CZ	2:C:3:AMU:H82	2.27	0.64
1:A:101:THR:HG22	1:A:101:THR:O	1.98	0.64
1:B:101:THR:O	1:B:101:THR:HG22	1.97	0.63
1:A:81:ASN:HD21	1:B:81:ASN:CG	2.06	0.63
1:B:112:TYR:CE2	2:C:3:AMU:H81	2.33	0.62
1:A:81:ASN:ND2	1:B:81:ASN:OD1	2.28	0.62
1:B:44:THR:O	1:B:109:ASN:O	2.17	0.62
1:A:146:THR:HG22	1:A:147:PRO:HD2	1.82	0.61
1:B:162:VAL:HG12	2:C:2:NAG:H61	1.82	0.61
1:A:95:ARG:HD2	1:B:97:TRP:CD1	2.36	0.60
1:A:201:VAL:HG13	1:A:221:LEU:HD22	1.82	0.60
1:B:159:GLY:C	2:C:3:AMU:H62	2.27	0.60
1:A:207:GLN:HB2	1:A:253:LEU:HD22	1.85	0.59
1:B:223:LYS:HG3	1:B:224:ASP:N	2.19	0.57
2:C:4:NAG:H3	2:C:4:NAG:H83	1.87	0.57
1:A:187:LEU:HD21	1:A:252:LEU:HD23	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:119:ARG:CD	2:C:3:AMU:H82	2.36	0.56
1:B:153:HIS:CE1	1:B:161:LYS:HE3	2.42	0.55
1:A:69:ASP:OD1	1:A:72:ALA:HA	2.07	0.55
1:A:186:GLU:O	1:A:187:LEU:O	2.25	0.55
1:B:161:LYS:NZ	2:C:1:AMV:H111	2.23	0.54
1:A:186:GLU:O	1:A:187:LEU:C	2.47	0.54
1:B:121:TRP:CH2	2:C:3:AMU:H83	2.42	0.54
1:A:187:LEU:O	1:A:189:ARG:N	2.42	0.53
1:B:82:ARG:NH1	4:B:2006:HOH:O	2.41	0.53
1:A:179:VAL:HG13	1:A:180:PRO:HD2	1.90	0.53
1:A:162:VAL:CG2	1:A:234:MET:HE1	2.36	0.53
1:A:246:ASP:CB	1:A:249:THR:OG1	2.57	0.53
1:B:119:ARG:HD2	2:C:3:AMU:H82	1.92	0.52
1:A:48:LEU:HB3	1:A:49:PRO:HD3	1.91	0.52
1:B:25:THR:O	1:B:25:THR:CG2	2.54	0.51
1:A:70:GLU:HG3	1:A:127:ALA:CB	2.41	0.51
1:A:246:ASP:HB3	1:A:249:THR:H	1.74	0.51
1:A:21:PRO:O	1:A:22:ARG:HG3	2.10	0.51
2:C:4:NAG:H3	2:C:4:NAG:C8	2.41	0.51
1:B:161:LYS:HZ3	2:C:1:AMV:H6C2	1.76	0.51
1:A:146:THR:CG2	1:A:147:PRO:HD2	2.41	0.50
2:C:3:AMU:H111	2:C:4:NAG:O5	2.11	0.50
1:B:225:THR:O	1:B:229:ILE:HG22	2.12	0.49
1:B:48:LEU:HB3	1:B:49:PRO:HD3	1.92	0.49
1:A:225:THR:O	1:A:229:ILE:HG22	2.12	0.49
1:A:79:ASP:C	1:A:81:ASN:H	2.20	0.49
1:B:218:THR:HB	1:B:220:GLU:H	1.78	0.49
1:B:229:ILE:O	1:B:233:GLN:HG3	2.13	0.49
1:A:211:HIS:CE1	1:A:256:VAL:HG13	2.48	0.48
2:C:4:NAG:H83	2:C:4:NAG:C3	2.43	0.48
1:B:190:ARG:NH1	1:B:193:GLU:OE1	2.46	0.48
1:A:83:ARG:HB3	1:B:33:ASP:HB3	1.95	0.48
1:A:110:GLN:HB3	1:A:122:TYR:CD1	2.49	0.48
2:C:1:AMV:H2	2:C:1:AMV:C11	2.44	0.47
1:A:187:LEU:HD21	1:A:252:LEU:CD2	2.45	0.47
1:B:153:HIS:CE1	1:B:161:LYS:CE	2.97	0.47
1:B:41:LEU:HD13	1:B:168:PHE:CZ	2.51	0.46
1:B:82:ARG:HH11	1:B:82:ARG:HG2	1.81	0.46
1:A:211:HIS:CE1	1:A:256:VAL:CG1	3.00	0.45
1:B:45:SER:H	2:C:1:AMV:H8C3	1.81	0.45
1:B:80:GLU:HB3	1:B:103:ILE:HD11	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:112:TYR:HB3	1:B:121:TRP:CE3	2.53	0.44
1:A:119:ARG:NH2	4:A:2017:HOH:O	2.49	0.44
1:B:110:GLN:HB3	1:B:122:TYR:CD1	2.52	0.44
1:A:70:GLU:HG3	1:A:127:ALA:HB2	2.00	0.44
1:B:112:TYR:CD1	1:B:119:ARG:HD2	2.53	0.43
1:B:112:TYR:CZ	2:C:3:AMU:H81	2.52	0.43
1:A:203:TRP:CZ2	1:A:254:LEU:HD11	2.53	0.43
2:C:4:NAG:C8	2:C:4:NAG:C3	2.96	0.43
1:A:171:LYS:HG2	1:A:181:TRP:CZ2	2.53	0.43
1:A:80:GLU:HB3	1:A:103:ILE:HD11	1.99	0.43
1:A:41:LEU:HD13	1:A:168:PHE:CZ	2.54	0.43
1:A:235:LYS:HD3	1:A:236:TYR:CE2	2.54	0.43
1:A:36:VAL:HG12	1:A:145:ILE:HD11	2.01	0.42
1:A:237:ARG:NH2	1:A:240:ARG:HG3	2.34	0.42
1:B:36:VAL:HG12	1:B:145:ILE:HD11	2.01	0.42
1:A:187:LEU:O	1:A:188:ALA:C	2.62	0.42
1:A:148:ASP:HB3	1:A:259:SER:HA	2.00	0.42
1:B:79:ASP:C	1:B:81:ASN:H	2.28	0.42
1:B:114:ASP:OD1	1:B:119:ARG:HD3	2.19	0.42
1:B:161:LYS:HA	2:C:2:NAG:H5	2.01	0.42
1:B:171:LYS:HD3	1:B:181:TRP:CE2	2.55	0.42
1:B:258:THR:HG22	1:B:258:THR:O	2.20	0.42
1:A:201:VAL:HG22	1:A:221:LEU:CD2	2.38	0.41
1:A:221:LEU:HD12	1:A:226:ARG:HH11	1.86	0.41
1:B:237:ARG:HB2	1:B:252:LEU:HD11	2.03	0.41
1:A:242:ASP:OD1	1:A:242:ASP:N	2.45	0.40
1:A:126:GLU:CD	1:A:172:ARG:HH22	2.29	0.40
1:A:236:TYR:CZ	1:A:256:VAL:HG21	2.55	0.40
1:A:101:THR:O	1:A:101:THR:CG2	2.67	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	239/259 (92%)	226 (95%)	11 (5%)	2 (1%)	16	37
1	B	239/259 (92%)	231 (97%)	8 (3%)	0	100	100
All	All	478/518 (92%)	457 (96%)	19 (4%)	2 (0%)	30	54

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	187	LEU
1	A	21	PRO

5.3.2 Protein sidechains ⓘ

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	200/215 (93%)	192 (96%)	8 (4%)	28	56
1	B	199/215 (93%)	189 (95%)	10 (5%)	22	48
All	All	399/430 (93%)	381 (96%)	18 (4%)	24	52

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

Mol	Chain	Res	Type
1	A	74	VAL
1	A	162	VAL
1	A	221	LEU
1	A	242	ASP
1	A	244	GLU
1	A	246	ASP
1	A	253	LEU
1	A	254	LEU
1	B	52	LEU
1	B	81	ASN
1	B	108	VAL
1	B	117	GLN
1	B	132	LEU
1	B	146	THR

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Mol	Chain	Res	Type
1	B	162	VAL
1	B	183	LYS
1	B	218	THR
1	B	234	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	HIS
1	A	81	ASN
1	A	211	HIS
1	B	57	HIS

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 ⓘ

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	AMV	C	1	2	21,21,21	1.86	6 (28%)	25,29,29	1.47	6 (24%)
2	NAG	C	2	2	14,14,15	2.43	4 (28%)	17,19,21	2.72	5 (29%)
2	AMU	C	3	2	19,19,20	2.14	5 (26%)	23,26,28	3.37	9 (39%)
2	NAG	C	4	2	14,14,15	1.07	0	17,19,21	2.50	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
2	AMV	C	1	2	-	6/16/36/36	0/1/1/1
2	NAG	C	2	2	-	5/6/23/26	0/1/1/1
2	AMU	C	3	2	-	5/14/31/34	0/1/1/1
2	NAG	C	4	2	-	4/6/23/26	0/1/1/1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	2	NAG	C1-C2	7.35	1.62	1.52
2	C	3	AMU	O5-C5	4.93	1.53	1.43
2	C	3	AMU	C3-C2	4.74	1.58	1.52
2	C	1	AMV	C3-C2	3.74	1.59	1.53
2	C	1	AMV	O1-C1	3.67	1.46	1.40
2	C	2	NAG	C3-C2	3.60	1.60	1.52
2	C	3	AMU	O5-C1	3.31	1.49	1.43
2	C	1	AMV	C4-C3	2.99	1.60	1.52
2	C	2	NAG	O4-C4	2.79	1.49	1.43
2	C	3	AMU	O4-C4	2.79	1.49	1.43
2	C	1	AMV	O4-C4	2.72	1.49	1.43
2	C	1	AMV	O3-C3	2.65	1.50	1.43
2	C	3	AMU	C4-C5	2.30	1.57	1.53
2	C	2	NAG	C2-N2	2.26	1.50	1.46
2	C	1	AMV	C2-N2	2.18	1.49	1.45

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	3	AMU	C1-O5-C5	8.64	123.77	112.19
2	C	3	AMU	C4-C3-C2	-7.58	97.96	111.18
2	C	2	NAG	C2-N2-C7	7.04	132.34	122.90
2	C	2	NAG	C1-C2-N2	6.78	121.12	110.43
2	C	3	AMU	C3-C2-N2	6.40	121.01	110.57
2	C	4	NAG	O5-C1-C2	-5.86	102.22	111.29
2	C	4	NAG	C2-N2-C7	5.06	129.68	122.90
2	C	3	AMU	O3-C3-C2	4.78	120.06	108.85
2	C	4	NAG	C1-C2-N2	4.40	117.37	110.43
2	C	1	AMV	C1-O5-C5	3.33	120.22	113.72
2	C	3	AMU	O3-C3-C4	3.21	115.38	107.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	3	AMU	O4-C4-C3	3.02	117.67	109.94
2	C	3	AMU	O5-C1-C2	-2.94	106.74	111.29
2	C	2	NAG	C4-C3-C2	2.94	115.32	111.02
2	C	4	NAG	C8-C7-N2	2.74	120.65	116.12
2	C	3	AMU	C3-C4-C5	-2.69	103.83	109.57
2	C	2	NAG	C8-C7-N2	2.68	120.56	116.12
2	C	1	AMV	O3-C9-C11	2.57	114.71	107.54
2	C	1	AMV	O5-C5-C6	2.57	112.80	106.44
2	C	2	NAG	O7-C7-C8	-2.42	117.75	122.05
2	C	3	AMU	O5-C5-C4	2.39	116.64	110.83
2	C	4	NAG	O7-C7-C8	-2.33	117.90	122.05
2	C	1	AMV	C3-C2-N2	2.18	114.38	110.91
2	C	1	AMV	O6-C6-C5	-2.17	103.94	111.33
2	C	1	AMV	O4-C4-C3	2.04	115.18	109.94

There are no chirality outliers.

All (20) torsion outliers are listed below:

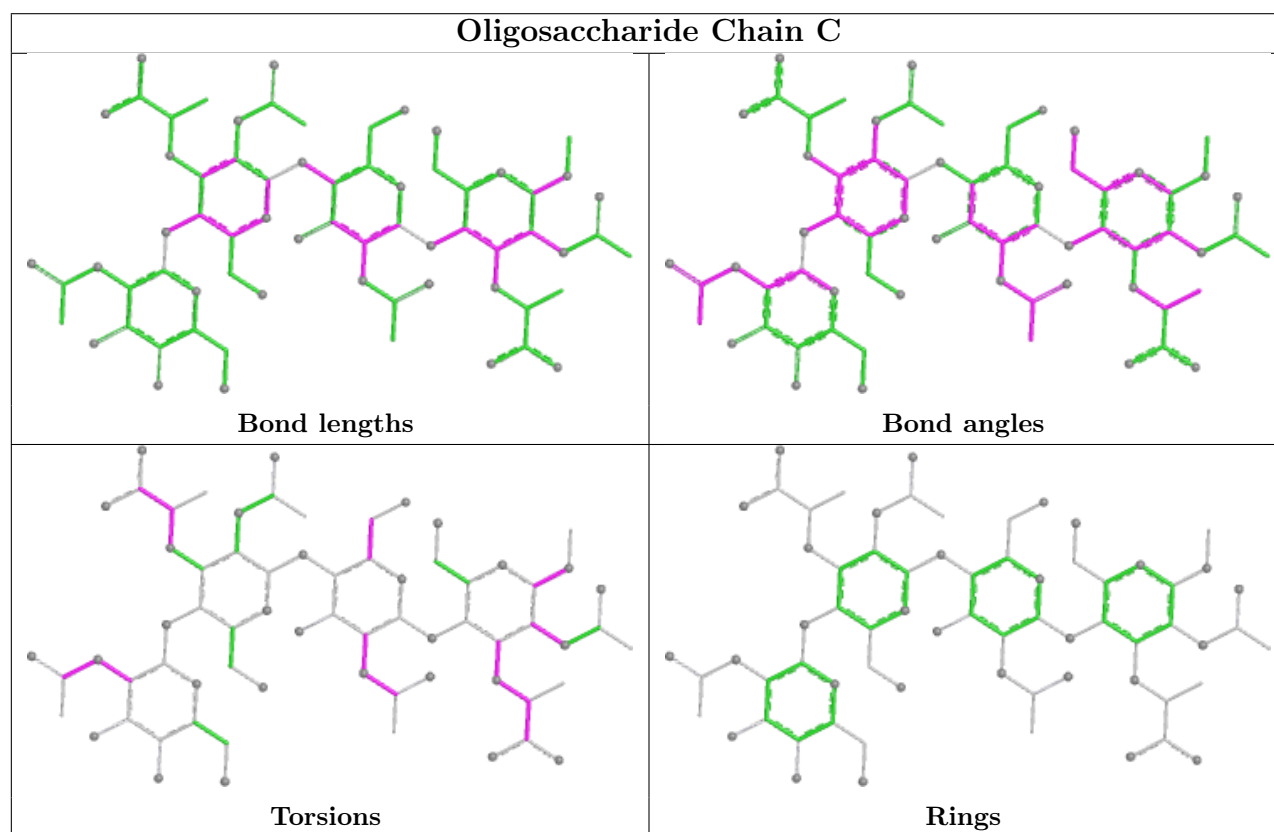
Mol	Chain	Res	Type	Atoms
2	C	1	AMV	C2-C3-O3-C9
2	C	1	AMV	O11-C10-C9-C11
2	C	1	AMV	C11-C9-O3-C3
2	C	2	NAG	C1-C2-N2-C7
2	C	4	NAG	C3-C2-N2-C7
2	C	2	NAG	O5-C5-C6-O6
2	C	2	NAG	C4-C5-C6-O6
2	C	2	NAG	C8-C7-N2-C2
2	C	2	NAG	O7-C7-N2-C2
2	C	4	NAG	C8-C7-N2-C2
2	C	4	NAG	O7-C7-N2-C2
2	C	1	AMV	OXT-C10-C9-C11
2	C	3	AMU	C11-C9-O3-C3
2	C	1	AMV	C3-C2-N2-C7
2	C	3	AMU	O11-C10-C9-C11
2	C	4	NAG	C1-C2-N2-C7
2	C	3	AMU	O10-C10-C9-C11
2	C	3	AMU	O11-C10-C9-O3
2	C	1	AMV	O5-C1-O1-C12
2	C	3	AMU	C10-C9-O3-C3

There are no ring outliers.

4 monomers are involved in 24 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	AMV	5	0
2	C	3	AMU	9	0
2	C	4	NAG	5	0
2	C	2	NAG	6	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](#)

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	241/259 (93%)	0.53	6 (2%) 58 55	33, 53, 88, 107	0
1	B	241/259 (93%)	0.16	4 (1%) 69 67	31, 40, 62, 91	0
All	All	482/518 (93%)	0.35	10 (2%) 63 61	31, 44, 82, 107	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	19	SER	2.8
1	A	243	GLY	2.4
1	A	219	GLY	2.3
1	B	198	LEU	2.3
1	B	21	PRO	2.3
1	A	199	PRO	2.2
1	A	201	VAL	2.2
1	A	194	LEU	2.1
1	B	187	LEU	2.1
1	A	200	ASP	2.0

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

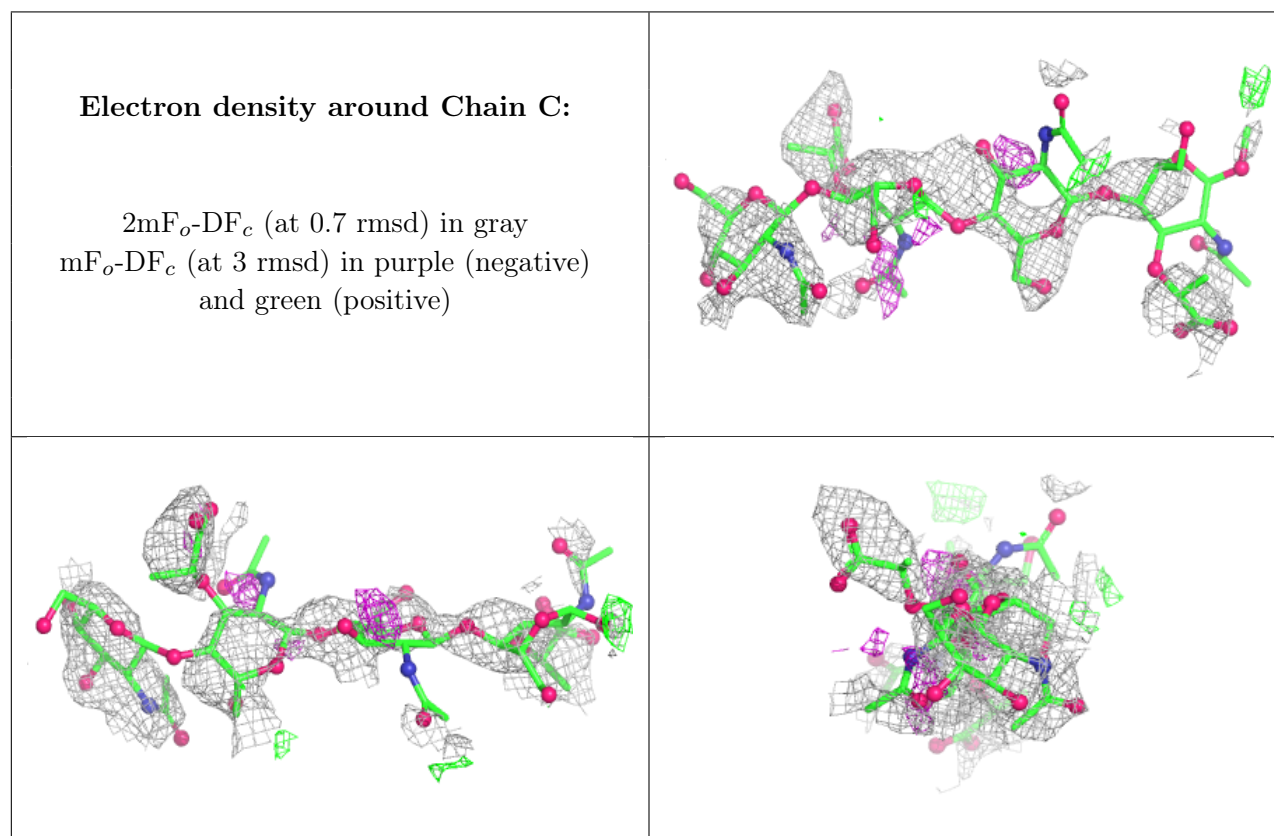
There are no non-standard protein/DNA/RNA residues in this entry.

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	AMV	C	1	21/21	0.61	0.20	100,111,119,125	0
2	NAG	C	2	14/15	0.65	0.20	93,96,105,108	0
2	AMU	C	3	19/20	0.71	0.19	94,98,107,108	0
2	NAG	C	4	14/15	0.79	0.15	100,108,120,120	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 ⓘ

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	ZN	A	1260	1/1	0.89	0.09	102,102,102,102	0
3	ZN	B	1260	1/1	0.99	0.04	50,50,50,50	0

6.5 Other polymers [i](#)

There are no such residues in this entry.