wwPDB X-ray Structure Validation Summary Report (i)

Nov 2, 2026 – 03:23 pm GMT

PDB ID	:	28NP / pdb_000028np
Title	:	Crystal structure of the Norovirus GII.4 P domain in complex with 3 sialyl fucosyllactose
Authors	:	Pisapia, L.; Varrot, A.
Deposited on	:	2026-02-10
Resolution	:	1.95 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

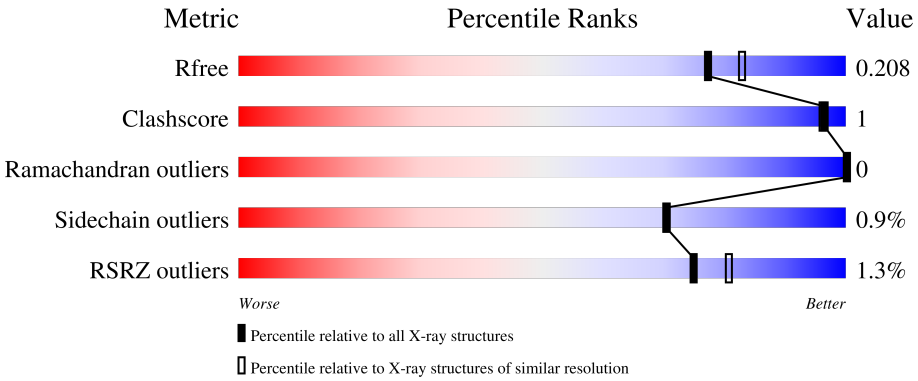
MolProbity	:	4-5-2 with Phenix2.2
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.2
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

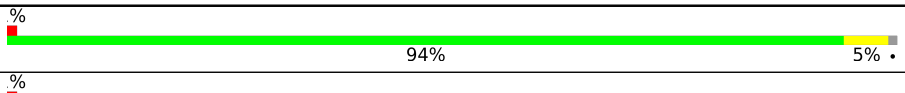
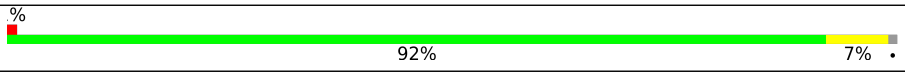
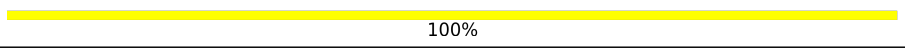
The reported resolution of this entry is 1.95 Å.

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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	310	
1	B	310	
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
3	EDO	B	602	-	X	-	-

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 5551 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.

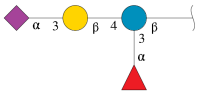
- Molecule 1 is a protein called VP1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	308	Total	C	N	O	S	0	6	0
			2445	1549	416	468	12			
1	B	308	Total	C	N	O	S	0	1	0
			2401	1524	406	460	11			

There are 8 discrepancies between the modelled and reference sequences:

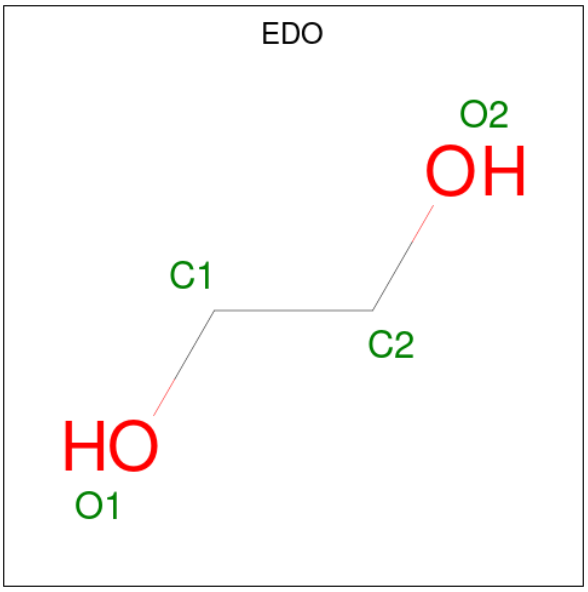
Chain	Residue	Modelled	Actual	Comment	Reference
A	221	GLY	-	expression tag	UNP A0A4D6BSS1
A	222	ALA	-	expression tag	UNP A0A4D6BSS1
A	223	MET	-	expression tag	UNP A0A4D6BSS1
A	224	ALA	-	expression tag	UNP A0A4D6BSS1
B	221	GLY	-	expression tag	UNP A0A4D6BSS1
B	222	ALA	-	expression tag	UNP A0A4D6BSS1
B	223	MET	-	expression tag	UNP A0A4D6BSS1
B	224	ALA	-	expression tag	UNP A0A4D6BSS1

- Molecule 2 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galacto pyranose-(1-4)-[alpha-L-fucopyranose-(1-3)]beta-D-glucopyranose.



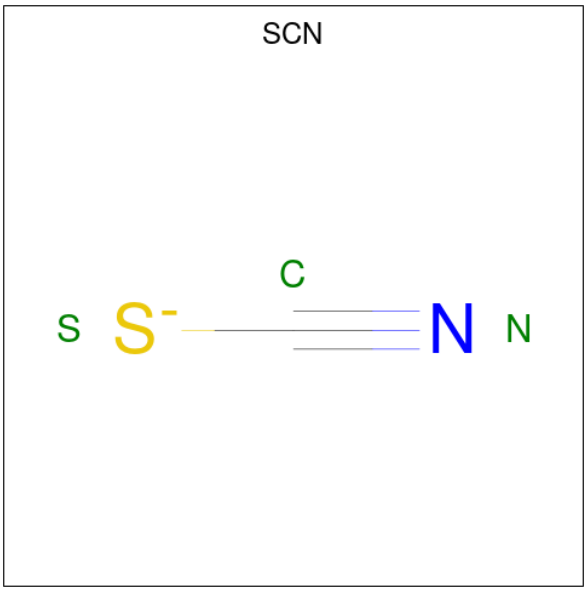
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	4	Total	C	N	O	0	0	0
			53	29	1	23			

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is THIOCYANATE ION (CCD ID: SCN) (formula: CNS).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N S 3 1 1 1	0	0
4	A	1	Total C N S 3 1 1 1	0	0
4	B	1	Total C N S 3 1 1 1	0	0

- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Na 1 1	0	0

- Molecule 6 is water.

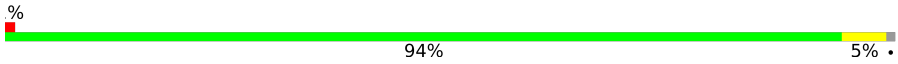
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	306	Total O 308 308	0	2
6	B	310	Total O 310 310	0	0

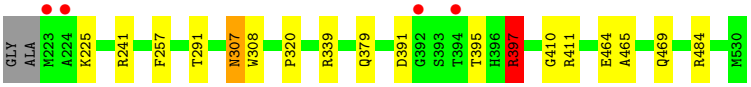
3 Residue-property plots (i)

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: VP1

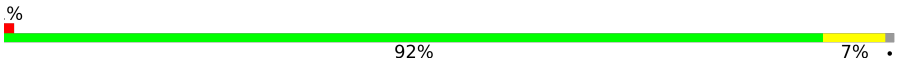
Chain A:

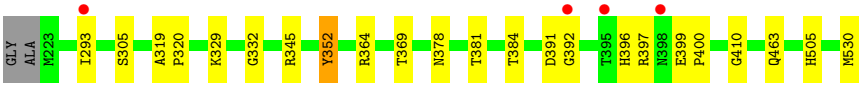




• Molecule 1: VP1

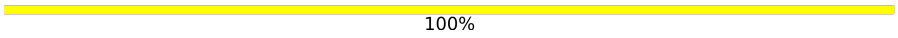
Chain B:





• Molecule 2: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-[alpha-L-fucopyranose-(1-3)]beta-D-glucopyranose

Chain C:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α, β, γ	86.92Å 100.23Å 130.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.28 – 1.95 46.28 – 1.95	Depositor EDS
% Data completeness (in resolution range)	100.0 (46.28-1.95) 100.0 (46.28-1.95)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.89 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.8.0430	Depositor
R, R_{free}	0.144 , 0.200 0.157 , 0.208	Depositor DCC
R_{free} test set	2079 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å^2)	23.6	Xtriage
Anisotropy	0.160	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 54.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	5551	wwPDB-VP
Average B, all atoms (Å^2)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.17% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NA, EDO, SCN, FUC, BGC, GAL, SIA

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	1.08	1/2526 (0.0%)	1.33	11/3453 (0.3%)
1	B	1.07	1/2472 (0.0%)	1.32	7/3381 (0.2%)
All	All	1.07	2/4998 (0.0%)	1.32	18/6834 (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	3
1	B	0	2
All	All	0	5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	411	ARG	NE-CZ	7.05	1.40	1.33
1	B	505	HIS	CE1-NE2	5.34	1.37	1.32

The worst 5 of 18 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	391	ASP	CA-CB-CG	6.51	119.11	112.60
1	A	397	ARG	CG-CD-NE	6.24	125.73	112.00
1	B	378	ASN	CA-CB-CG	-6.08	106.52	112.60
1	B	463	GLN	CB-CA-C	-5.98	100.68	110.85
1	A	257	PHE	CA-CB-CG	5.78	119.58	113.80

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	241	ARG	Sidechain
1	A	339[A]	ARG	Sidechain
1	A	397	ARG	Sidechain
1	B	345	ARG	Sidechain
1	B	364	ARG	Sidechain

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	2445	0	2333	3	0
1	B	2401	0	2278	7	1
2	C	53	0	45	0	0
3	A	8	0	8	1	0
3	B	16	0	21	1	0
4	A	6	0	0	0	0
4	B	3	0	0	1	0
5	A	1	0	0	0	0
6	A	308	0	0	0	0
6	B	310	0	0	0	0
All	All	5551	0	4685	10	1

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 1.

The worst 5 of 10 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:329:LYS:HG2	1:B:352:TYR:CD2	2.36	0.60
1:B:399:GLU:HA	1:B:400:PRO:C	2.31	0.54
1:B:391:ASP:C	1:B:391:ASP:OD1	2.51	0.53
1:B:329:LYS:HE3	3:B:602:EDO:H22	1.93	0.50
1:B:305:SER:HB3	4:B:601:SCN:S	2.54	0.47

All (1) 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:B:392:GLY:O	1:B:396:HIS:CE1[4_565]	2.00	0.20

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	312/310 (101%)	303 (97%)	9 (3%)	0	100	100
1	B	307/310 (99%)	297 (97%)	10 (3%)	0	100	100
All	All	619/620 (100%)	600 (97%)	19 (3%)	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	273/268 (102%)	271 (99%)	2 (1%)	76	76
1	B	264/268 (98%)	261 (99%)	3 (1%)	65	64
All	All	537/536 (100%)	532 (99%)	5 (1%)	70	70

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	225	LYS
1	A	397	ARG
1	B	293	ILE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	352	TYR
1	B	530	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	396	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

4 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	BGC	C	1	2	12,12,12	0.92	1 (8%)	17,17,17	1.22	3 (17%)
2	GAL	C	2	2	11,11,12	1.19	1 (9%)	15,15,17	1.17	1 (6%)
2	SIA	C	3	2	20,20,21	1.97	2 (10%)	24,28,31	0.97	2 (8%)
2	FUC	C	4	2	10,10,11	1.32	1 (10%)	14,14,16	1.44	1 (7%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BGC	C	1	2	-	1/2/22/22	0/1/1/1
2	GAL	C	2	2	-	0/2/19/22	0/1/1/1
2	SIA	C	3	2	-	4/18/34/38	0/1/1/1
2	FUC	C	4	2	-	-	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	3	SIA	C2-C1	7.39	1.59	1.52
2	C	3	SIA	O1B-C1	-3.25	1.19	1.30
2	C	4	FUC	C2-C3	-2.97	1.48	1.52
2	C	2	GAL	O5-C5	2.24	1.48	1.43
2	C	1	BGC	O5-C5	-2.11	1.39	1.44

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	4	FUC	C1-C2-C3	4.04	114.64	109.67
2	C	2	GAL	C1-C2-C3	2.93	113.27	109.67
2	C	1	BGC	C1-C2-C3	-2.57	104.97	110.31
2	C	1	BGC	O5-C5-C4	-2.40	105.34	109.69
2	C	1	BGC	O2-C2-C3	2.40	115.89	110.35

There are no chirality outliers.

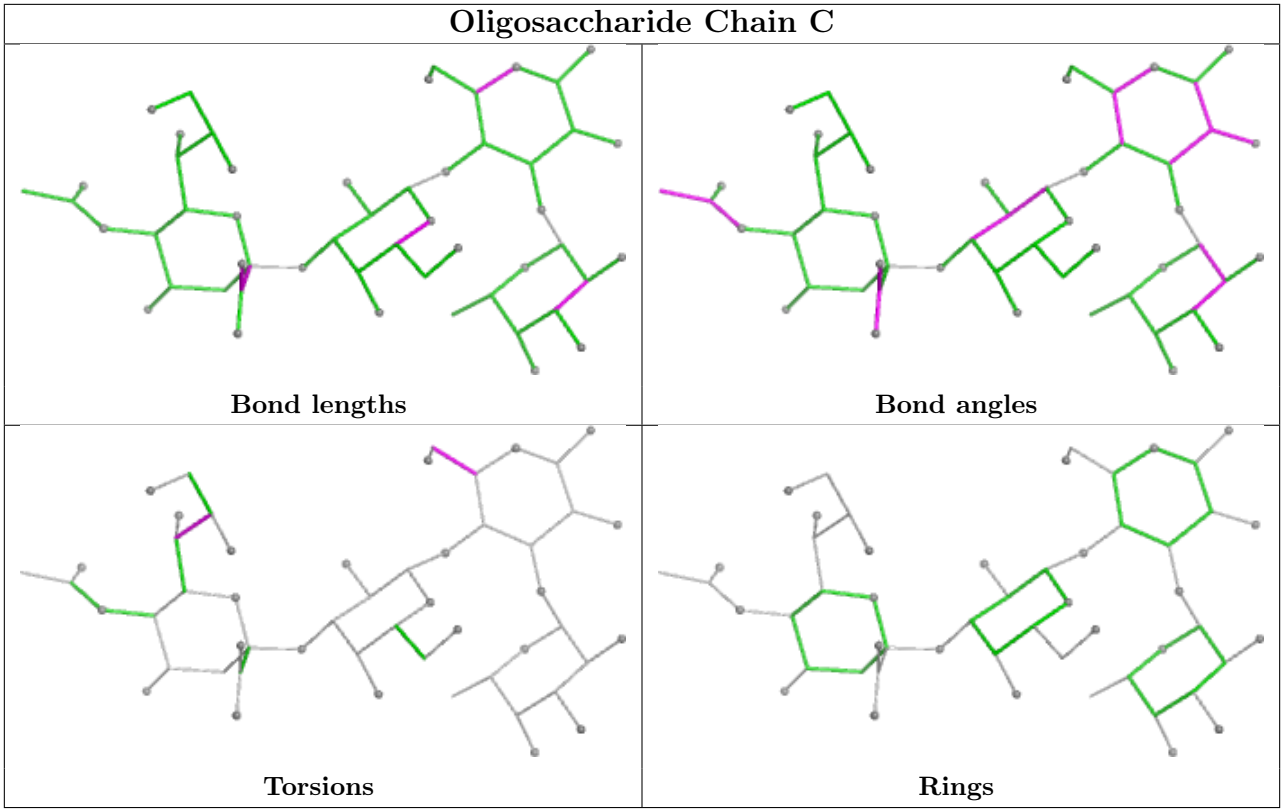
All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	3	SIA	C6-C7-C8-C9
2	C	3	SIA	O7-C7-C8-C9
2	C	3	SIA	O7-C7-C8-O8
2	C	3	SIA	C6-C7-C8-O8
2	C	1	BGC	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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 10 ligands modelled in this entry, 1 is monoatomic - leaving 9 for Mogul analysis.

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	$\#\|Z\| > 2$
4	SCN	B	601	-	1,2,2	4.39	1 (100%)	0,1,1	-	-
4	SCN	A	603	-	1,2,2	2.06	1 (100%)	0,1,1	-	-
3	EDO	B	604	-	3,3,3	0.52	0	2,2,2	0.19	0
3	EDO	B	605	-	3,3,3	0.72	0	2,2,2	0.31	0
3	EDO	B	603	-	3,3,3	0.34	0	2,2,2	0.31	0
3	EDO	A	602	-	3,3,3	2.07	1 (33%)	2,2,2	1.00	0
3	EDO	A	601	-	3,3,3	2.30	1 (33%)	2,2,2	0.59	0
3	EDO	B	602	-	3,3,3	2.18	2 (66%)	2,2,2	0.87	0
4	SCN	A	604	-	1,2,2	0.18	0	0,1,1	-	-

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	EDO	B	604	-	-	1/1/1/1	-
3	EDO	B	605	-	-	1/1/1/1	-
3	EDO	B	603	-	-	1/1/1/1	-
3	EDO	A	602	-	-	1/1/1/1	-
3	EDO	A	601	-	-	1/1/1/1	-
3	EDO	B	602	-	-	1/1/1/1	-

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	601	SCN	C-N	-4.39	1.00	1.15
3	A	601	EDO	O1-C1	-3.42	1.24	1.42
3	A	602	EDO	O2-C2	-3.02	1.26	1.42
3	B	602	EDO	O2-C2	-2.83	1.27	1.42
3	B	602	EDO	O1-C1	-2.47	1.29	1.42

There are no bond angle outliers.

There are no chirality outliers.

5 of 6 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	602	EDO	O1-C1-C2-O2
3	B	605	EDO	O1-C1-C2-O2
3	A	601	EDO	O1-C1-C2-O2
3	B	603	EDO	O1-C1-C2-O2
3	B	602	EDO	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	601	SCN	1	0
3	A	602	EDO	1	0
3	B	602	EDO	1	0

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

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	308/310 (99%)	-0.40	4 (1%)	75	81	11, 21, 42, 69	6 (1%)
1	B	308/310 (99%)	-0.39	4 (1%)	75	81	12, 22, 44, 77	1 (0%)
All	All	616/620 (99%)	-0.40	8 (1%)	75	81	11, 22, 44, 77	7 (1%)

The worst 5 of 8 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	293	ILE	3.0
1	A	223	MET	3.0
1	B	395	THR	2.8
1	A	392	GLY	2.4
1	A	224	ALA	2.4

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates

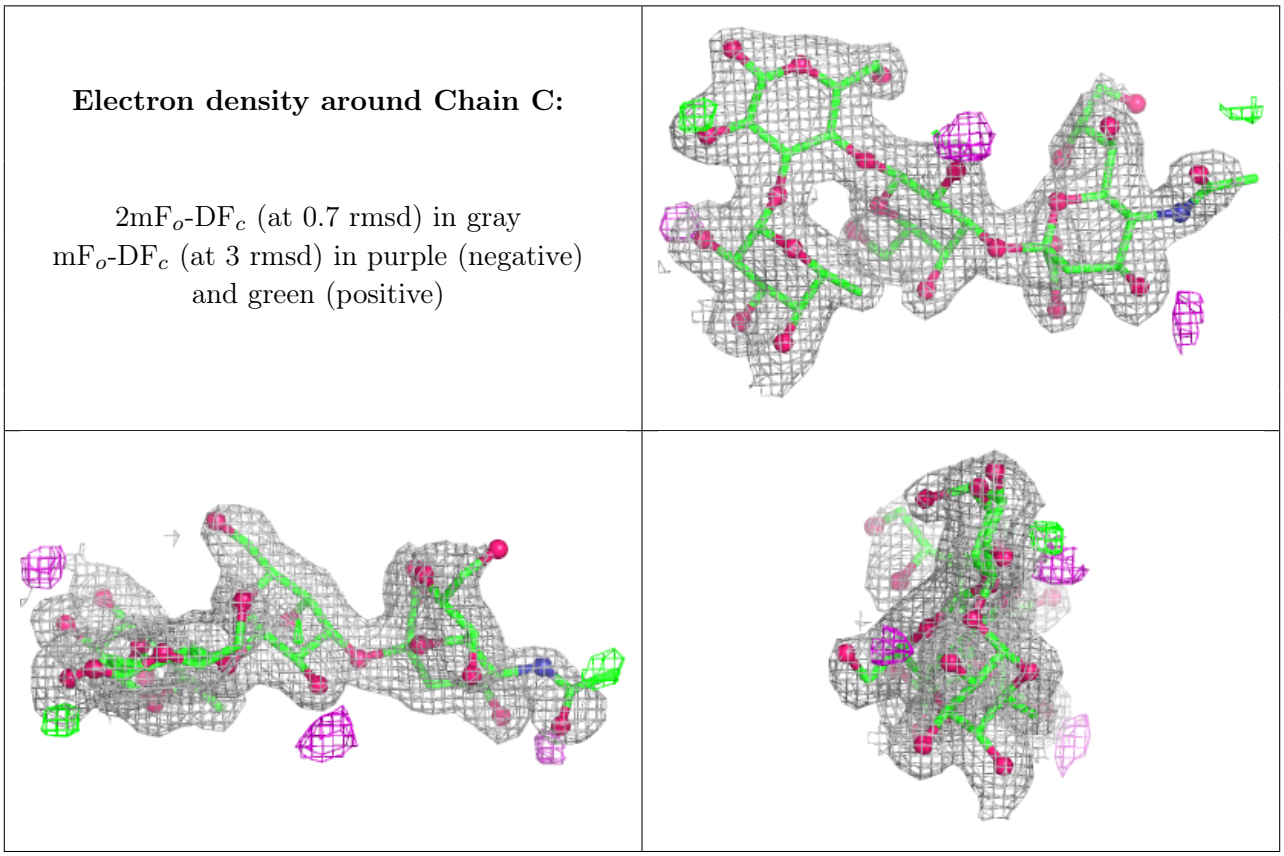
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	SIA	C	3	20/21	0.87	0.10	37,46,66,70	0
2	GAL	C	2	11/12	0.88	0.09	28,33,44,45	0
2	BGC	C	1	12/12	0.94	0.07	23,30,42,44	0
2	FUC	C	4	10/11	0.97	0.05	18,20,23,28	0

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

Electron density around Chain C:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



6.4 Ligands

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	EDO	B	604	4/4	0.80	0.15	47,50,53,64	0
3	EDO	B	605	4/4	0.82	0.17	39,48,52,54	0
3	EDO	B	603	4/4	0.86	0.10	43,50,54,59	0
3	EDO	B	602	4/4	0.87	0.12	28,31,39,43	0
3	EDO	A	602	4/4	0.88	0.14	35,37,41,43	0
3	EDO	A	601	4/4	0.91	0.10	20,32,35,36	0
4	SCN	A	604	3/3	0.92	0.15	41,41,47,61	0
4	SCN	B	601	3/3	0.95	0.10	13,13,23,51	0
4	SCN	A	603	3/3	0.96	0.11	24,24,35,42	0
5	NA	A	605	1/1	0.99	0.02	21,21,21,21	0

6.5 Other polymers (i)

There are no such residues in this entry.