

wwPDB X-ray Structure Validation Summary Report (i)

Sep 24, 2023 – 08:59 AM EDT

PDB ID : 5U74

Title: Structure of human Niemann-Pick C1 protein

Authors : Li, X.

Deposited on : 2016-12-11

Resolution : 3.33 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 : 4.02b-467

Mogul: 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.35.1

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove) roteins) : Engh & Huber (2001)

Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

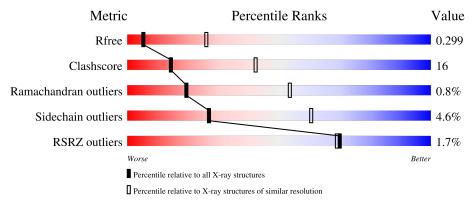
Validation Pipeline (wwPDB-VP) : 2.35.1

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 3.33 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries,\ resolution\ range(\mathring{A})}) \end{array}$
R_{free}	130704	1060 (3.38-3.30)
Clashscore	141614	1111 (3.38-3.30)
Ramachandran outliers	138981	1090 (3.38-3.30)
Sidechain outliers	138945	1089 (3.38-3.30)
RSRZ outliers	127900	1028 (3.38-3.30)

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	1278	45%	24%	• 30%			
2	В	2	50%		50%			
3	С	3	33%	67 ⁻	%			
4	D	3	67%		33%			

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
4	NAG	D	1	-	-	-	X
5	NAG	A	1301	-	-	=	X



2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 7311 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 Niemann-Pick C1 protein.

Mol	Chain	Residues		A	toms			ZeroOcc	AltConf	Trace
1	A	900	Total 7081	C 4604	N 1144	O 1287	S 46	0	3	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	642	ILE	MET	conflict	UNP O15118

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



Mo	Chain	Residues	1	Ator	ns	ZeroOcc	AltConf	Trace
2	В	2	Total 28		N 2	0	0	0

• Molecule 3 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	A	Aton	ns		ZeroOcc	AltConf	Trace
3	С	3	Total 39	C 22	N 2	O 15	0	0	0

• Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-

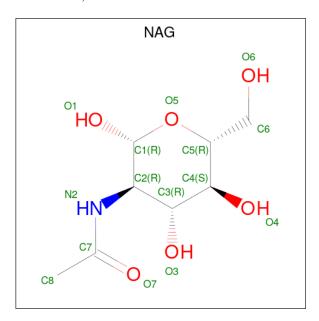


beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	A	Aton	ns		ZeroOcc	AltConf	Trace
4	D	3	10001	С	N	0	0	0	0
			39	22	2	15			

 \bullet Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $\rm C_8H_{15}NO_6).$



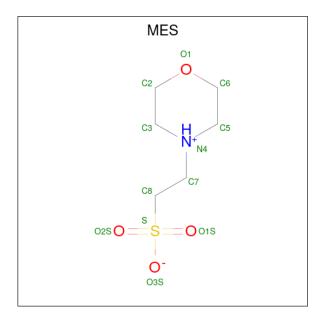
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C N O 14 8 1 5	0	0
5	A	1	Total C N O	0	0
F	Δ.	1	14 8 1 5 Total C N O	0	0
5	A	1	14 8 1 5	0	0
5	A	1	Total C N O 14 8 1 5	0	0
5	A	1	Total C N O 14 8 1 5	0	0
5	A	1	Total C N O 14 8 1 5	0	0
5	A	1	Total C N O 14 8 1 5	0	0



 $Continued\ from\ previous\ page...$

Mol	Chain	Residues	A	ton	$\mathbf{n}\mathbf{s}$		ZeroOcc	AltConf
5	A	1	Total 14	C 8	N 1	O 5	0	0

• Molecule 6 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: $C_6H_{13}NO_4S$).



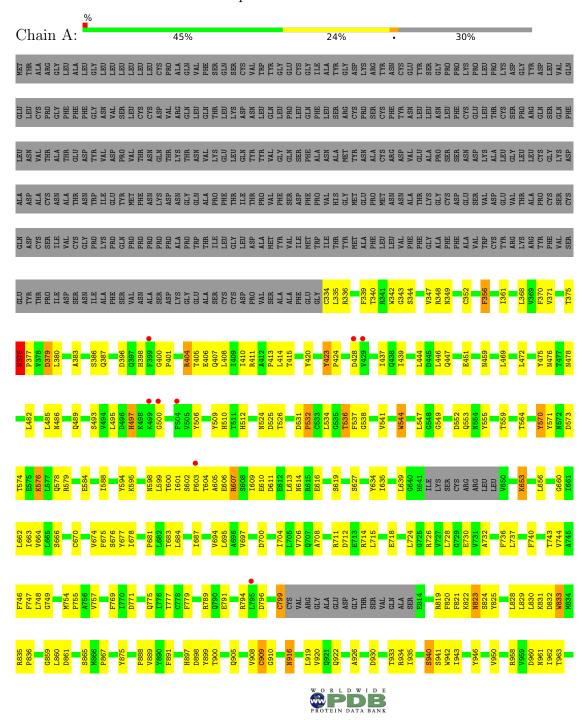
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total 12	C 6	N 1	O 4	S 1	0	0

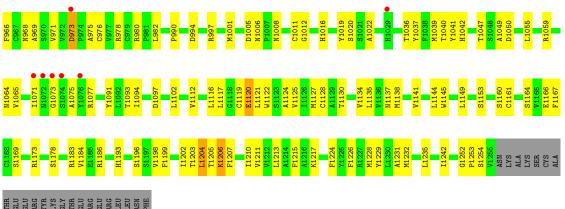


3 Residue-property plots (i)

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: Niemann-Pick C1 protein





THR GLU GLU GRU ARG TYR LYS GLY THR GLU ARG GLU ARG

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

Chain B: 50% 50%



• Molecule 3: alpha-D-mannopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-a cetamido-2-deoxy-beta-D-glucopyranose

Chain C: 67%

• Molecule 4: alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-a cetamido-2-deoxy-beta-D-glucopyranose

Chain D: 33%



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants	174.87Å 222.12Å 107.85Å	Donogitor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	67.92 - 3.33	Depositor
Resolution (A)	67.92 - 3.33	EDS
% Data completeness	96.1 (67.92-3.33)	Depositor
(in resolution range)	96.1 (67.92-3.33)	EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.24 (at 3.33Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
D D.	0.246 , 0.299	Depositor
R, R_{free}	0.246 , 0.299	DCC
R_{free} test set	1492 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	102.6	Xtriage
Anisotropy	0.501	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.28, 56.6	EDS
L-test for twinning ²	$ < L > = 0.45, < L^2> = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	7311	wwPDB-VP
Average B, all atoms (Å ²)	102.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.69% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



¹Intensities estimated from amplitudes.

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: MAN, NAG, MES

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).

Mal	Chain	Bond lengths		Bond angles	
Mol Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.30	0/7259	0.48	1/9879 (0.0%)

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 maintain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$Ideal(^{o})$
1	A	973	ASP	CB-CG-OD1	7.47	125.03	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	536	THR	Peptide

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	7081	0	6994	221	0
2	В	28	0	25	1	0
3	С	39	0	34	1	0
4	D	39	0	34	2	0
5	A	112	0	102	5	0
6	A	12	0	12	1	0
All	All	7311	0	7201	225	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 16.

The worst 5 of 225 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
1:A:1231:ALA:O	1:A:1235:LEU:HB2	1.80	0.81
1:A:771:ASP:O	1:A:775:GLN:HB2	1.82	0.79
1:A:371:VAL:HB	1:A:676:SER:HB2	1.66	0.78
1:A:410:ILE:HD11	1:A:599:LEU:HD13	1.66	0.76
1:A:405:THR:H	1:A:606:GLU:HG3	1.51	0.75

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	897/1278 (70%)	774 (86%)	116 (13%)	7 (1%)	19 53

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	376	ASN
1	A	532	PRO



Continued from previous page...

Mol	Chain	Res	Type
1	A	423	TYR
1	A	571	TYR
1	A	832	ASP

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	Analysed Rotameric Outliers		Percentiles	
1	A	786/1109 (71%)	750 (95%)	36 (5%)	27 60	

5 of 36 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1077	ARG
1	A	1215	PHE
1	A	1128	CYS
1	A	1161	CYS
1	A	576	LYS

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	823	ASN
1	A	953	GLN
1	A	1006	ASN
1	A	1008	ASN

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)

8 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	Во	ond leng	ths	Bond angles		
MIOI	Type	Chain	nes	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	В	1	1,2	14,14,15	0.75	1 (7%)	17,19,21	0.64	1 (5%)
2	NAG	В	2	2	14,14,15	0.27	0	17,19,21	0.63	0
3	NAG	С	1	1,3	14,14,15	0.60	1 (7%)	17,19,21	0.76	1 (5%)
3	NAG	С	2	3	14,14,15	0.40	0	17,19,21	1.00	1 (5%)
3	MAN	С	3	3	11,11,12	1.92	1 (9%)	15,15,17	3.53	5 (33%)
4	NAG	D	1	1,4	14,14,15	1.40	1 (7%)	17,19,21	1.68	2 (11%)
4	NAG	D	2	4	14,14,15	1.78	2 (14%)	17,19,21	1.82	4 (23%)
4	MAN	D	3	4	11,11,12	1.83	2 (18%)	15,15,17	2.22	5 (33%)

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	В	1	1,2	-	4/6/23/26	0/1/1/1
2	NAG	В	2	2	-	2/6/23/26	0/1/1/1
3	NAG	С	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	С	2	3	-	2/6/23/26	0/1/1/1
3	MAN	С	3	3	-	1/2/19/22	0/1/1/1
4	NAG	D	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	D	2	4	-	1/6/23/26	0/1/1/1
4	MAN	D	3	4	-	2/2/19/22	1/1/1/1

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\operatorname{\AA})$	$[Ideal(\AA)]$
4	D	2	NAG	O5-C1	5.52	1.52	1.43



Continued from previous page...

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\text{\AA})$	Ideal(A)
3	С	3	MAN	O5-C5	5.40	1.54	1.43
4	D	1	NAG	O5-C1	4.80	1.51	1.43
4	D	3	MAN	C1-C2	4.58	1.62	1.52
4	D	2	NAG	C1-C2	3.49	1.57	1.52

The worst 5 of 19 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$Ideal(^{o})$
3	С	3	MAN	C1-C2-C3	-8.91	98.72	109.67
4	D	1	NAG	C1-O5-C5	6.28	120.70	112.19
3	С	3	MAN	C1-O5-C5	6.23	120.64	112.19
4	D	2	NAG	C1-O5-C5	5.66	119.87	112.19
3	С	3	MAN	O3-C3-C4	5.32	122.65	110.35

There are no chirality outliers.

5 of 16 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	D	1	NAG	C4-C5-C6-O6
4	D	1	NAG	O5-C5-C6-O6
2	В	1	NAG	O5-C5-C6-O6
2	В	1	NAG	C4-C5-C6-O6
4	D	2	NAG	C1-C2-N2-C7

All (1) ring outliers are listed below:

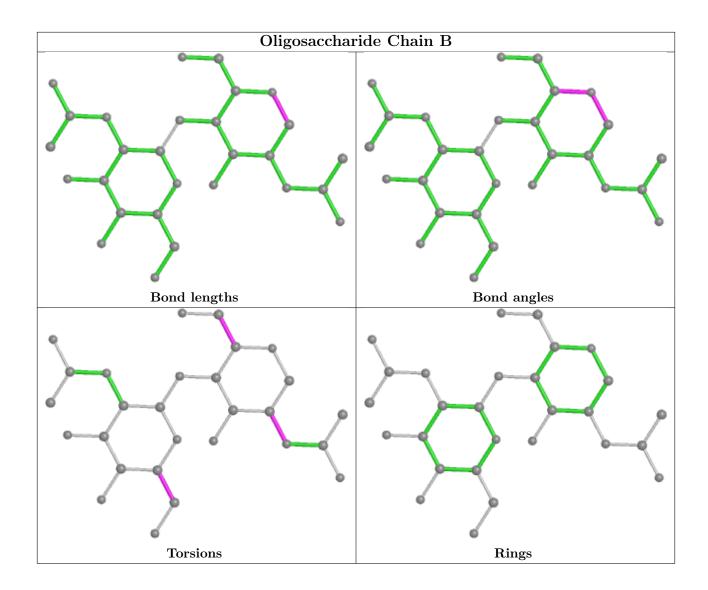
Mol	Chain	Res	Type	Atoms
4	D	3	MAN	C1-C2-C3-C4-C5-O5

4 monomers are involved in 4 short contacts:

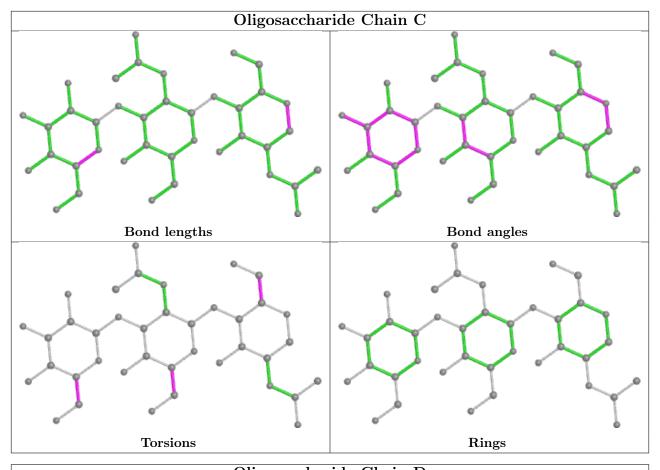
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	С	3	MAN	1	0
2	В	1	NAG	1	0
4	D	2	NAG	2	0
3	С	2	NAG	1	0

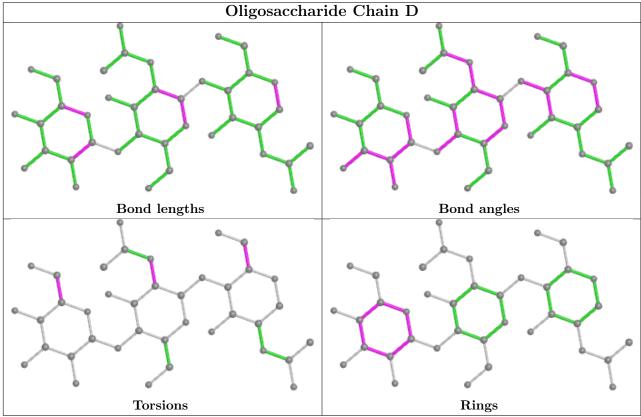
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.













5.6 Ligand geometry (i)

9 ligands 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	Во	ond leng	ths	Bond angles		
IVIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	A	1315	1	14,14,15	0.34	0	17,19,21	0.49	0
5	NAG	A	1309	1	14,14,15	0.15	0	17,19,21	0.42	0
6	MES	A	1317	-	12,12,12	2.26	1 (8%)	14,16,16	1.94	5 (35%)
5	NAG	A	1316	1	14,14,15	0.57	0	17,19,21	1.21	1 (5%)
5	NAG	A	1302	1	14,14,15	0.36	0	17,19,21	0.52	0
5	NAG	A	1310	1	14,14,15	0.77	1 (7%)	17,19,21	0.78	1 (5%)
5	NAG	A	1314	1	14,14,15	0.48	0	17,19,21	0.57	0
5	NAG	A	1301	1	14,14,15	0.21	0	17,19,21	0.36	0
5	NAG	A	1305	1	14,14,15	0.30	0	17,19,21	0.48	0

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
5	NAG	A	1315	1	-	2/6/23/26	0/1/1/1
5	NAG	A	1309	1	-	2/6/23/26	0/1/1/1
6	MES	A	1317	-	-	5/6/14/14	0/1/1/1
5	NAG	A	1316	1	-	5/6/23/26	0/1/1/1
5	NAG	A	1302	1	-	3/6/23/26	0/1/1/1
5	NAG	A	1310	1	-	3/6/23/26	0/1/1/1
5	NAG	A	1314	1	-	1/6/23/26	0/1/1/1
5	NAG	A	1301	1	-	1/6/23/26	0/1/1/1
5	NAG	A	1305	1	-	0/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(A)	$Ideal(\AA)$
6	A	1317	MES	C8-S	-7.53	1.66	1.77
5	A	1310	NAG	O5-C1	2.11	1.47	1.43

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

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
6	A	1317	MES	C5-N4-C3	4.70	119.40	108.83
5	A	1316	NAG	C2-N2-C7	4.13	128.79	122.90
5	A	1310	NAG	C1-O5-C5	2.59	115.70	112.19
6	A	1317	MES	O2S-S-C8	2.43	109.84	106.92
6	A	1317	MES	O1S-S-C8	2.32	109.71	106.92

There are no chirality outliers.

5 of 22 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	1317	MES	N4-C7-C8-S
5	A	1302	NAG	O5-C5-C6-O6
5	A	1309	NAG	O5-C5-C6-O6
5	A	1316	NAG	C8-C7-N2-C2
5	A	1316	NAG	O7-C7-N2-C2

There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	1317	MES	1	0
5	A	1316	NAG	2	0
5	A	1302	NAG	1	0
5	A	1310	NAG	1	0
5	A	1305	NAG	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 (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, 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 $>$	$\# \mathrm{RSRZ}{>}2$	$OWAB(Å^2)$	Q < 0.9
1	A	900/1278 (70%)	0.03	15 (1%) 70 69	58, 96, 143, 248	0

The worst 5 of 15 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1072[A]	ASN	6.3
1	A	973	ASP	4.3
1	A	1071[A]	ILE	3.7
1	A	1073[A]	GLY	3.2
1	A	504	PHE	2.7

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, 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.

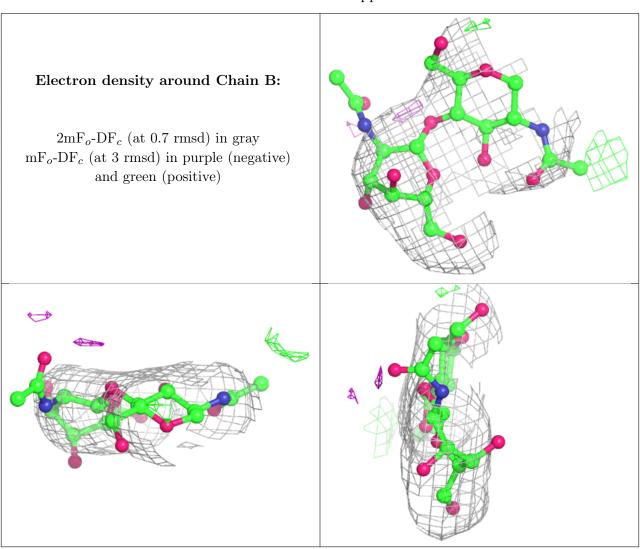
Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathring{\mathbf{A}}^2)$	Q<0.9
4	MAN	D	3	11/12	0.55	0.29	131,162,174,175	0
3	MAN	С	3	11/12	0.76	0.24	119,127,144,152	0
4	NAG	D	1	14/15	0.77	0.53	156,170,179,184	0
2	NAG	В	2	14/15	0.80	0.38	91,146,152,156	0
3	NAG	С	2	14/15	0.84	0.21	78,99,116,143	0
4	NAG	D	2	14/15	0.86	0.27	153,178,184,189	0
2	NAG	В	1	14/15	0.92	0.20	86,112,133,134	0



Continued from previous page...

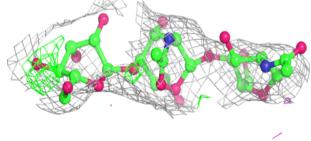
Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
3	NAG	С	1	14/15	0.96	0.13	74,94,106,107	0

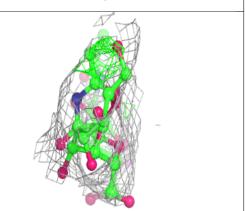
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.





Electron density around Chain C: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray ${ m mF}_o{ m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive) Electron density around Chain D: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray mF_o -DF_c (at 3 rmsd) in purple (negative) and green (positive)







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, 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
5	NAG	A	1301	14/15	0.69	0.41	131,162,171,172	0
5	NAG	A	1315	14/15	0.80	0.30	105,136,153,153	0
5	NAG	A	1314	14/15	0.83	0.19	125,132,142,148	0
5	NAG	A	1302	14/15	0.83	0.34	168,175,181,183	0
5	NAG	A	1316	14/15	0.83	0.53	142,158,171,178	0
5	NAG	A	1310	14/15	0.86	0.22	123,155,166,166	0
5	NAG	A	1305	14/15	0.86	0.16	117,142,148,149	0
5	NAG	A	1309	14/15	0.88	0.20	94,125,139,139	0
6	MES	A	1317	12/12	0.91	0.41	104,110,145,147	0

6.5 Other polymers (i)

There are no such residues in this entry.

