

Full wwPDB X-ray Structure Validation Report (i)

Jan 4, 2021 – 12:11 pm GMT

PDB ID : 6Y6T

Title: Mouse Galactocerebrosidase complexed with galacto-noeurostegine GNS at

pH 4.6

Authors : Deane, J.E.; McLoughlin, J.

Deposited on : 2020-02-27

Resolution : 2.25 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 (i) symbol.

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 : 2.16

buster-report : 1.1.7 (2018)

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

Refmac: 5.8.0158

CCP4 : 7.0.044 (Gargrove)

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

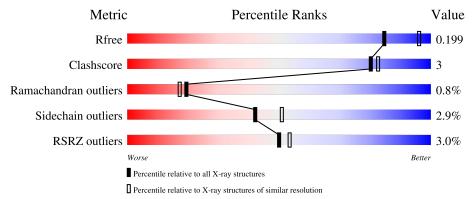
Validation Pipeline (wwPDB-VP) : 2.16

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 2.25 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} \textbf{Similar resolution} \\ (\#\textbf{Entries, resolution range}(\text{\r{A}})) \end{array}$
R_{free}	130704	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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	654	89% 8% ••	
2	В	2	100%	•
2	С	2	100%	•
2	D	2	100%	•



2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 5460 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 Galactocerebrosidase.

\mathbf{Mol}	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	A	641	Total 5100	C 3302	N 849	O 934	S 15	0	1	0

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	15	HIS	_	expression tag	UNP P54818
A	16	HIS	-	expression tag	UNP P54818
A	17	HIS	_	expression tag	UNP P54818
A	18	HIS	-	expression tag	UNP P54818
A	19	HIS	_	expression tag	UNP P54818
A	20	HIS	_	expression tag	UNP P54818
A	21	ILE	_	expression tag	UNP P54818
A	22	GLU	-	expression tag	UNP P54818
A	23	GLY	_	expression tag	UNP P54818
A	24	ARG	_	expression tag	UNP P54818

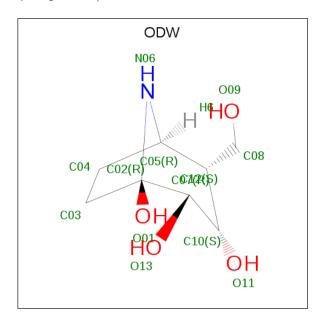
• Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-a cetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
2	В	2	Total C N O 28 16 2 10	0	0	0
2	С	2	Total C N O 28 16 2 10	0	0	0
2	D	2	Total C N O 28 16 2 10	0	0	0

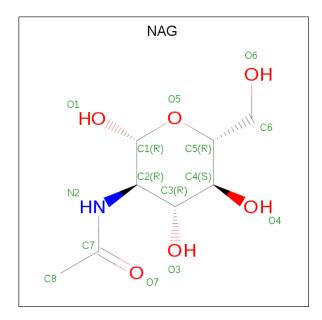


• Molecule 3 is $(1 \{R\}, 2 \{S\}, 3 \{S\}, 4 \{R\}, 5 \{R\})$ -4-(hydroxymethyl)-8-azabicyclo[3.2.1]octane -1,2,3-triol (three-letter code: ODW) (formula: $C_8H_{15}NO_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total 13	C 8	N 1	O 4	0	0

• Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).





\mathbf{Mol}	Chain	Residues	Atoms				ZeroOcc	AltConf
1	Λ	1	Total	С	N	О	0	0
4	Α	1	14	8	1	5	0	0

• Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

• Molecule 6 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total Ni 1 1	0	0

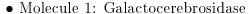
• Molecule 7 is water.

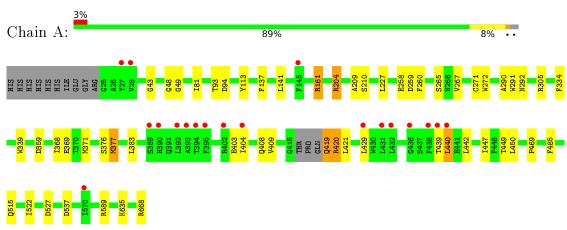
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	247	Total O 247 247	0	0



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

Chain B: 100%

NAG1 NAG2

 $\bullet \ \, \text{Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2$

Chain C:

NAG1 NAG2

 $\bullet \ \, \text{Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2$

Chain D:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants	$250.83 \text{\AA} 250.83 \text{Å} 77.90 \text{Å}$	Danagitan
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	41.98 - 2.25	Depositor
Resolution (A)	41.98 - 2.25	EDS
% Data completeness	99.9 (41.98-2.25)	Depositor
(in resolution range)	99.9 (41.98-2.25)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.41 (at 2.24Å)	Xtriage
Refinement program	PHENIX 1.16_3549, PHENIX 1.16_3549	Depositor
D D.	0.159 , 0.198	Depositor
R, R_{free}	0.160 , 0.199	DCC
R_{free} test set	2278 reflections (5.15%)	wwPDB-VP
Wilson B-factor (Å ²)	46.8	Xtriage
Anisotropy	0.389	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.31 , 45.4	EDS
L-test for twinning ²	$< L > = 0.49, < L^2> = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5460	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.36% 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: NI, CA, NAG, ODW

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	0.39	0/5266	0.56	0/7185	

There are no bond length outliers.

There are no bond angle outliers.

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	5100	0	4820	27	0
2	В	28	0	25	0	0
2	С	28	0	25	0	0
2	D	28	0	25	0	0
3	A	13	0	0	1	0
4	A	14	0	13	0	0
5	A	1	0	0	0	0
6	A	1	0	0	0	0
7	A	247	0	0	4	0
All	All	5460	0	4908	27	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 3.



All (27) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
4 4 055 1 155 1154	4 4 200 7 777 77	distance (Å)	overlap (Å)
1:A:377:MET:HE1	1:A:383:LEU:H	1.56	0.69
1:A:305:ARG:NH1	7:A:803:HOH:O	2.32	0.62
1:A:161:ARG:NH2	7:A:801:HOH:O	2.21	0.61
1:A:339:TRP:CE2	1:A:359:ASP:HB3	2.37	0.59
1:A:377:MET:CE	1:A:383:LEU:H	2.19	0.56
1:A:537:ASP:OD2	1:A:668:ARG:HD2	2.12	0.49
1:A:409:VAL:HG21	1:A:429:LEU:HG	1.95	0.48
1:A:419:GLN:CD	1:A:420:ARG:H	2.17	0.48
1:A:43:GLY:HA3	1:A:81:ILE:HB	1.95	0.48
1:A:334:PHE:CD2	1:A:449:THR:HG21	2.50	0.47
1:A:408:GLN:O	1:A:450:LEU:HA	2.14	0.47
1:A:137:PHE:HB3	1:A:141:LEU:HD12	1.96	0.46
1:A:209:ALA:HA	1:A:210:SER:HA	1.71	0.46
1:A:291:TRP:HA	1:A:292:ASN:HA	1.69	0.46
1:A:371:MET:HB3	1:A:376:SER:HB3	1.98	0.45
1:A:469:PRO:HB3	1:A:668:ARG:NH1	2.32	0.45
1:A:204:ARG:HD2	7:A:996:HOH:O	2.18	0.43
1:A:272:TRP:CH2	1:A:290:ALA:HB2	2.53	0.43
1:A:48:GLY:N	1:A:49:GLY:HA2	2.32	0.43
1:A:49:GLY:O	1:A:94:ASP:HB2	2.20	0.42
1:A:137:PHE:HB2	1:A:522:ILE:HB	2.02	0.41
1:A:227:LEU:HD12	1:A:227:LEU:HA	1.90	0.41
1:A:265:SER:OG	1:A:267:VAL:HG22	2.20	0.41
1:A:485:PHE:HB3	7:A:1035:HOH:O	2.21	0.41
1:A:258:GLU:OE2	3:A:701:ODW:N06	2.54	0.41
1:A:440:LEU:HD12	1:A:442:LEU:HD21	2.02	0.40
1:A:368:ILE:O	1:A:447:ILE:HA	2.21	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	638/654 (98%)	612 (96%)	21 (3%)	5 (1%)	19 17	

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	93	THR
1	A	403	GLU
1	A	420	ARG
1	A	260	PHE
1	A	404	ILE

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	526/555~(95%)	511 (97%)	15 (3%)	42 51	

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	113	TYR
1	A	161	ARG
1	A	204	ARG
1	A	259	ASP
1	A	271	CYS
1	A	369	GLU
1	A	377	MET
1	A	419	GLN
1	A	421	LEU
1	A	439	THR
1	A	440	LEU
1	A	515	GLN
1	A	527	ASP
1	A	589	ARG
1	A	635	LYS



Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	${f Res}$	Type
1	A	538	HIS
1	A	644	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)

6 monosaccharides are modelled in this entry.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

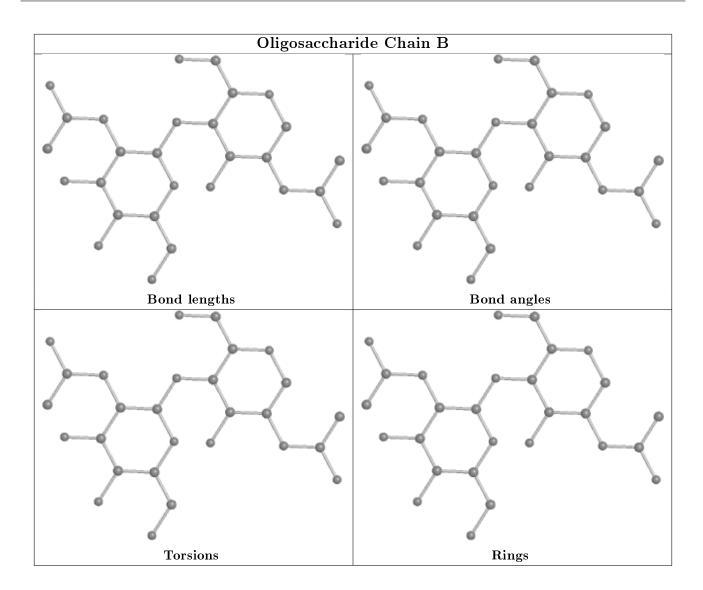
There are no torsion outliers.

There are no ring outliers.

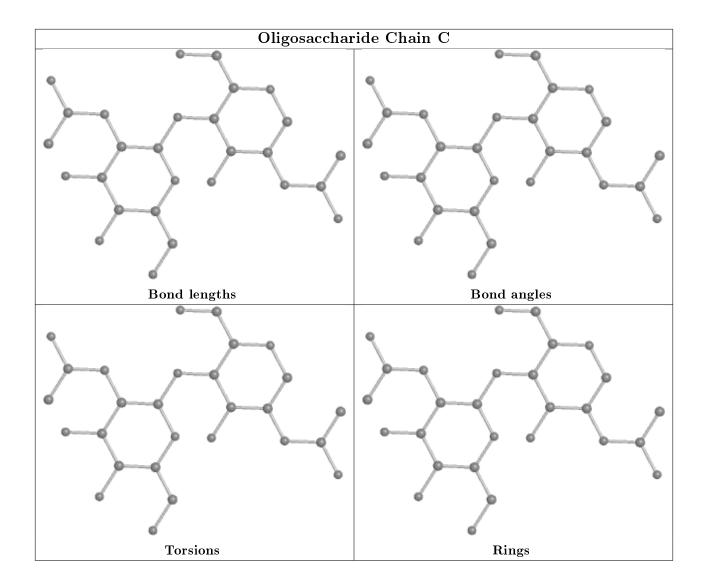
No monomer is involved in short contacts.

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

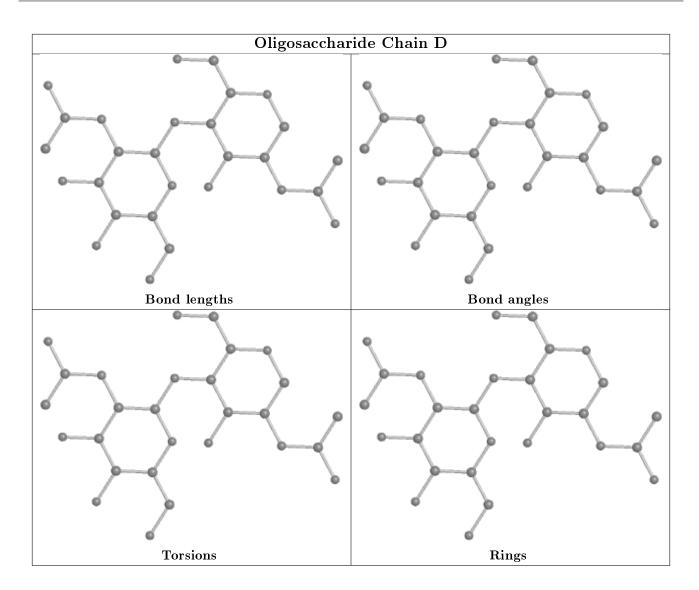












5.6 Ligand geometry (i)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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.

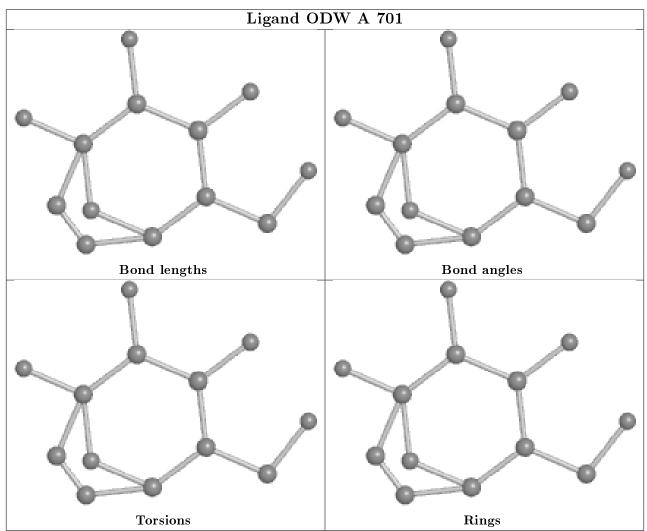
There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier.



Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	$\langle { m RSRZ} \rangle$	$\#\mathrm{RSRZ}{>}2$	$OWAB(A^2)$	Q < 0.9
1	A	641/654 (98%)	-0.24	19 (2%) 50 53	34, 54, 95, 136	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	431	LEU	4.5
1	A	438	PHE	3.4
1	A	393	ALA	3.2
1	A	395	PHE	3.2
1	A	436	GLY	3.1
1	A	392	LEU	3.1
1	A	390	HIS	3.0
1	A	404	ILE	2.9
1	A	394	THR	2.7
1	A	440	LEU	2.7
1	A	432	LEU	2.6
1	A	28	VAL	2.3
1	A	145	PHE	2.2
1	A	27	TYR	2.1
1	A	402	ARG	2.1
1	A	389	SER	2.1
1	A	570	ILE	2.0
1	A	429	LEU	2.0
1	A	439	THR	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.

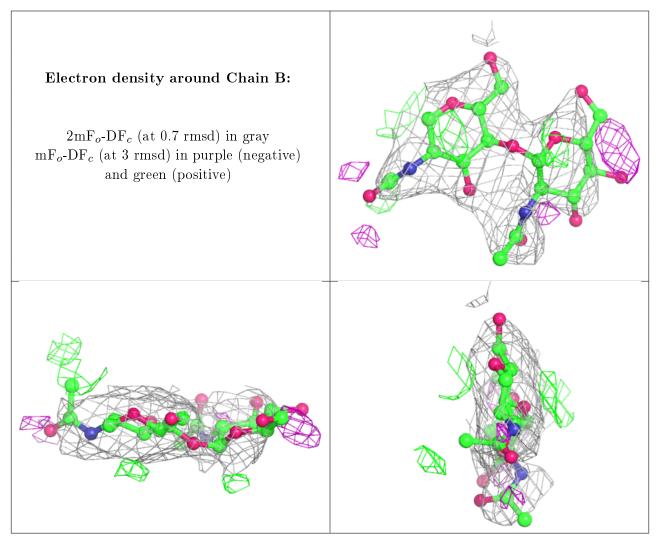


6.3 Carbohydrates (i)

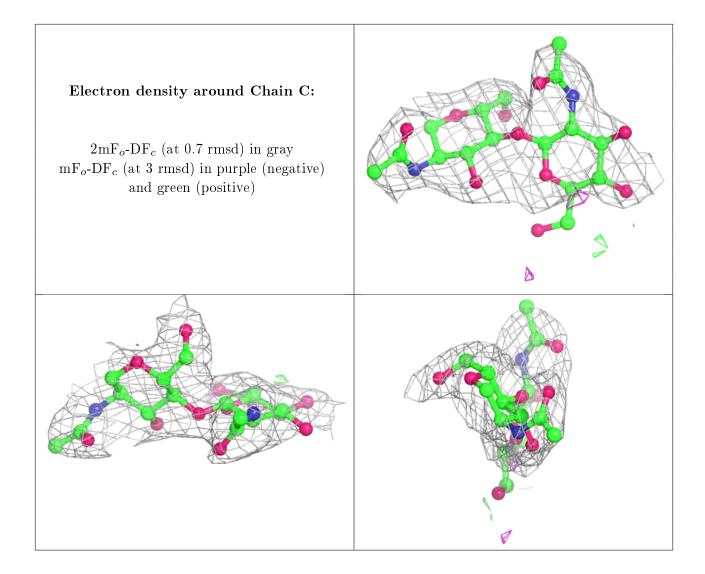
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	$ extbf{B-factors}(extbf{A}^2)$	Q<0.9
2	NAG	В	2	14/15	0.79	0.33	124,137,142,146	0
2	NAG	D	2	14/15	0.83	0.29	82,108,117,120	0
2	NAG	С	2	14/15	0.86	0.37	102,122,127,127	0
2	NAG	С	1	14/15	0.89	0.23	81,92,100,109	0
2	NAG	В	1	14/15	0.94	0.13	68,89,108,122	0
2	NAG	D	1	14/15	0.95	0.19	62,73,88,101	0

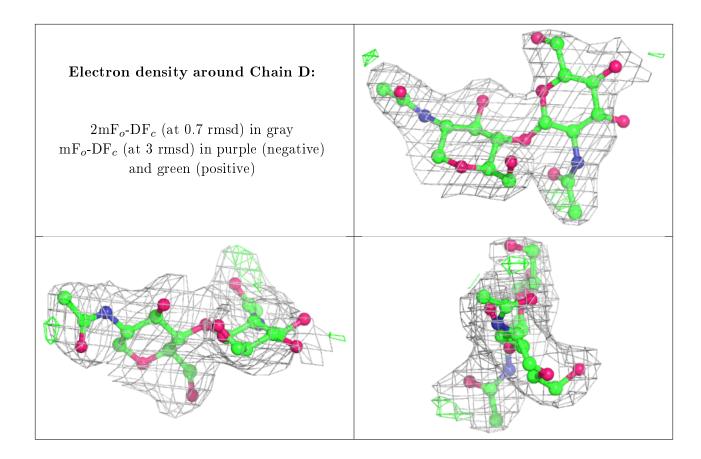
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.











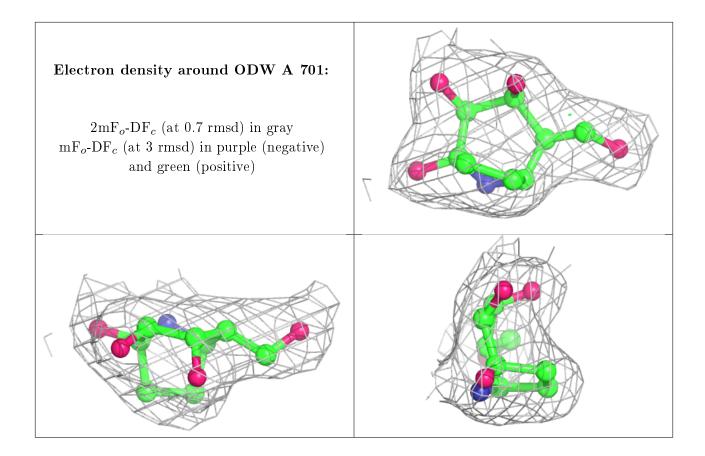
6.4 Ligands (i)

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	\mathbf{Res}	Atoms	RSCC	RSR	${f B-factors}({f A}^2)$	Q<0.9
4	NAG	A	706	14/15	0.81	0.33	95,109,119,122	0
6	NI	A	710	1/1	0.96	0.12	107,107,107,107	0
3	ODW	A	701	13/13	0.96	0.18	36,41,48,50	0
5	CA	A	709	1/1	0.99	0.03	69,69,69,69	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.





6.5 Other polymers (i)

There are no such residues in this entry.

