

# wwPDB X-ray Structure Validation Summary Report (i)

#### Sep 23, 2023 – 08:16 PM EDT

PDB ID : 5T0D

Title: Crystal structure of H6 hemagglutinin G225D mutant from Taiwan (2013)

H6N1 influenza virus in complex with 3'-SLN

Authors: Wilson, I.A.; Tzarum, N.; Zhu, X.

Deposited on : 2016-08-15

Resolution : 2.86 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:

 $Mol Probity \quad : \quad 4.02b\text{--}467$ 

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)

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

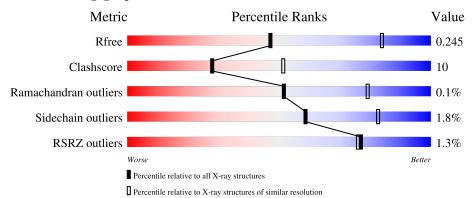
Validation Pipeline (wwPDB-VP) : 2.35.1

## 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 2.86 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	Similar resolution
Metric	$(\# \mathrm{Entries})$	$(\#  ext{Entries},  ext{ resolution range}(\AA))$
$R_{free}$	130704	3168 (2.90-2.82)
Clashscore	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)
RSRZ outliers	127900	3103 (2.90-2.82)

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	Δ.	222	2%		
1	A	333	70%	26%	• •
1	С	333	75%	22%	• •
	_	222	<u>%</u>		
1	Е	333	75%	21%	•••
	D	100	<u>%</u>		
2	В	180	72%	23%	• •
	_		2%		
2	D	180	81%	14%	• •

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_			F			
	Mol	$\mathbf{Chain}$	Length	Quality of chain		
L						
				2%		
	0	-	100			
	2	F.	180	74%	21%	<del></del>
$\vdash$						
	9	$\alpha$	9			
	3	G	3	100%		



## 2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 12070 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 Hemagglutinin.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	330	Total	С	N	О	S	0	0	0
1	A	330	2604	1650	443	498	13	0		
1	С	325	Total	С	N	О	S	0	0	0
1		329	2566	1627	436	490	13	U	U	
1	Е	205	Total	С	N	О	S	0	0	0
1	E	E 325	2569	1629	437	490	13		U	

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	225	ASP	GLY	engineered mutation	UNP A0A0J9X268
С	225	ASP	GLY	engineered mutation	UNP A0A0J9X268
Е	225	ASP	GLY	engineered mutation	UNP A0A0J9X268

• Molecule 2 is a protein called Hemagglutinin HA2 chain.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
2	В	173	Total	С	N	О	S	0	0	0
2	Ъ	175	1393	868	244	274	7	0	U	U
2	D	173	Total	С	N	О	S	0	0	0
2	D	175	1393	868	244	274	7	0	0	
9	E	172	Total	С	N	О	S	0	0	0
	Г	112	1385	862	243	273	7	0		

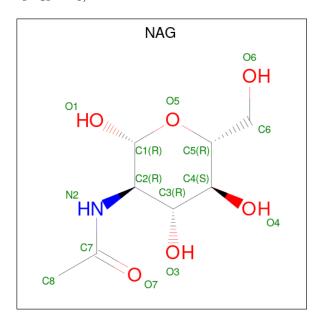
• Molecule 3 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galacto pyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.





Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace		
3	G	3	Total 46	C 25	N 2	O 19	0	0	0

 $\bullet$  Molecule 4 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
4	A	1	Total C N O 14 8 1 5	0	0
4	A	1	Total C N O 14 8 1 5	0	0
4	Е	1	Total C N O 14 8 1 5	0	0

• Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	22	Total O 22 22	0	0
5	В	9	Total O 9 9	0	0
5	С	9	Total O 9 9	0	0
5	D	5	Total O 5 5	0	0
5	Е	24	Total O 24 24	0	0

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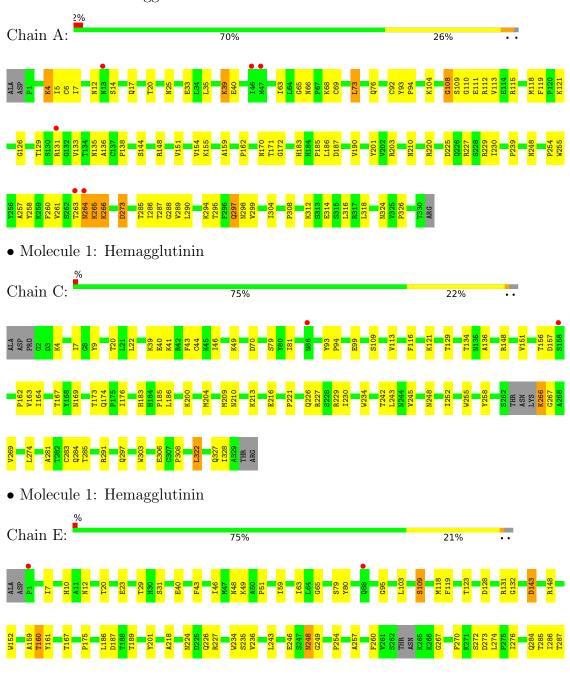
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	F	3	Total O 3 3	0	0



## 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: Hemagglutinin







• Molecule 2: Hemagglutinin HA2 chain

Chain B: 72% 23% . .



• Molecule 2: Hemagglutinin HA2 chain

Chain D: 2%

14% • •



ARG

• Molecule 2: Hemagglutinin HA2 chain

Chain F: 74% 21% . .



 $\bullet$  Molecule 3: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:

NAG1 GAL2 SIA3



## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants	185.06Å 98.73Å 132.79Å	Donositon
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $126.76^{\circ}$ $90.00^{\circ}$	Depositor
Resolution (Å)	49.38 - 2.86	Depositor
Resolution (A)	49.38 - 2.86	EDS
% Data completeness	97.2 (49.38-2.86)	Depositor
(in resolution range)	97.2 (49.38-2.86)	EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.08	Depositor
$< I/\sigma(I) > 1$	3.33 (at 2.86Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
D D.	0.206 , 0.245	Depositor
$R, R_{free}$	0.207 , $0.245$	DCC
$R_{free}$ test set	2197 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	43.8	Xtriage
Anisotropy	0.951	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.35, 49.6	EDS
L-test for twinning <sup>2</sup>	$ < L >=0.50, < L^2>=0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	12070	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.94% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>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.



<sup>&</sup>lt;sup>1</sup>Intensities 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: SIA, NAG, GAL

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		
IVIOI		RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	0.48	0/2666	0.65	3/3628~(0.1%)	
1	С	0.48	0/2626	0.62	0/3572	
1	Е	0.47	0/2630	0.61	0/3576	
2	В	0.47	0/1421	0.61	1/1913~(0.1%)	
2	D	0.46	0/1421	0.59	1/1913 (0.1%)	
2	F	0.45	0/1413	0.62	1/1902~(0.1%)	
All	All	0.47	0/12177	0.62	6/16504~(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 mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	Е	0	1
All	All	0	2

There are no bond length outliers.

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

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$Observed(^o)$	$\mathrm{Ideal}(^{o})$
2	D	1	GLY	N-CA-C	-7.23	95.02	113.10
1	A	264	ASN	N-CA-C	-6.37	93.80	111.00
1	A	4	LYS	CD-CE-NZ	6.25	126.07	111.70
2	В	1	GLY	N-CA-C	-5.70	98.84	113.10
1	A	108	GLY	N-CA-C	-5.46	99.44	113.10

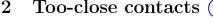
There are no chirality outliers.

All (2) planarity outliers are listed below:



Mol	Chain	Res	Type	Group
1	A	172	GLY	Peptide
1	Е	109	SER	Peptide

#### Too-close contacts (i) 5.2



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	2604	0	2552	74	0
1	С	2566	0	2510	51	0
1	Е	2569	0	2516	54	0
2	В	1393	0	1309	32	0
2	D	1393	0	1309	21	0
2	F	1385	0	1298	27	0
3	G	46	0	40	4	0
4	A	28	0	26	0	0
4	Е	14	0	13	0	0
5	A	22	0	0	1	0
5	В	9	0	0	4	0
5	С	9	0	0	0	0
5	D	5	0	0	0	0
5	Е	24	0	0	4	0
5	F	3	0	0	0	0
All	All	12070	0	11573	233	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 10.

The worst 5 of 233 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
1:A:264:ASN:O	1:A:265:LYS:CG	1.79	1.30
1:A:264:ASN:O	1:A:265:LYS:HG3	0.95	1.10
1:A:73:LEU:HD11	1:A:148:ARG:HG2	1.52	0.92
1:A:266:LYS:O	1:A:304:ILE:HD12	1.70	0.92
1:C:134:THR:HG22	1:C:136:ALA:H	1.33	0.90

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	Perce	ntiles
1	A	328/333 (98%)	316 (96%)	11 (3%)	1 (0%)	41	68
1	$\mathbf{C}$	321/333 (96%)	312 (97%)	9 (3%)	0	100	100
1	E	321/333 (96%)	303 (94%)	18 (6%)	0	100	100
2	В	171/180 (95%)	169 (99%)	2 (1%)	0	100	100
2	D	171/180 (95%)	163 (95%)	8 (5%)	0	100	100
2	F	170/180 (94%)	165 (97%)	5 (3%)	0	100	100
All	All	1482/1539 (96%)	1428 (96%)	53 (4%)	1 (0%)	51	79

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	265	LYS

### 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	nalysed Rotameric Outliers		Percentiles		
1	A	$290/292\ (99\%)$	282 (97%)	8 (3%)	43	73	
1	С	285/292 (98%)	282 (99%)	3 (1%)	73	90	
1	E	286/292 (98%)	282 (99%)	4 (1%)	67	86	
2	В	$147/153\ (96\%)$	144 (98%)	3 (2%)	55	80	
2	D	$147/153\ (96\%)$	143 (97%)	4 (3%)	44	74	
2	F	146/153 (95%)	144 (99%)	2 (1%)	67	86	

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Mol	Chain	Analysed	Analysed Rotameric Outlie		Percentiles
All	All	$1301/1335\ (98\%)$	1277 (98%)	24 (2%)	59 82

5 of 24 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	22	TYR
2	D	145	ASP
2	D	144	CYS
1	Е	123	THR
1	A	273	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
2	В	163	GLN
1	С	226	GLN
2	D	169	ASN
1	Ε	244	ASN
1	Е	248	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)

3 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	Tuno	Chain	Chain	Chain	Chain	Chain	Chain	Chain	nain Res	Link	Bond lengths			Bond angles		
Mol Type Chain	Res Lilik		Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2								
3	NAG	G	1	3	15,15,15	0.57	0	21,21,21	1.68	3 (14%)						
3	GAL	G	2	3	11,11,12	1.65	2 (18%)	15,15,17	1.95	4 (26%)						
3	SIA	G	3	3	20,20,21	0.76	0	24,28,31	1.17	3 (12%)						

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	G	1	3	-	5/6/26/26	0/1/1/1
3	GAL	G	2	3	-	2/2/19/22	0/1/1/1
3	SIA	G	3	3	-	0/18/34/38	0/1/1/1

#### All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$Observed(\AA)$	$\operatorname{Ideal}( ext{\AA})$
3	G	2	GAL	O5-C1	-3.71	1.37	1.43
3	G	2	GAL	C1-C2	3.30	1.59	1.52

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
3	G	2	GAL	O2-C2-C3	-4.41	101.30	110.14
3	G	1	NAG	O5-C1-C2	-4.06	105.43	109.52
3	G	1	NAG	O4-C4-C3	-3.42	102.44	110.35
3	G	2	GAL	C1-O5-C5	-3.25	107.79	112.19
3	G	2	GAL	O2-C2-C1	-3.19	102.63	109.15

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	G	1	NAG	C1-C2-N2-C7
3	G	1	NAG	C8-C7-N2-C2
3	G	1	NAG	O7-C7-N2-C2
3	G	1	NAG	O5-C5-C6-O6
3	G	2	GAL	O5-C5-C6-O6

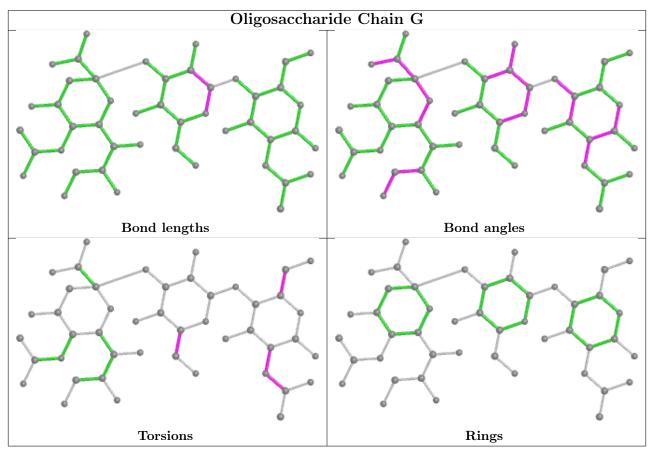
There are no ring outliers.



0			•	11	•	4	1 .	1 1
3	monomers	are	invo	lved	ın	4 S	hort	contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	3	SIA	2	0
3	G	2	GAL	1	0
3	G	1	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)

#### 3 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	Bo	nd leng	$ ag{ths}$	В	ond ang	les
	Type				Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z >2
4	NAG	A	401	1	14,14,15	0.54	0	17,19,21	0.66	0
4	NAG	A	402	1	14,14,15	0.48	0	17,19,21	1.23	3 (17%)
4	NAG	Е	401	1	14,14,15	0.83	1 (7%)	17,19,21	0.71	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
4	NAG	A	401	1	-	4/6/23/26	0/1/1/1
4	NAG	A	402	1	-	2/6/23/26	0/1/1/1
4	NAG	Е	401	1	-	1/6/23/26	0/1/1/1

#### All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	$\mathbf{E}$	401	NAG	O5-C1	-2.55	1.39	1.43

#### All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
4	A	402	NAG	C1-O5-C5	2.74	115.90	112.19
4	A	402	NAG	C2-N2-C7	-2.70	119.06	122.90
4	A	402	NAG	C1-C2-N2	2.07	114.03	110.49

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	401	NAG	C8-C7-N2-C2
4	A	401	NAG	O7-C7-N2-C2
4	A	401	NAG	O5-C5-C6-O6
4	A	401	NAG	C1-C2-N2-C7
4	A	402	NAG	C8-C7-N2-C2

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,  $95^{th}$  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>	$\#\mathrm{RSRZ}{>}2$	$OWAB(A^2)$	Q < 0.9
1	A	330/333~(99%)	0.05	6 (1%) 68 66	29, 44, 63, 83	0
1	С	325/333 (97%)	-0.10	2 (0%) 89 89	28, 40, 58, 80	0
1	E	325/333~(97%)	-0.10	2 (0%) 89 89	28, 42, 55, 80	0
2	В	173/180 (96%)	-0.08	2 (1%) 79 78	21, 40, 58, 73	0
2	D	173/180 (96%)	0.00	3 (1%) 70 68	26, 41, 61, 92	0
2	F	172/180 (95%)	-0.05	4 (2%) 60 57	26, 42, 61, 75	0
All	All	1498/1539 (97%)	-0.05	19 (1%) 77 76	21, 42, 60, 92	0

The worst 5 of 19 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	В	1	GLY	4.7
2	D	1	GLY	4.7
2	F	1	GLY	4.4
1	С	158	SER	4.1
1	A	47	MET	3.8

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

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

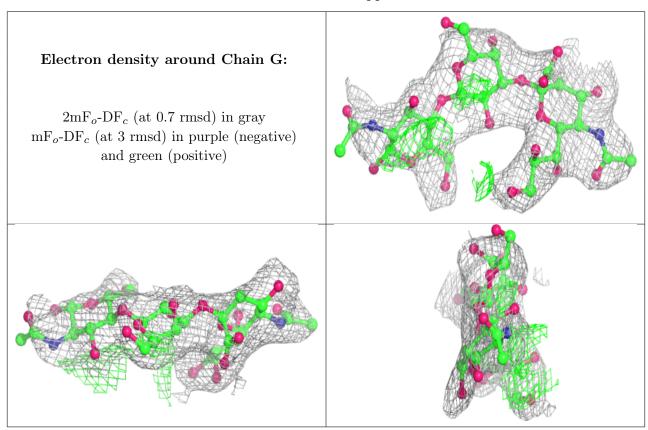
### 6.3 Carbohydrates (i)

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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
3	NAG	G	1	15/15	0.82	0.31	67,88,94,95	0
3	GAL	G	2	11/12	0.86	0.19	62,74,83,85	0
3	SIA	G	3	20/21	0.91	0.22	50,61,71,73	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	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
4	NAG	A	401	14/15	0.77	0.26	60,73,77,86	0
4	NAG	A	402	14/15	0.83	0.18	46,59,67,72	0
4	NAG	E	401	14/15	0.85	0.22	66,77,88,90	0



## 6.5 Other polymers (i)

There are no such residues in this entry.

