



# wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 16, 2016 – 07:44 AM EDT

PDB ID : 4YY0  
Title : The structure of hemagglutinin from a H6N1 influenza virus (A/chicken/Taiwan/A2837/2013)  
Authors : Wang, F.; Qi, J.; Bi, Y.; Zhang, W.; Wang, M.; Wang, M.; Liu, J.; Yan, J.; Shi, Y.; Gao, G.F.  
Deposited on : 2015-03-23  
Resolution : 2.59 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20027939  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20027939

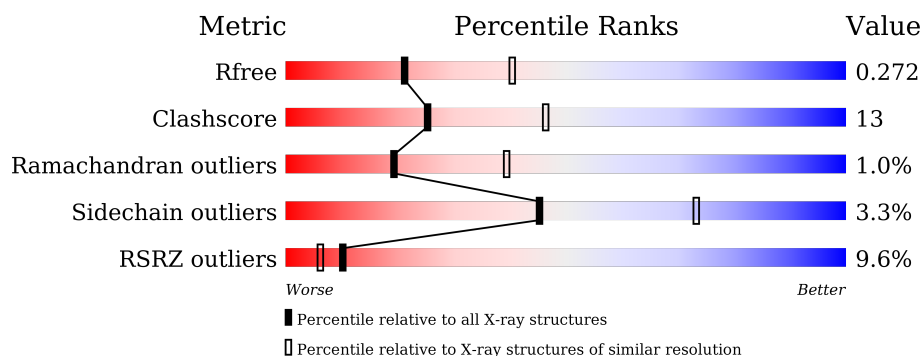
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.59 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))
$R_{free}$	91344	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	325	<div> <div>2%</div> <div>85%</div> <div>13%</div> <div>..</div> </div>
1	C	325	<div> <div>5%</div> <div>83%</div> <div>15%</div> <div>.</div> </div>
1	E	325	<div> <div>6%</div> <div>80%</div> <div>18%</div> <div>..</div> </div>
2	B	159	<div> <div>15%</div> <div>74%</div> <div>24%</div> <div>..</div> </div>
2	D	159	<div> <div>24%</div> <div>67%</div> <div>27%</div> <div>5%</div> <div>.</div> </div>
2	F	159	<div> <div>23%</div> <div>69%</div> <div>22%</div> <div>8%</div> <div>.</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11681 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 HA1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	325	Total	C	N	O	S	0	0	0
			2563	1625	436	489	13			
1	C	325	Total	C	N	O	S	0	0	0
			2563	1625	436	489	13			
1	E	325	Total	C	N	O	S	0	0	0
			2563	1625	436	489	13			

- Molecule 2 is a protein called HA2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	159	Total	C	N	O	S	0	0	0
			1275	794	221	253	7			
2	D	159	Total	C	N	O	S	0	0	0
			1275	794	221	253	7			
2	F	159	Total	C	N	O	S	0	0	0
			1275	794	221	253	7			

- Molecule 3 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	D	1	Total	C	N	O	0	0
			14	8	1	5		
3	E	1	Total	C	N	O	0	0
			14	8	1	5		
3	E	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	30	Total	O	0	0
			30	30		
4	B	1	Total	O	0	0
			1	1		
4	C	15	Total	O	0	0
			15	15		
4	D	5	Total	O	0	0
			5	5		
4	E	13	Total	O	0	0
			13	13		

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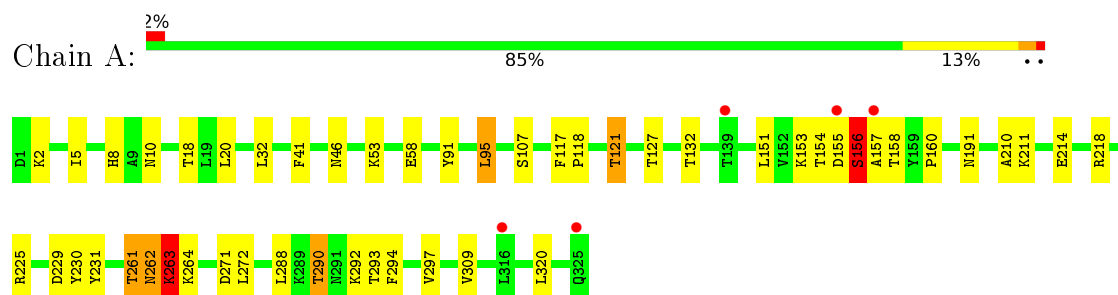
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	F	5	Total	O	0	0
			5	5		

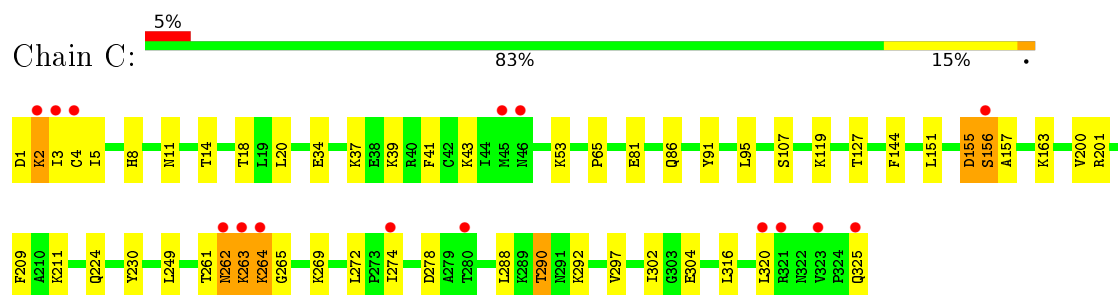
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ( $\text{RSRZ} > 2$ ). 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.

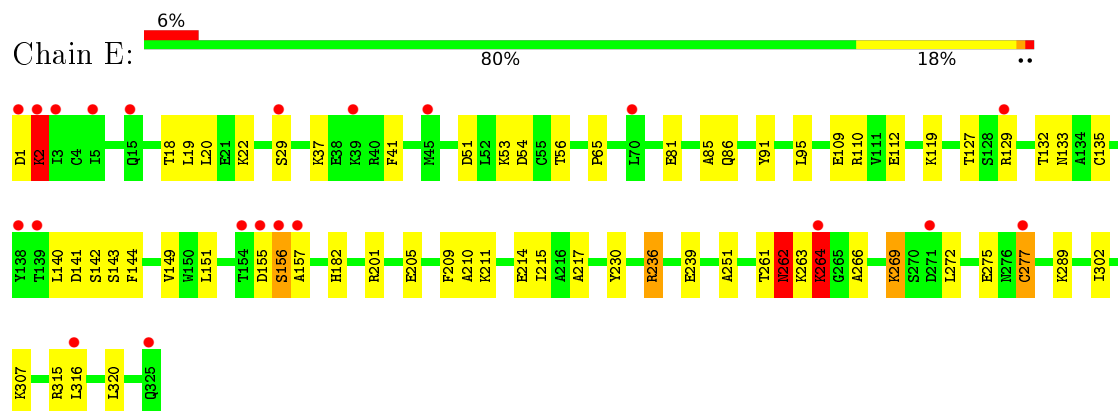
#### • Molecule 1: HA1



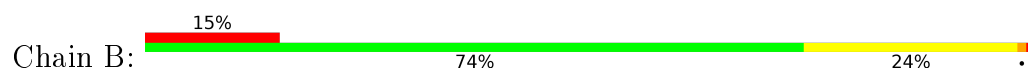
#### • Molecule 1: HA1

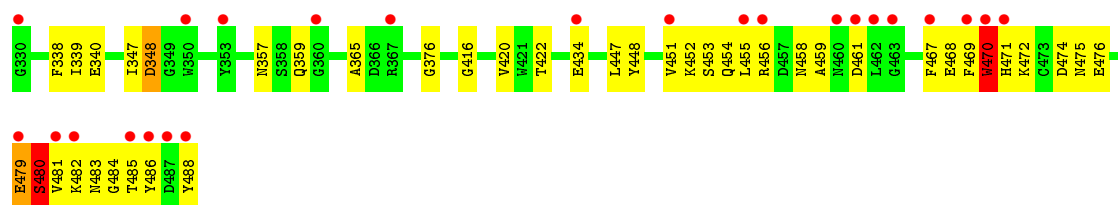


#### • Molecule 1: HA1

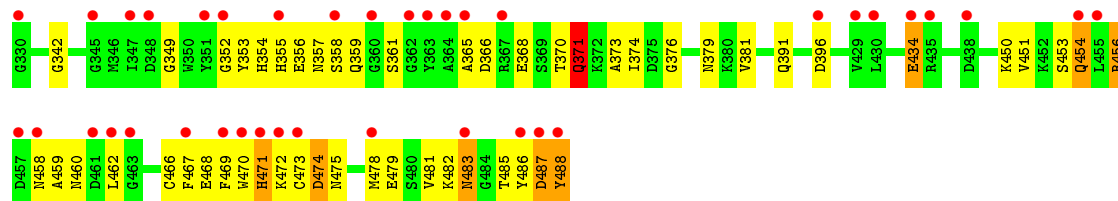


#### • Molecule 2: HA2

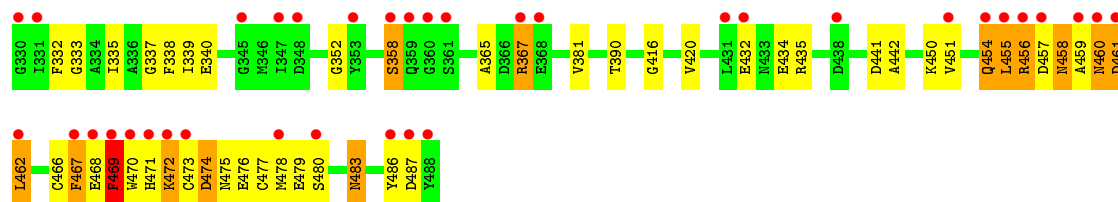




• Molecule 2: HA2



• Molecule 2: HA2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	67.61Å 106.26Å 125.28Å 90.00° 102.75° 90.00°	Depositor
Resolution (Å)	41.37 – 2.59 41.37 – 2.59	Depositor EDS
% Data completeness (in resolution range)	99.2 (41.37-2.59) 99.2 (41.37-2.59)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.63 (at 2.58Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.8.4_1496)	Depositor
R, $R_{free}$	0.220 , 0.271 0.222 , 0.272	Depositor DCC
$R_{free}$ test set	2711 reflections (5.09%)	DCC
Wilson B-factor (Å <sup>2</sup> )	51.1	Xtriage
Anisotropy	0.402	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 41.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.024 for h,-k,-h-l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	11681	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.95% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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 ⓘ

### 5.1 Standard geometry ⓘ

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

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	0.31	0/2625	0.61	4/3571 (0.1%)
1	C	0.31	0/2625	0.56	1/3571 (0.0%)
1	E	0.34	0/2625	0.63	1/3571 (0.0%)
2	B	0.48	0/1301	0.85	2/1754 (0.1%)
2	D	0.52	1/1301 (0.1%)	0.76	3/1754 (0.2%)
2	F	0.56	3/1301 (0.2%)	0.87	6/1754 (0.3%)
All	All	0.40	4/11778 (0.0%)	0.68	17/15975 (0.1%)

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
2	B	0	2
2	F	0	1
All	All	0	3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	469	PHE	CG-CD1	-7.08	1.28	1.38
2	D	434	GLU	CB-CG	-5.66	1.41	1.52
2	F	467	PHE	CD1-CE1	-5.56	1.28	1.39
2	F	469	PHE	CD1-CE1	-5.09	1.29	1.39

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	262	ASN	N-CA-CB	-12.28	88.49	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	261	THR	N-CA-C	-11.74	79.31	111.00
1	A	262	ASN	N-CA-CB	-9.79	92.98	110.60
2	F	462	LEU	CA-CB-CG	8.92	135.81	115.30
1	C	265	GLY	N-CA-C	-7.33	94.77	113.10

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	470	TRP	Peptide
2	B	479	GLU	Peptide
2	F	358	SER	Peptide

## 5.2 Too-close contacts ⓘ

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	2563	0	2508	41	0
1	C	2563	0	2508	52	0
1	E	2563	0	2507	72	0
2	B	1275	0	1180	50	0
2	D	1275	0	1178	57	0
2	F	1275	0	1180	54	0
3	A	28	0	26	0	0
3	C	28	0	26	1	0
3	D	14	0	13	3	0
3	E	28	0	26	0	0
4	A	30	0	0	4	0
4	B	1	0	0	0	0
4	C	15	0	0	2	0
4	D	5	0	0	0	0
4	E	13	0	0	3	0
4	F	5	0	0	0	0
All	All	11681	0	11152	286	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.

The worst 5 of 286 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:E:109:GLU:HB2	1:E:261:THR:CG2	1.86	1.04
1:A:261:THR:O	1:A:261:THR:CG2	2.03	1.02
1:A:261:THR:HG22	1:A:261:THR:O	1.63	0.96
2:B:471:HIS:HD2	2:B:472:LYS:H	1.11	0.94
1:C:2:LYS:NZ	1:C:4:CYS:SG	2.41	0.92

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	323/325 (99%)	306 (95%)	15 (5%)	2 (1%)	30	56
1	C	323/325 (99%)	306 (95%)	13 (4%)	4 (1%)	16	33
1	E	323/325 (99%)	308 (95%)	12 (4%)	3 (1%)	21	42
2	B	157/159 (99%)	148 (94%)	6 (4%)	3 (2%)	10	19
2	D	157/159 (99%)	147 (94%)	8 (5%)	2 (1%)	15	30
2	F	157/159 (99%)	150 (96%)	6 (4%)	1 (1%)	30	56
All	All	1440/1452 (99%)	1365 (95%)	60 (4%)	15 (1%)	19	39

5 of 15 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	155	ASP
2	B	480	SER
1	C	156	SER
1	C	263	LYS
2	D	483	ASN

### 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	285/285 (100%)	279 (98%)	6 (2%)	61	85
1	C	285/285 (100%)	279 (98%)	6 (2%)	61	85
1	E	285/285 (100%)	277 (97%)	8 (3%)	51	78
2	B	134/134 (100%)	131 (98%)	3 (2%)	60	83
2	D	134/134 (100%)	126 (94%)	8 (6%)	24	47
2	F	134/134 (100%)	124 (92%)	10 (8%)	17	33
All	All	1257/1257 (100%)	1216 (97%)	41 (3%)	45	73

5 of 41 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	396	ASP
2	D	487	ASP
2	F	461	ASP
2	D	450	LYS
2	D	454	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
2	D	458	ASN
2	D	471	HIS
1	E	262	ASN
1	C	224	GLN
2	D	475	ASN

### 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 ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

7 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	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	NAG	A	401	1	14,14,15	0.26	0	15,19,21	0.61	1 (6%)
3	NAG	A	402	1	14,14,15	0.22	0	15,19,21	0.54	0
3	NAG	C	401	1	14,14,15	0.50	0	15,19,21	0.55	0
3	NAG	C	402	1	14,14,15	0.31	0	15,19,21	0.81	1 (6%)
3	NAG	D	501	2	14,14,15	0.59	0	15,19,21	0.91	1 (6%)
3	NAG	E	401	1	14,14,15	1.31	2 (14%)	15,19,21	1.01	1 (6%)
3	NAG	E	402	1	14,14,15	0.22	0	15,19,21	0.71	1 (6%)

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	A	401	1	-	0/6/23/26	0/1/1/1
3	NAG	A	402	1	-	0/6/23/26	0/1/1/1
3	NAG	C	401	1	-	0/6/23/26	0/1/1/1
3	NAG	C	402	1	-	0/6/23/26	0/1/1/1
3	NAG	D	501	2	-	0/6/23/26	0/1/1/1
3	NAG	E	401	1	-	0/6/23/26	0/1/1/1
3	NAG	E	402	1	-	0/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	401	NAG	O5-C1	2.98	1.48	1.43
3	E	401	NAG	C1-C2	3.62	1.57	1.52

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	501	NAG	C1-O5-C5	-2.42	108.58	112.14
3	A	401	NAG	C1-O5-C5	2.14	115.28	112.14
3	E	401	NAG	O3-C3-C4	2.35	115.66	110.36
3	E	402	NAG	C1-O5-C5	2.36	115.61	112.14
3	C	402	NAG	C1-O5-C5	2.70	116.11	112.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	402	NAG	1	0
3	D	501	NAG	3	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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	325/325 (100%)	0.20	5 (1%) 76 71	32, 54, 92, 147	0
1	C	325/325 (100%)	0.29	15 (4%) 36 29	33, 57, 106, 154	0
1	E	325/325 (100%)	0.50	21 (6%) 22 16	39, 71, 110, 168	0
2	B	159/159 (100%)	0.99	24 (15%) 3 1	40, 85, 164, 215	0
2	D	159/159 (100%)	1.39	38 (23%) 1 0	37, 96, 166, 232	0
2	F	159/159 (100%)	1.34	36 (22%) 1 0	37, 95, 174, 220	0
All	All	1452/1452 (100%)	0.63	139 (9%) 10 6	32, 67, 146, 232	0

The worst 5 of 139 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	488	TYR	15.9
2	B	488	TYR	9.4
2	B	487	ASP	7.9
2	F	467	PHE	7.5
2	F	455	LEU	7.3

### 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](#)

There are no carbohydrates in this entry.

## 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	NAG	C	402	14/15	0.86	0.18	0.43	78,93,104,107	0
3	NAG	E	402	14/15	0.90	0.20	0.25	64,80,88,88	0
3	NAG	A	401	14/15	0.73	0.36	-	119,124,128,129	0
3	NAG	C	401	14/15	0.42	0.37	-	140,159,164,164	0
3	NAG	D	501	14/15	0.71	0.35	-	166,171,173,174	0
3	NAG	A	402	14/15	0.86	0.15	-	66,80,93,95	0
3	NAG	E	401	14/15	0.79	0.20	-	140,148,152,154	0

## 6.5 Other polymers

There are no such residues in this entry.