



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:02 PM GMT

PDB ID : 4UO8  
Title : Structure of the A\_Canine\_Colorado\_17864\_06 H3 haemagglutinin in complex with 6SO4-3SLN  
Authors : Vachieri, S.G.; Collins, P.J.; Haire, L.F.; Ogrodowicz, R.W.; Martin, S.R.; Walker, P.A.; Xiong, X.; Gamblin, S.J.; Skehel, J.J.  
Deposited on : 2014-05-31  
Resolution : 3.00 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](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 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
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) : trunk26865

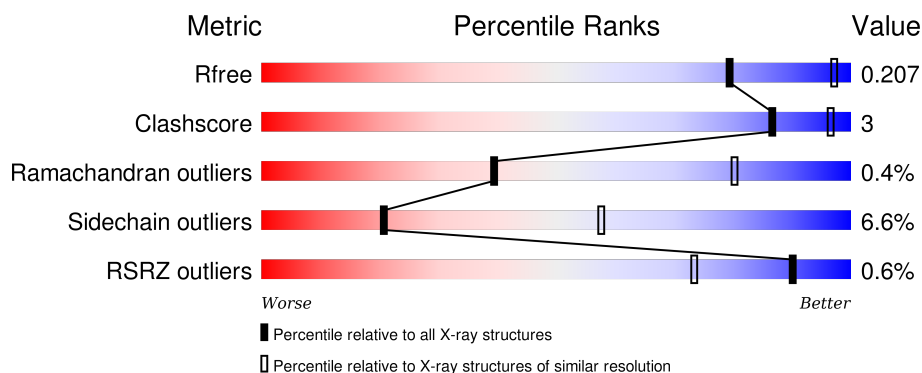
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*



The reported resolution of this entry is 3.00 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	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

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	328	 85% 11% . .
2	B	175	 % 89% 10% .

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
5	NAG	A	621	X	-	-	-
5	MAN	A	625	X	-	-	-
6	NAG	A	641	-	-	-	X

## 2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 4190 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	319	Total	C	N	O	S	0	0	0
			2468	1541	434	477	16			

- Molecule 2 is a protein called HA2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	175	Total	C	N	O	S	0	0	0
			1425	886	250	283	6			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	173	SER	-	EXPRESSION TAG	UNP E0UVR5
B	174	GLY	-	EXPRESSION TAG	UNP E0UVR5
B	175	ARG	-	EXPRESSION TAG	UNP E0UVR5
B	131	GLU	ASP	CONFLICT	UNP E0UVR5

- Molecule 3 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



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

- Molecule 4 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 5 is a polymer of unknown type called SUGAR (5-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	5	Total	C	N	O	0	0
			61	34	2	25		
5	A	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 6 is a polymer of unknown type called SUGAR (5-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 7 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	3	Total	C	N	O	S	0	0
			50	25	2	22	1		

- Molecule 8 is SULFATE ION (three-letter code: SO<sub>4</sub>) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	O	S	0	0
			5	4	1		
8	A	1	Total	O	S	0	0
			5	4	1		
8	B	1	Total	O	S	0	0
			5	4	1		

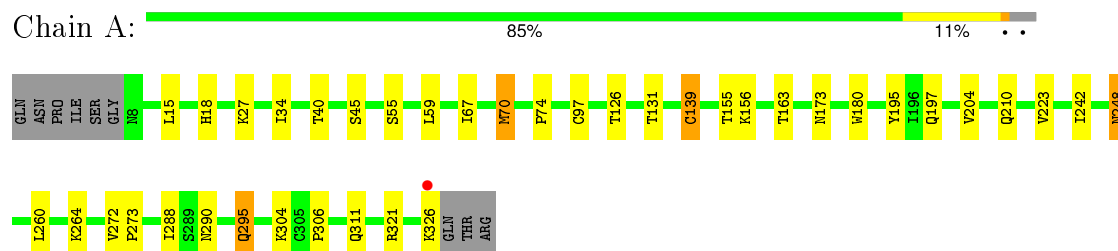
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	4	Total	O	0	0
			4	4		
9	B	3	Total	O	0	0
			3	3		

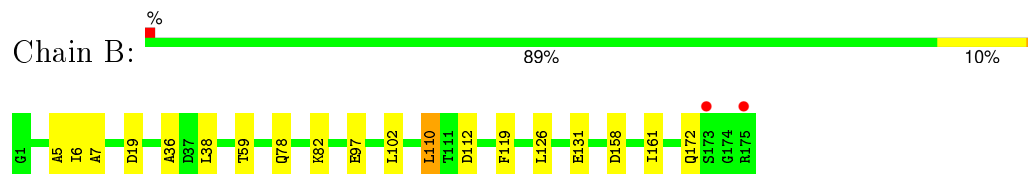
### 3 Residue-property plots

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 ( $RSRZ > 2$ ). 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 2: HA2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	97.29 Å 97.29 Å 348.27 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.69 – 3.00 48.64 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.69-3.00) 99.3 (48.64-3.00)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.56 (at 3.01 Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, $R_{free}$	0.207 , 0.251 0.216 , 0.207	Depositor DCC
$R_{free}$ test set	1050 reflections (5.41%)	DCC
Wilson B-factor (Å <sup>2</sup> )	64.2	Xtriage
Anisotropy	0.868	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 45.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	3 of 20455 reflections (0.015%)	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	4190	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.99% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: BMA, NAG, SIA, GAL, SO4, NGS, MAN

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	0.36	0/2519	0.58	0/3418
2	B	0.37	0/1450	0.61	1/1949 (0.1%)
All	All	0.36	0/3969	0.59	1/5367 (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
5	A	2	0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
2	B	110	LEU	CA-CB-CG	5.62	128.24	115.30

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	A	621	NAG	C1
5	A	625	MAN	C1

There are no planarity outliers.

## 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	2468	0	2412	13	0
2	B	1425	0	1344	7	0
3	A	14	0	13	0	0
4	A	28	0	25	0	0
5	A	122	0	104	2	0
6	A	61	0	52	0	0
7	A	50	0	40	2	0
8	A	10	0	0	0	0
8	B	5	0	0	0	0
9	A	4	0	0	0	0
9	B	3	0	0	0	0
All	All	4190	0	3990	22	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 (22) 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:A:290:ASN:OD1	1:A:304:LYS:O	2.16	0.64
1:A:15:LEU:HD22	2:B:119:PHE:HA	1.83	0.60
7:A:701:SIA:H113	7:A:701:SIA:O7	2.06	0.55
1:A:27:LYS:NZ	2:B:97:GLU:OE2	2.40	0.54
1:A:34:ILE:HD11	1:A:321:ARG:HD2	1.92	0.51
1:A:180:TRP:CE2	1:A:204:VAL:HG21	2.46	0.50
5:A:633:BMA:H61	5:A:637:MAN:H5	1.94	0.49
2:B:19:ASP:HB2	2:B:36:ALA:HB2	1.94	0.48
1:A:295:GLN:HB3	1:A:306:PRO:HB2	1.97	0.47
1:A:74:PRO:HG2	1:A:139:CYS:HB2	1.97	0.47
2:B:158:ASP:HB3	2:B:161:ILE:HD12	1.97	0.46
1:A:272:VAL:HG13	1:A:273:PRO:HD2	1.98	0.45
7:A:703:NGS:O3	7:A:703:NGS:HH3A	2.17	0.45
5:A:633:BMA:C6	5:A:637:MAN:H5	2.47	0.45
2:B:6:ILE:O	2:B:7:ALA:HB3	2.16	0.44
1:A:197:GLN:NE2	1:A:248:ASN:O	2.49	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:CYS:HA	1:A:139:CYS:HB3	2.00	0.44
1:A:67:ILE:O	1:A:70:MET:HG3	2.17	0.43
2:B:78:GLN:HE21	2:B:82:LYS:NZ	2.16	0.42
1:A:288:ILE:HD12	1:A:295:GLN:HG3	2.03	0.41
2:B:5:ALA:HB3	2:B:112:ASP:OD1	2.20	0.41
1:A:248:ASN:H	1:A:248:ASN:HD22	1.68	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	317/328 (97%)	296 (93%)	20 (6%)	1 (0%)	46	84
2	B	173/175 (99%)	162 (94%)	10 (6%)	1 (1%)	30	72
All	All	490/503 (97%)	458 (94%)	30 (6%)	2 (0%)	39	80

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	172	GLN
1	A	163	THR

### 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	279/289 (96%)	257 (92%)	22 (8%)	15	48
2	B	148/148 (100%)	142 (96%)	6 (4%)	37	76
All	All	427/437 (98%)	399 (93%)	28 (7%)	21	57

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	HIS
1	A	40	THR
1	A	45	SER
1	A	55	SER
1	A	59	LEU
1	A	70	MET
1	A	126	THR
1	A	131	THR
1	A	139	CYS
1	A	155	THR
1	A	156	LYS
1	A	173	ASN
1	A	195	TYR
1	A	210	GLN
1	A	223	VAL
1	A	242	ILE
1	A	248	ASN
1	A	260	LEU
1	A	264	LYS
1	A	295	GLN
1	A	311	GLN
1	A	326	LYS
2	B	38	LEU
2	B	59	THR
2	B	102	LEU
2	B	110	LEU
2	B	126	LEU
2	B	131	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	173	ASN
1	A	183	HIS

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Mol	Chain	Res	Type
1	A	248	ASN
1	A	290	ASN
1	A	311	GLN
2	B	53	ASN
2	B	65	GLN
2	B	78	GLN
2	B	125	GLN

### 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 ⓘ

20 carbohydrates 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	$\# Z  > 2$
4	NAG	A	611	1,4	14,14,15	0.80	0	15,19,21	1.25	1 (6%)
4	NAG	A	612	4	14,14,15	0.42	0	15,19,21	1.79	1 (6%)
5	NAG	A	621	1,5	14,14,15	0.51	0	15,19,21	0.85	0
5	NAG	A	622	5	14,14,15	0.53	0	15,19,21	0.99	1 (6%)
5	BMA	A	623	5	11,11,12	0.44	0	14,15,17	1.13	1 (7%)
5	MAN	A	624	5	11,11,12	0.54	0	14,15,17	1.16	1 (7%)
5	MAN	A	625	5	11,11,12	0.61	0	14,15,17	1.71	3 (21%)
5	NAG	A	631	1,5	14,14,15	0.50	0	15,19,21	1.86	3 (20%)
5	NAG	A	632	5	14,14,15	0.55	0	15,19,21	0.89	1 (6%)
5	BMA	A	633	5	11,11,12	0.55	0	14,15,17	1.39	1 (7%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	MAN	A	634	5	11,11,12	0.90	1 (9%)	14,15,17	2.20	3 (21%)
5	MAN	A	637	5	11,11,12	0.66	0	14,15,17	1.45	2 (14%)
6	NAG	A	641	1,6	14,14,15	0.69	0	15,19,21	0.92	0
6	NAG	A	642	6	14,14,15	0.51	0	15,19,21	0.63	0
6	BMA	A	643	6	11,11,12	0.42	0	14,15,17	1.20	2 (14%)
6	MAN	A	644	6	11,11,12	0.91	1 (9%)	14,15,17	1.94	4 (28%)
6	BMA	A	645	6	11,11,12	0.60	0	14,15,17	2.06	2 (14%)
7	SIA	A	701	7	16,20,21	0.32	0	18,28,31	1.21	2 (11%)
7	GAL	A	702	7	11,11,12	0.52	0	14,15,17	0.95	1 (7%)
7	NGS	A	703	7	19,19,19	0.79	1 (5%)	22,28,28	1.35	2 (9%)

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	611	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	612	4	-	0/6/23/26	0/1/1/1
5	NAG	A	621	1,5	1/1/5/7	0/6/23/26	0/1/1/1
5	NAG	A	622	5	-	0/6/23/26	0/1/1/1
5	BMA	A	623	5	-	0/2/19/22	0/1/1/1
5	MAN	A	624	5	-	0/2/19/22	1/1/1/1
5	MAN	A	625	5	1/1/4/5	0/2/19/22	0/1/1/1
5	NAG	A	631	1,5	-	0/6/23/26	0/1/1/1
5	NAG	A	632	5	-	0/6/23/26	0/1/1/1
5	BMA	A	633	5	-	0/2/19/22	0/1/1/1
5	MAN	A	634	5	-	0/2/19/22	0/1/1/1
5	MAN	A	637	5	-	0/2/19/22	0/1/1/1
6	NAG	A	641	1,6	-	0/6/23/26	0/1/1/1
6	NAG	A	642	6	-	0/6/23/26	0/1/1/1
6	BMA	A	643	6	-	0/2/19/22	0/1/1/1
6	MAN	A	644	6	-	0/2/19/22	0/1/1/1
6	BMA	A	645	6	-	0/2/19/22	0/1/1/1
7	SIA	A	701	7	-	0/14/34/38	0/1/1/1
7	GAL	A	702	7	-	0/2/19/22	0/1/1/1
7	NGS	A	703	7	-	0/10/30/30	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	634	MAN	C2-C3	2.18	1.55	1.52
6	A	644	MAN	C2-C3	2.30	1.55	1.52
7	A	703	NGS	O8-S	2.69	1.64	1.50

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	632	NAG	C4-C3-C2	-2.34	107.59	111.23
7	A	702	GAL	O5-C1-C2	-2.21	107.27	110.86
6	A	643	BMA	C1-O5-C5	2.03	114.82	112.25
5	A	631	NAG	C3-C4-C5	2.04	113.76	110.20
7	A	701	SIA	C5-N5-C10	2.11	128.51	123.10
7	A	703	NGS	O6-C6-C5	2.14	112.02	107.90
6	A	644	MAN	O2-C2-C3	2.18	114.51	110.12
5	A	625	MAN	C1-O5-C5	2.23	115.08	112.25
5	A	637	MAN	C1-O5-C5	2.28	115.14	112.25
5	A	634	MAN	C3-C4-C5	2.31	114.22	110.20
5	A	631	NAG	C8-C7-N2	2.41	120.72	116.11
5	A	623	BMA	C3-C4-C5	2.45	114.48	110.20
5	A	622	NAG	C4-C3-C2	2.61	115.29	111.23
7	A	701	SIA	C11-C10-N5	2.80	121.46	116.11
6	A	644	MAN	C1-C2-C3	2.98	113.06	109.54
7	A	703	NGS	CH3-C-N	3.05	121.94	116.11
6	A	644	MAN	C2-C3-C4	3.15	116.40	111.04
5	A	625	MAN	C1-C2-C3	3.16	113.28	109.54
5	A	637	MAN	C3-C4-C5	3.23	115.82	110.20
6	A	643	BMA	C1-C2-C3	3.27	113.41	109.54
5	A	634	MAN	C2-C3-C4	3.37	116.76	111.04
5	A	624	MAN	C1-O5-C5	3.60	116.82	112.25
4	A	611	NAG	C4-C3-C2	3.90	117.30	111.23
5	A	633	BMA	C1-C2-C3	4.25	114.57	109.54
6	A	644	MAN	C3-C4-C5	4.32	117.72	110.20
5	A	625	MAN	O5-C1-C2	4.65	118.41	110.86
6	A	645	BMA	C1-C2-C3	4.67	115.07	109.54
6	A	645	BMA	C1-O5-C5	5.13	118.75	112.25
5	A	631	NAG	C1-O5-C5	5.61	119.37	112.25
4	A	612	NAG	C1-O5-C5	6.43	120.41	112.25
5	A	634	MAN	C1-C2-C3	6.61	117.36	109.54

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	A	625	MAN	C1

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Mol	Chain	Res	Type	Atom
5	A	621	NAG	C1

There are no torsion outliers.

All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	624	MAN	C1-C2-C3-C4-C5-O5

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	633	BMA	2	0
5	A	637	MAN	2	0
7	A	701	SIA	1	0
7	A	703	NGS	1	0

## 5.6 Ligand geometry [i](#)

4 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	$\# Z  > 2$
8	SO4	A	1327	-	4,4,4	0.38	0	6,6,6	0.09	0
8	SO4	A	1328	-	4,4,4	0.42	0	6,6,6	0.09	0
3	NAG	A	601	1	14,14,15	0.50	0	15,19,21	1.47	3 (20%)
8	SO4	B	1176	-	4,4,4	0.38	0	6,6,6	0.11	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
8	SO4	A	1327	-	-	0/0/0/0	0/0/0/0
8	SO4	A	1328	-	-	0/0/0/0	0/0/0/0
3	NAG	A	601	1	-	0/6/23/26	0/1/1/1
8	SO4	B	1176	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	NAG	O5-C5-C6	2.11	111.91	107.35
3	A	601	NAG	C8-C7-N2	2.53	120.94	116.11
3	A	601	NAG	C2-N2-C7	2.95	126.83	123.04

There are no chirality outliers.

There are no torsion outliers.

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<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	319/328 (97%)	-0.26	1 (0%) 94 84	57, 80, 119, 144	0
2	B	175/175 (100%)	-0.36	2 (1%) 82 58	56, 75, 96, 129	0
All	All	494/503 (98%)	-0.30	3 (0%) 90 73	56, 79, 115, 144	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	326	LYS	3.8
2	B	173	SER	3.3
2	B	175	ARG	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. 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
6	NAG	A	641	14/15	0.90	0.28	3.49	105,113,119,128	0
6	BMA	A	645	11/12	0.86	0.25	1.20	118,127,134,138	0
5	NAG	A	632	14/15	0.94	0.18	-0.35	100,112,119,119	0
7	SIA	A	701	20/21	0.96	0.17	-0.75	93,101,106,108	0
6	BMA	A	643	11/12	0.69	0.52	-	154,163,167,168	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
7	NGS	A	703	19/19	0.90	0.14	-	127,134,145,148	0
5	NAG	A	631	14/15	0.95	0.14	-	110,113,116,118	0
5	BMA	A	623	11/12	0.70	0.37	-	141,152,158,158	0
6	MAN	A	644	11/12	0.81	0.28	-	140,148,151,154	0
7	GAL	A	702	11/12	0.94	0.15	-	108,115,124,124	0
4	NAG	A	611	14/15	0.91	0.19	-	110,117,129,136	0
5	MAN	A	637	11/12	0.78	0.39	-	153,164,167,167	0
5	BMA	A	633	11/12	0.83	0.30	-	124,131,137,138	0
5	MAN	A	634	11/12	0.71	0.43	-	130,139,141,142	0
5	NAG	A	622	14/15	0.73	0.28	-	127,137,145,153	0
5	NAG	A	621	14/15	0.88	0.27	-	92,103,117,130	0
6	NAG	A	642	14/15	0.86	0.38	-	123,140,146,153	0
4	NAG	A	612	14/15	0.82	0.34	-	138,144,152,154	0
5	MAN	A	624	11/12	0.80	0.53	-	142,155,158,158	0
5	MAN	A	625	11/12	0.77	0.59	-	144,153,156,156	0

## 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. 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( $\text{\AA}^2$ )	Q<0.9
8	SO4	B	1176	5/5	0.97	0.12	-3.51	81,84,92,92	0
3	NAG	A	601	14/15	0.79	0.28	-	104,118,122,123	0
8	SO4	A	1327	5/5	0.91	0.20	-	139,141,145,146	0
8	SO4	A	1328	5/5	0.80	0.31	-	123,132,135,135	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.