



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 4, 2023 – 11:35 AM EDT

PDB ID : 6PCX
Title : Crystal Structure of a H5N1 influenza virus hemagglutinin at pH 6.0
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Deposited on : 2019-06-18
Resolution : 2.11 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.35.1
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) : 2.35.1

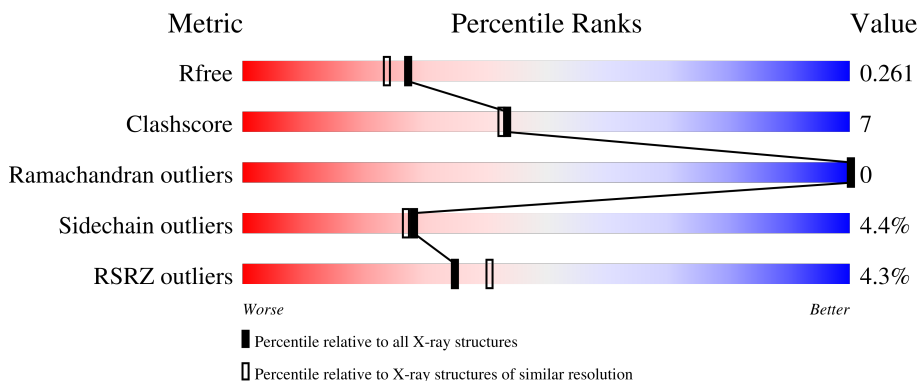
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.11 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}	130704	6241 (2.14-2.10)
Clashscore	141614	6778 (2.14-2.10)
Ramachandran outliers	138981	6705 (2.14-2.10)
Sidechain outliers	138945	6706 (2.14-2.10)
RSRZ outliers	127900	6112 (2.14-2.10)

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 $\leq 5\%$. 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	526	

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
2	SO4	A	1108	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	A	1110	-	-	-	X

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 4206 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
			Total	C	N	O	S			
1	A	491	3943	2476	682	761	24	0	2	0

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	LEU	-	expression tag	UNP Q1KHJ9
A	-3	LEU	-	expression tag	UNP Q1KHJ9
A	-2	ALA	-	expression tag	UNP Q1KHJ9
A	-1	ALA	-	expression tag	UNP Q1KHJ9
A	0	ALA	-	expression tag	UNP Q1KHJ9
A	1	ALA	-	expression tag	UNP Q1KHJ9
A	2	HIS	-	expression tag	UNP Q1KHJ9
A	3	SER	-	expression tag	UNP Q1KHJ9
A	4	ALA	-	expression tag	UNP Q1KHJ9
A	5	PHE	-	expression tag	UNP Q1KHJ9
A	6	ALA	-	expression tag	UNP Q1KHJ9
A	7	ALA	-	expression tag	UNP Q1KHJ9
A	8	ASP	-	expression tag	UNP Q1KHJ9
A	9	PRO	-	expression tag	UNP Q1KHJ9
A	10	GLY	-	expression tag	UNP Q1KHJ9
A	1042	SER	-	expression tag	UNP Q1KHJ9
A	1043	ARG	-	expression tag	UNP Q1KHJ9
A	1044	LEU	-	expression tag	UNP Q1KHJ9
A	1045	VAL	-	expression tag	UNP Q1KHJ9
A	1046	PRO	-	expression tag	UNP Q1KHJ9
A	1047	ARG	-	expression tag	UNP Q1KHJ9

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0

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



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

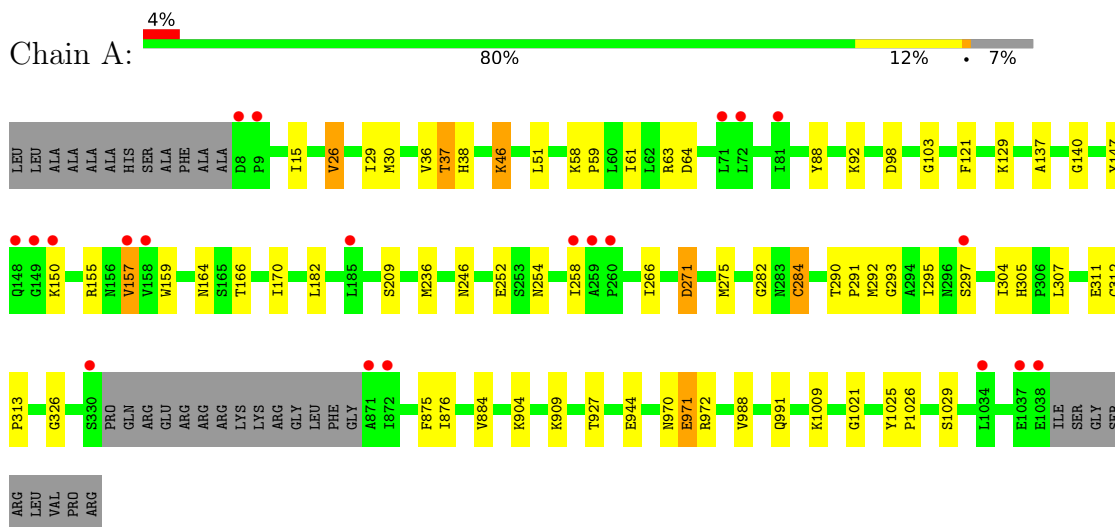
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	129	Total	O	0	0
			129	129		

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



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	108.54Å 108.54Å 419.79Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.44 – 2.11 29.42 – 2.11	Depositor EDS
% Data completeness (in resolution range)	98.1 (29.44-2.11) 98.1 (29.42-2.11)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.67 (at 2.10Å)	Xtrriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.207 , 0.255 0.213 , 0.261	Depositor DCC
R_{free} test set	2763 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	54.9	Xtrriage
Anisotropy	0.140	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 54.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4206	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.48% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, NAG, GOL

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	0.63	0/4037	0.81	0/5461

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	3943	0	3794	49	0
2	A	60	0	0	5	0
3	A	18	0	24	3	0
4	A	56	0	52	6	0
5	A	129	0	0	2	0
All	All	4206	0	3870	53	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 7.

All (53) 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:THR:HG22	1:A:292:MET:H	1.33	0.92
1:A:290:THR:HB	1:A:293:GLY:O	1.72	0.90
1:A:98:ASP:OD2	3:A:1115:GOL:H11	1.72	0.89
1:A:59:PRO:HG2	1:A:61:ILE:HD11	1.63	0.80
1:A:971:GLU:OE1	1:A:972:ARG:NH2	2.18	0.76
1:A:37:THR:HG22	1:A:326:GLY:HA3	1.67	0.74
1:A:164:ASN:HD21	4:A:1119:NAG:H61	1.58	0.69
1:A:295:ILE:HD11	1:A:304:ILE:HD12	1.79	0.64
1:A:98:ASP:CG	3:A:1115:GOL:H11	2.19	0.63
1:A:147:TYR:CD2	1:A:155:ARG:NH2	2.70	0.59
1:A:291:PRO:HD3	1:A:307:LEU:O	2.02	0.59
1:A:1009[A]:LYS:HA	1:A:1009[A]:LYS:HE2	1.85	0.59
1:A:297:SER:CB	1:A:313:PRO:HD3	2.37	0.54
1:A:63:ARG:O	1:A:92:LYS:HD3	2.08	0.53
1:A:103:GLY:HA3	1:A:236:MET:O	2.09	0.53
1:A:290:THR:HG23	1:A:305:HIS:HB3	1.92	0.52
1:A:311:GLU:HB3	1:A:927:THR:HG22	1.90	0.52
1:A:297:SER:HB2	1:A:313:PRO:HD3	1.93	0.51
1:A:312:CYS:O	2:A:1106:SO4:O4	2.28	0.51
1:A:295:ILE:HD11	1:A:304:ILE:CD1	2.42	0.50
1:A:209:SER:OG	1:A:252:GLU:HB3	2.12	0.50
1:A:121:PHE:HD1	1:A:266:ILE:HG22	1.76	0.50
4:A:1119:NAG:C8	5:A:1305:HOH:O	2.60	0.49
1:A:37:THR:HG23	2:A:1108:SO4:S	2.54	0.46
1:A:1025:TYR:N	1:A:1026:PRO:CD	2.78	0.46
2:A:1112:SO4:O3	4:A:1117:NAG:O4	2.33	0.46
1:A:46:LYS:HE2	1:A:46:LYS:HA	1.96	0.46
4:A:1119:NAG:H82	5:A:1305:HOH:O	2.15	0.46
1:A:157:VAL:HG13	1:A:258:ILE:HG22	1.98	0.46
1:A:170:ILE:O	1:A:252:GLU:HA	2.17	0.45
1:A:59:PRO:HB3	1:A:88:TYR:CZ	2.52	0.44
1:A:311:GLU:HB3	1:A:927:THR:CG2	2.48	0.44
1:A:15:ILE:HD11	1:A:988:VAL:HG21	2.00	0.44
1:A:155:ARG:HD3	1:A:155:ARG:HA	1.77	0.44
1:A:129:LYS:HE2	1:A:137:ALA:O	2.18	0.43
1:A:38:HIS:HB2	2:A:1108:SO4:O1	2.19	0.43
1:A:98:ASP:OD1	3:A:1115:GOL:H11	2.17	0.43
1:A:290:THR:HG21	1:A:304:ILE:HG22	2.01	0.43
1:A:971:GLU:OE1	1:A:972:ARG:CZ	2.66	0.42
1:A:140:GLY:HA3	1:A:159:TRP:O	2.20	0.42
1:A:271:ASP:OD1	1:A:271:ASP:N	2.52	0.42
1:A:991:GLN:OE1	1:A:1021:GLY:HA2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:121:PHE:CD1	1:A:266:ILE:HG22	2.53	0.42
1:A:51:LEU:HD22	1:A:275:MET:CE	2.49	0.42
2:A:1111:SO4:S	4:A:1118:NAG:O4	2.72	0.41
1:A:1025:TYR:HB3	1:A:1026:PRO:HD3	2.01	0.41
1:A:875:PHE:CE1	1:A:876:ILE:HG13	2.55	0.41
1:A:29:ILE:HG23	1:A:30:MET:HG3	2.02	0.41
1:A:140:GLY:HA3	1:A:159:TRP:HB3	2.03	0.41
1:A:164:ASN:ND2	4:A:1119:NAG:H61	2.29	0.40
1:A:282:GLY:C	1:A:284:CYS:H	2.25	0.40
1:A:26:VAL:HG13	1:A:970:ASN:ND2	2.36	0.40
1:A:182:LEU:HD23	1:A:182:LEU:HA	1.86	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	489/526 (93%)	461 (94%)	28 (6%)	0	100 100

There are no Ramachandran outliers to report.

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	437/460 (95%)	418 (96%)	19 (4%)	29 28

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	VAL
1	A	36	VAL
1	A	37	THR
1	A	46	LYS
1	A	58	LYS
1	A	64	ASP
1	A	150	LYS
1	A	157	VAL
1	A	166	THR
1	A	246	ASN
1	A	254	ASN
1	A	271	ASP
1	A	284	CYS
1	A	884	VAL
1	A	904	LYS
1	A	909	LYS
1	A	944	GLU
1	A	971	GLU
1	A	1029	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry

19 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	GOL	A	1114	-	5,5,5	0.12	0	5,5,5	0.36	0
2	SO4	A	1112	-	4,4,4	0.38	0	6,6,6	0.04	0
4	NAG	A	1118	1	14,14,15	0.37	0	17,19,21	1.57	2 (11%)
2	SO4	A	1110	-	4,4,4	0.37	0	6,6,6	0.05	0
2	SO4	A	1104	-	4,4,4	0.37	0	6,6,6	0.08	0
2	SO4	A	1108	-	4,4,4	0.39	0	6,6,6	0.05	0
2	SO4	A	1105	-	4,4,4	0.38	0	6,6,6	0.07	0
2	SO4	A	1109	-	4,4,4	0.37	0	6,6,6	0.05	0
2	SO4	A	1103	-	4,4,4	0.38	0	6,6,6	0.12	0
2	SO4	A	1101	-	4,4,4	0.32	0	6,6,6	0.07	0
2	SO4	A	1106	-	4,4,4	0.35	0	6,6,6	0.08	0
4	NAG	A	1119	1	14,14,15	0.49	0	17,19,21	2.32	4 (23%)
2	SO4	A	1111	-	4,4,4	0.37	0	6,6,6	0.06	0
3	GOL	A	1115	-	5,5,5	0.09	0	5,5,5	0.39	0
4	NAG	A	1116	1	14,14,15	0.43	0	17,19,21	2.18	7 (41%)
4	NAG	A	1117	1	14,14,15	0.37	0	17,19,21	1.24	2 (11%)
3	GOL	A	1113	-	5,5,5	0.10	0	5,5,5	0.34	0
2	SO4	A	1102	-	4,4,4	0.36	0	6,6,6	0.07	0
2	SO4	A	1107	-	4,4,4	0.37	0	6,6,6	0.09	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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	A	1119	1	-	0/6/23/26	0/1/1/1
3	GOL	A	1114	-	-	0/4/4/4	-
3	GOL	A	1115	-	-	2/4/4/4	-
4	NAG	A	1118	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	A	1116	1	-	3/6/23/26	0/1/1/1
4	NAG	A	1117	1	-	0/6/23/26	0/1/1/1
3	GOL	A	1113	-	-	2/4/4/4	-

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1119	NAG	C1-C2-N2	7.03	122.50	110.49
4	A	1118	NAG	O5-C1-C2	4.89	119.01	111.29
4	A	1116	NAG	C1-O5-C5	4.43	118.19	112.19
4	A	1119	NAG	O5-C5-C6	4.31	113.96	107.20
4	A	1116	NAG	C2-N2-C7	3.85	128.38	122.90
4	A	1116	NAG	C1-C2-N2	3.62	116.67	110.49
4	A	1116	NAG	O5-C1-C2	3.51	116.83	111.29
4	A	1117	NAG	O5-C1-C2	2.86	115.81	111.29
4	A	1119	NAG	O5-C1-C2	2.72	115.58	111.29
4	A	1116	NAG	C3-C4-C5	2.20	114.17	110.24
4	A	1117	NAG	C3-C4-C5	-2.19	106.33	110.24
4	A	1118	NAG	C1-O5-C5	2.19	115.16	112.19
4	A	1119	NAG	C1-O5-C5	2.12	115.07	112.19
4	A	1116	NAG	C6-C5-C4	-2.05	108.21	113.00
4	A	1116	NAG	O7-C7-N2	2.04	125.70	121.95

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1115	GOL	O1-C1-C2-C3
4	A	1116	NAG	C4-C5-C6-O6
3	A	1113	GOL	O1-C1-C2-O2
3	A	1113	GOL	O1-C1-C2-C3
3	A	1115	GOL	O1-C1-C2-O2
4	A	1116	NAG	O5-C5-C6-O6
4	A	1116	NAG	C3-C2-N2-C7

There are no ring outliers.

8 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1112	SO4	1	0
4	A	1118	NAG	1	0
2	A	1108	SO4	2	0
2	A	1106	SO4	1	0
4	A	1119	NAG	4	0
2	A	1111	SO4	1	0
3	A	1115	GOL	3	0
4	A	1117	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, 95th 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 ²)	Q<0.9
1	A	491/526 (93%)	0.01	21 (4%) 35 41	49, 62, 92, 130	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	8	ASP	4.4
1	A	871	ALA	4.2
1	A	157	VAL	3.6
1	A	297	SER	3.6
1	A	1034	LEU	3.4
1	A	185	LEU	3.3
1	A	72	LEU	3.3
1	A	9	PRO	3.1
1	A	71	LEU	3.1
1	A	258	ILE	3.0
1	A	259	ALA	2.8
1	A	330	SER	2.7
1	A	1037	GLU	2.7
1	A	149	GLY	2.6
1	A	260	PRO	2.6
1	A	1038	GLU	2.5
1	A	158	VAL	2.5
1	A	150	LYS	2.3
1	A	148	GLN	2.3
1	A	81	ILE	2.1
1	A	872	ILE	2.1

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 monosaccharides in this entry.

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, 95th 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	B-factors(Å ²)	Q<0.9
2	SO4	A	1110	5/5	0.44	0.50	185,185,193,194	5
4	NAG	A	1119	14/15	0.64	0.36	105,140,148,152	0
2	SO4	A	1112	5/5	0.70	0.36	148,155,165,171	0
2	SO4	A	1106	5/5	0.70	0.24	119,140,157,171	0
3	GOL	A	1114	6/6	0.71	0.34	87,97,105,111	0
2	SO4	A	1111	5/5	0.74	0.38	136,152,156,158	0
4	NAG	A	1116	14/15	0.83	0.37	96,117,127,129	0
2	SO4	A	1101	5/5	0.88	0.14	90,100,128,134	0
2	SO4	A	1107	5/5	0.88	0.29	115,125,138,139	0
3	GOL	A	1113	6/6	0.90	0.25	107,110,115,129	0
3	GOL	A	1115	6/6	0.93	0.12	88,90,91,92	0
2	SO4	A	1105	5/5	0.93	0.20	95,118,120,136	0
4	NAG	A	1118	14/15	0.93	0.29	87,98,110,111	0
2	SO4	A	1104	5/5	0.93	0.15	107,115,122,135	0
2	SO4	A	1109	5/5	0.94	0.18	86,91,104,104	5
2	SO4	A	1102	5/5	0.95	0.17	94,94,99,105	0
2	SO4	A	1108	5/5	0.95	0.14	125,135,141,155	0
4	NAG	A	1117	14/15	0.96	0.12	67,72,78,82	0
2	SO4	A	1103	5/5	0.99	0.08	65,69,79,80	0

6.5 Other polymers [i](#)

There are no such residues in this entry.