



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:56 PM GMT

PDB ID : 1DBO  
Title : CRYSTAL STRUCTURE OF CHONDROITINASE B  
Authors : Huang, W.; Matte, A.; Li, Y.; Kim, Y.S.; Linhardt, R.J.; Su, H.; Cygler, M.  
Deposited on : 1999-11-03  
Resolution : 1.70 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

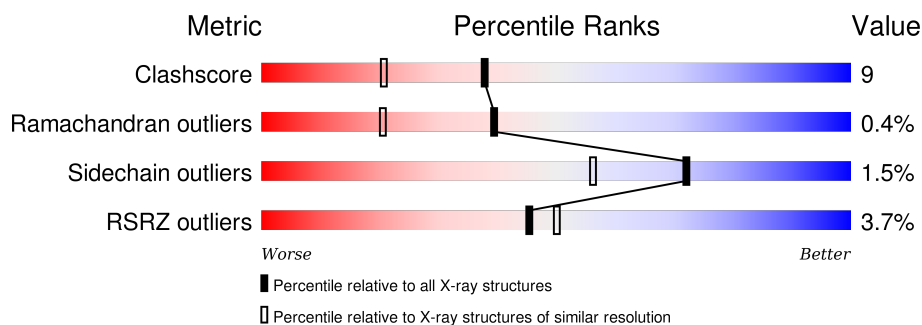
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.70 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))
Clashscore	102246	3585 (1.70-1.70)
Ramachandran outliers	100387	3527 (1.70-1.70)
Sidechain outliers	100360	3527 (1.70-1.70)
RSRZ outliers	91569	3200 (1.70-1.70)

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	506	

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	GCU	A	509	-	-	-	X
2	MXV	A	511	X	-	-	-
3	GC4	A	516	X	-	-	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4365 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 CHONDROITINASE B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	481	Total	C	N	O	S	0	1	0
			3787	2413	662	695	17			

- Molecule 2 is a polymer of unknown type called SUGAR (7-MER).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	7	Total	C	O	0	0
			73	41	32		

- Molecule 3 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	2	Total	C	N	O	S	0	0
			30	14	1	14	1		

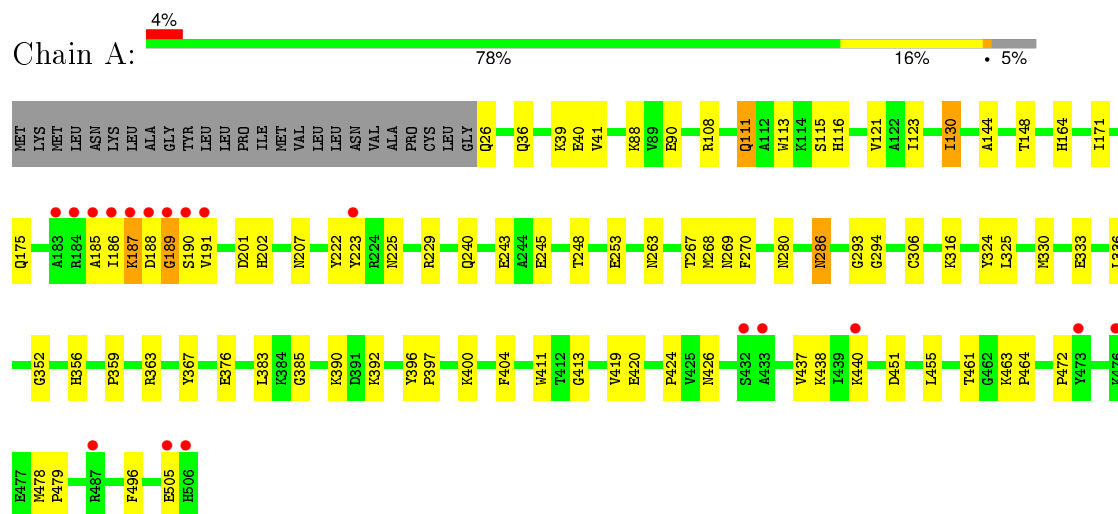
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	475	Total	O	0	0
			475	475		

### 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 ( $RSRZ > 2$ ). 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: CHONDROITINASE B



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	50.60Å 74.40Å 58.70Å 90.00° 92.90° 90.00°	Depositor
Resolution (Å)	19.95 – 1.70 19.95 – 1.69	Depositor EDS
% Data completeness (in resolution range)	(Not available) (19.95-1.70) 91.5 (19.95-1.69)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.03 (at 1.69Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.181 , 0.217 0.189 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	10.7	Xtriage
Anisotropy	0.069	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 43.6	EDS
Estimated twinning fraction	0.032 for h,-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 44133 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4365	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	12.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.55% 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: XYP, BGC, G4D, MXY, RAM, GCU, GC4, ASG, 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	0.35	0/3882	0.73	4/5249 (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
1	A	0	1
2	A	1	0
3	A	1	0
All	All	2	1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	148	THR	N-CA-C	-6.06	94.64	111.00
1	A	189	GLY	N-CA-C	-5.80	98.60	113.10
1	A	270	PHE	N-CA-C	-5.74	95.49	111.00
1	A	26	GLN	CB-CG-CD	-5.09	98.36	111.60

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	511	MXY	C1
3	A	516	GC4	C5

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	293	GLY	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	3787	0	3719	71	0
2	A	73	0	57	1	0
3	A	30	0	19	1	0
4	A	475	0	0	13	1
All	All	4365	0	3795	71	1

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 9.

All (71) 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:187:LYS:HG3	1:A:188:ASP:H	1.08	1.18
1:A:280:ASN:HD21	1:A:461:THR:H	1.12	0.94
1:A:111:GLN:H	1:A:111:GLN:HE21	1.10	0.92
1:A:207:ASN:HD22	1:A:240:GLN:HE22	1.25	0.84
1:A:187:LYS:HG3	1:A:188:ASP:N	1.92	0.82
1:A:111:GLN:N	1:A:111:GLN:HE21	1.78	0.81
1:A:189:GLY:O	1:A:191:VAL:N	2.15	0.80
1:A:111:GLN:H	1:A:111:GLN:NE2	1.82	0.78
1:A:39:LYS:HE3	4:A:917:HOH:O	1.84	0.78
1:A:187:LYS:HE2	4:A:811:HOH:O	1.86	0.74
1:A:40:GLU:HB2	4:A:962:HOH:O	1.87	0.73
1:A:201:ASP:OD2	1:A:202:HIS:HD2	1.76	0.69
1:A:202:HIS:HE1	2:A:507:MAN:O6	1.78	0.66
1:A:121:VAL:HG11	1:A:130[A]:ILE:HD12	1.79	0.64
1:A:187:LYS:CG	1:A:188:ASP:H	1.85	0.58
1:A:397:PRO:HB3	1:A:420:GLU:HG3	1.85	0.58
1:A:245:GLU:HG2	1:A:269:ASN:ND2	2.19	0.58
1:A:363:ARG:HG3	4:A:704:HOH:O	2.04	0.58
1:A:187:LYS:HG2	1:A:223:TYR:CG	2.40	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:GLN:O	1:A:40:GLU:HG3	2.05	0.57
1:A:263:ASN:HD21	1:A:286:ASN:ND2	2.03	0.56
1:A:463:LYS:HB2	1:A:464:PRO:CD	2.35	0.56
1:A:248:THR:HG23	1:A:248:THR:O	2.05	0.56
1:A:280:ASN:HA	1:A:306:CYS:O	2.06	0.56
1:A:404:PHE:CE2	1:A:424:PRO:HG2	2.41	0.55
1:A:223:TYR:HB2	4:A:974:HOH:O	2.06	0.55
1:A:363:ARG:NH1	3:A:516:GC4:H41	2.21	0.55
1:A:115:SER:O	1:A:116:HIS:HB2	2.06	0.55
1:A:333:GLU:HB3	1:A:367:TYR:OH	2.05	0.55
1:A:330:MET:HE2	4:A:743:HOH:O	2.06	0.55
1:A:187:LYS:CG	1:A:188:ASP:N	2.62	0.54
1:A:267:THR:HG22	1:A:294:GLY:HA3	1.89	0.54
1:A:189:GLY:O	1:A:191:VAL:HG22	2.07	0.54
1:A:164:HIS:HA	1:A:202:HIS:O	2.07	0.54
1:A:222:TYR:HB3	1:A:223:TYR:CE1	2.44	0.53
1:A:229:ARG:HA	1:A:253:GLU:O	2.08	0.53
1:A:438:LYS:HB3	1:A:440:LYS:HE3	1.91	0.52
1:A:330:MET:CE	4:A:743:HOH:O	2.59	0.51
1:A:356:HIS:NE2	1:A:400:LYS:HD3	2.25	0.51
1:A:330:MET:HE3	4:A:884:HOH:O	2.13	0.48
1:A:123:ILE:HD11	1:A:130[A]:ILE:HG13	1.95	0.48
1:A:437:VAL:HG12	1:A:438:LYS:N	2.29	0.48
1:A:263:ASN:HD21	1:A:286:ASN:HD21	1.61	0.47
1:A:286:ASN:HD22	1:A:286:ASN:C	2.17	0.47
1:A:390:LYS:HE2	1:A:392:LYS:O	2.15	0.45
1:A:383:LEU:HD22	1:A:411:TRP:CH2	2.51	0.45
1:A:352:GLY:HA2	1:A:396:TYR:CG	2.52	0.45
1:A:451:ASP:O	1:A:455:LEU:HG	2.17	0.45
1:A:185:ALA:O	1:A:186:ILE:C	2.55	0.45
1:A:352:GLY:HA2	1:A:396:TYR:CD1	2.53	0.44
1:A:419:VAL:HG22	1:A:420:GLU:N	2.32	0.44
1:A:171:ILE:HB	1:A:496:PHE:CD1	2.53	0.44
1:A:41:VAL:HA	4:A:934:HOH:O	2.16	0.44
1:A:187:LYS:HG2	1:A:223:TYR:CD1	2.53	0.43
1:A:88:LYS:HE3	1:A:90:GLU:OE1	2.18	0.43
1:A:207:ASN:HD22	1:A:240:GLN:NE2	2.05	0.43
1:A:316:LYS:HG3	4:A:991:HOH:O	2.18	0.43
1:A:383:LEU:HD22	1:A:411:TRP:CZ3	2.55	0.42
1:A:478:MET:HA	1:A:479:PRO:HD3	1.90	0.42
1:A:144:ALA:HA	1:A:175:GLN:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:GLY:C	1:A:191:VAL:N	2.73	0.42
1:A:225:ASN:HB2	4:A:974:HOH:O	2.19	0.41
1:A:336:LEU:HD22	1:A:376:GLU:HB2	2.01	0.41
1:A:505:GLU:HG3	4:A:885:HOH:O	2.21	0.41
1:A:437:VAL:HG13	4:A:979:HOH:O	2.20	0.41
1:A:108:ARG:HD2	1:A:113:TRP:CE2	2.56	0.41
1:A:164:HIS:O	1:A:472:PRO:HG3	2.21	0.40
1:A:324:TYR:HB3	1:A:359:PRO:HG3	2.03	0.40
1:A:325:LEU:O	1:A:359:PRO:HD2	2.22	0.40
1:A:385:GLY:HA2	1:A:413:GLY:O	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:710:HOH:O	4:A:781:HOH:O[2_656]	2.15	0.05

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	480/506 (95%)	450 (94%)	28 (6%)	2 (0%)	39 20

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	190	SER
1	A	187	LYS

### 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	398/419 (95%)	391 (98%)	7 (2%)	66 49

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	111	GLN
1	A	130[A]	ILE
1	A	130[B]	ILE
1	A	243	GLU
1	A	268	MET
1	A	286	ASN
1	A	426	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	GLN
1	A	111	GLN
1	A	202	HIS
1	A	213	ASN
1	A	240	GLN
1	A	280	ASN
1	A	286	ASN
1	A	320	ASN
1	A	379	HIS
1	A	426	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 ⓘ

9 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	$\# Z  > 2$
2	MAN	A	507	1,2	11,11,12	0.52	0	14,15,17	0.57	0
2	RAM	A	508	2	10,10,11	1.43	1 (10%)	14,14,16	1.27	2 (14%)
2	GCU	A	509	2	9,12,13	1.35	1 (11%)	13,17,19	1.00	1 (7%)
2	XYP	A	510	2	9,9,10	1.61	2 (22%)	12,12,14	2.04	4 (33%)
2	MXY	A	511	2	10,10,12	1.07	1 (10%)	12,13,17	1.14	2 (16%)
2	G4D	A	512	2	10,10,11	0.61	0	12,13,15	0.53	0
2	BGC	A	513	2	11,11,12	0.38	0	14,15,17	0.56	0
3	ASG	A	514	3	19,19,19	0.98	1 (5%)	21,28,28	1.35	2 (9%)
3	GC4	A	516	3	8,11,12	1.12	1 (12%)	8,15,17	2.43	4 (50%)

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
2	MAN	A	507	1,2	-	0/2/19/22	0/1/1/1
2	RAM	A	508	2	-	0/0/17/20	0/1/1/1
2	GCU	A	509	2	-	0/0/21/24	0/1/1/1
2	XYP	A	510	2	-	0/0/14/17	0/1/1/1
2	MXY	A	511	2	1/1/3/5	0/2/16/22	0/1/1/1
2	G4D	A	512	2	-	0/2/15/18	0/1/1/1
2	BGC	A	513	2	-	0/2/19/22	0/1/1/1
3	ASG	A	514	3	-	0/11/31/31	0/1/1/1
3	GC4	A	516	3	1/1/4/5	0/0/17/20	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	514	ASG	O4-S	-3.40	1.46	1.57
2	A	510	XYP	O5B-C5B	-3.28	1.36	1.42
2	A	510	XYP	O5B-C1B	-2.23	1.38	1.42
2	A	511	MXY	C4-C3	2.04	1.55	1.52
3	A	516	GC4	C3-C2	2.14	1.55	1.52
2	A	508	RAM	C4-C5	2.71	1.58	1.52
2	A	509	GCU	O5-C5	2.83	1.46	1.43

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	516	GC4	O3-C3-C4	-4.31	99.50	110.06
2	A	508	RAM	C1-C2-C3	-3.29	105.65	109.54
2	A	510	XYP	O3B-C3B-C2B	-2.13	106.15	110.00
2	A	509	GCU	C1-C2-C3	-2.08	107.08	109.54
2	A	511	MXY	C1-C2-C3	-2.04	106.36	109.39
3	A	514	ASG	O7-C7-C8	-2.03	118.33	122.06
3	A	516	GC4	C4-C3-C2	2.08	112.98	110.23
2	A	511	MXY	C5-O5-C1	2.19	115.11	111.57
2	A	508	RAM	O5-C5-C6	2.31	109.94	106.13
2	A	510	XYP	C5B-C4B-C3B	2.51	112.52	109.54
3	A	516	GC4	C1-O5-C5	2.81	118.03	113.47
2	A	510	XYP	O5B-C1B-C2B	2.99	115.89	110.31
3	A	516	GC4	O5-C5-C4	3.85	116.55	110.56
2	A	510	XYP	O5B-C5B-C4B	4.21	118.18	110.31
3	A	514	ASG	C4-O4-S	5.08	128.44	118.77

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	516	GC4	C5
2	A	511	MXY	C1

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	507	MAN	1	0
3	A	516	GC4	1	0

## 5.6 Ligand geometry

There are no ligands in this entry.

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	481/506 (95%)	0.24	18 (3%) 45 50	5, 9, 21, 42	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	189	GLY	8.9
1	A	188	ASP	8.3
1	A	186	ILE	7.1
1	A	191	VAL	5.3
1	A	185	ALA	5.1
1	A	187	LYS	4.9
1	A	223	TYR	3.8
1	A	432	SER	3.6
1	A	433	ALA	3.4
1	A	190	SER	3.2
1	A	487	ARG	3.0
1	A	440	LYS	2.8
1	A	505	GLU	2.8
1	A	506	HIS	2.5
1	A	476	LYS	2.4
1	A	184	ARG	2.2
1	A	183	ALA	2.2
1	A	473	TYR	2.2

### 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
2	GCU	A	509	12/13	0.95	0.10	3.18	9,11,13,14	0
2	RAM	A	508	10/11	0.94	0.11	0.85	7,8,9,9	0
3	ASG	A	514	19/19	0.95	0.10	-0.29	10,13,16,17	0
2	MAN	A	507	11/12	0.96	0.08	-0.85	7,8,9,9	0
2	BGC	A	513	11/12	0.68	0.23	-	25,28,29,29	0
2	XYP	A	510	9/10	0.91	0.17	-	15,18,19,20	0
3	GC4	A	516	11/12	0.92	0.14	-	17,18,19,20	0
2	G4D	A	512	10/11	0.82	0.27	-	26,28,30,30	0
2	MXY	A	511	10/12	0.88	0.12	-	20,21,23,24	0

### 6.4 Ligands [i](#)

There are no ligands in this entry.

### 6.5 Other polymers [i](#)

There are no such residues in this entry.