



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 3, 2018 – 11:41 PM EST

PDB ID : 6EJI  
Title : Structure of a glycosyltransferase  
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Deposited on : 2017-09-21  
Resolution : 2.30 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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20030736  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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-20030736

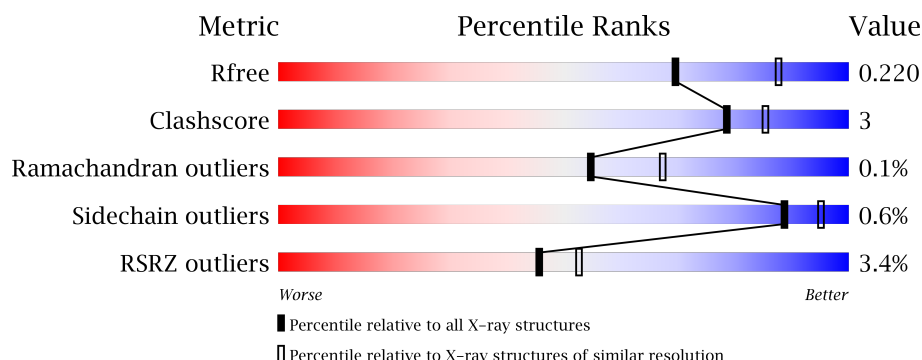
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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))
$R_{free}$	100719	4130 (2.30-2.30)
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (2.30-2.30)

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	373	<div> <div>2%</div> <div>89%</div> <div>8%</div> <div>.</div> </div>
1	B	373	<div> <div>5%</div> <div>87%</div> <div>9%</div> <div>.</div> </div>

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	UD2	A	401	X	-	-	-
4	GOL	B	403	-	-	-	X
5	NA	B	404	-	-	-	X

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 6130 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 WlaC protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	360	Total	C	N	O	S	Se	0	0	0
			2914	1886	486	529	9	4			
1	B	360	Total	C	N	O	S	Se	0	0	0
			2914	1886	486	529	9	4			

There are 30 discrepancies between the modelled and reference sequences:

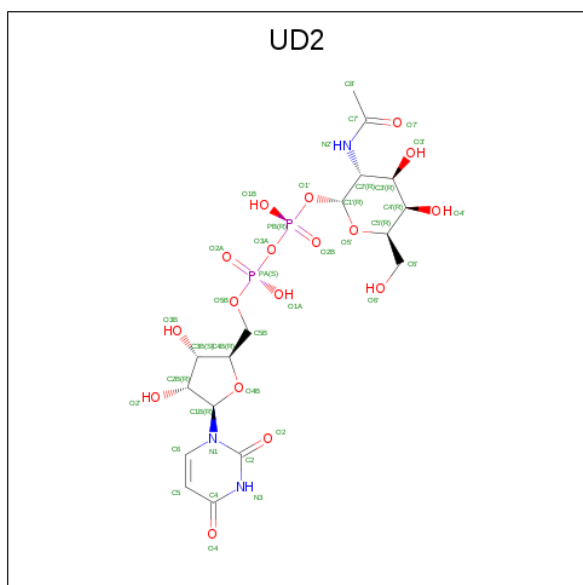
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	initiating methionine	UNP O86151
A	2	MSE	-	expression tag	UNP O86151
A	361	LEU	-	expression tag	UNP O86151
A	362	GLY	-	expression tag	UNP O86151
A	363	SER	-	expression tag	UNP O86151
A	364	HIS	-	expression tag	UNP O86151
A	365	HIS	-	expression tag	UNP O86151
A	366	HIS	-	expression tag	UNP O86151
A	367	HIS	-	expression tag	UNP O86151
A	368	HIS	-	expression tag	UNP O86151
A	369	HIS	-	expression tag	UNP O86151
A	370	HIS	-	expression tag	UNP O86151
A	371	HIS	-	expression tag	UNP O86151
A	372	HIS	-	expression tag	UNP O86151
A	373	HIS	-	expression tag	UNP O86151
B	1	MSE	-	initiating methionine	UNP O86151
B	2	MSE	-	expression tag	UNP O86151
B	361	LEU	-	expression tag	UNP O86151
B	362	GLY	-	expression tag	UNP O86151
B	363	SER	-	expression tag	UNP O86151
B	364	HIS	-	expression tag	UNP O86151
B	365	HIS	-	expression tag	UNP O86151
B	366	HIS	-	expression tag	UNP O86151
B	367	HIS	-	expression tag	UNP O86151
B	368	HIS	-	expression tag	UNP O86151

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Chain	Residue	Modelled	Actual	Comment	Reference
B	369	HIS	-	expression tag	UNP O86151
B	370	HIS	-	expression tag	UNP O86151
B	371	HIS	-	expression tag	UNP O86151
B	372	HIS	-	expression tag	UNP O86151
B	373	HIS	-	expression tag	UNP O86151

- Molecule 2 is URIDINE-DIPHOSPHATE-N-ACETYLGLACTOSAMINE (three-letter code: UD2) (formula:  $C_{17}H_{27}N_3O_{17}P_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			39	17	3	17	2		
2	B	1	Total	C	N	O	P	0	0
			39	17	3	17	2		

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	K	0	0
			1	1		
3	A	1	Total	K	0	0
			1	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



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

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	2	Total	Na	0	0
			2	2		
5	A	2	Total	Na	0	0
			2	2		

- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Cl	0	0
			1	1		
6	A	2	Total	Cl	0	0
			2	2		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	92	Total	O	0	0
			92	92		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	111	Total 111	O 111	0	0



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	61.62Å 127.03Å 71.58Å 90.00° 90.13° 90.00°	Depositor
Resolution (Å)	29.94 – 2.30 29.94 – 1.96	Depositor EDS
% Data completeness (in resolution range)	98.2 (29.94-2.30) 80.6 (29.94-1.96)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.10 (at 1.95Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155)	Depositor
R, $R_{free}$	0.196 , 0.223 0.197 , 0.220	Depositor DCC
$R_{free}$ test set	2389 reflections (4.98%)	DCC
Wilson B-factor (Å <sup>2</sup> )	35.8	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 37.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.118 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6130	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.62% 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: UD2, GOL, K, NA, CL

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	0.34	0/2965	0.52	1/3983 (0.0%)
1	B	0.31	0/2965	0.49	1/3983 (0.0%)
All	All	0.33	0/5930	0.51	2/7966 (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	1	0
1	B	1	0
All	All	2	0

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	290	ASN	N-CA-C	8.66	134.37	111.00
1	B	290	ASN	N-CA-C	7.47	131.16	111.00

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	290	ASN	CA
1	B	290	ASN	CA

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	2914	0	2992	19	0
1	B	2914	0	2992	25	0
2	A	39	0	21	0	0
2	B	39	0	21	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	6	0	8	0	0
4	B	6	0	8	1	0
5	A	2	0	0	0	0
5	B	2	0	0	0	0
6	A	2	0	0	0	0
6	B	1	0	0	0	0
7	A	92	0	0	2	0
7	B	111	0	0	2	0
All	All	6130	0	6042	40	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 (40) 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:248:GLU:CD	1:B:61:ARG:HH22	1.79	0.85
1:A:248:GLU:CD	1:B:61:ARG:NH2	2.39	0.76
1:B:38:LYS:NZ	1:B:42:GLY:O	2.28	0.67
1:B:125:LYS:NZ	4:B:403:GOL:O3	2.30	0.64
1:B:266:VAL:HA	7:B:503:HOH:O	1.98	0.63
1:A:20:VAL:HG23	7:A:515:HOH:O	2.00	0.61
1:A:248:GLU:OE1	1:B:61:ARG:NH2	2.34	0.60
1:B:51:GLU:OE1	1:B:51:GLU:N	2.33	0.59
1:B:40:HIS:HA	1:B:59:GLN:OE1	2.02	0.58
1:B:169:HIS:HA	1:B:343:ILE:HD11	1.87	0.57
1:A:81:LYS:O	1:A:85:GLU:HG3	2.05	0.57
1:A:266:VAL:HA	7:A:501:HOH:O	2.05	0.56
1:B:182:LYS:HB3	1:B:258:LYS:HG2	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:252:ALA:O	1:B:256:LYS:HE3	2.07	0.53
1:A:9:ILE:O	1:A:38:LYS:HA	2.08	0.52
1:B:290:ASN:HD22	1:B:290:ASN:C	2.14	0.51
1:B:9:ILE:O	1:B:38:LYS:HA	2.10	0.51
1:A:61:ARG:NH2	1:B:49:GLU:O	2.42	0.49
1:B:160:ARG:HD2	1:B:356:GLU:OE1	2.14	0.48
1:A:210:LYS:O	1:A:212:LEU:N	2.46	0.48
1:B:118:HIS:HA	1:B:146:LEU:HD11	1.96	0.47
1:B:199:VAL:HG22	1:B:203:LYS:HE3	1.97	0.47
1:A:40:HIS:HA	1:A:59:GLN:HE22	1.81	0.46
1:B:256:LYS:NZ	7:B:507:HOH:O	2.44	0.46
1:B:3:MSE:O	1:B:32:HIS:HB3	2.15	0.46
1:B:171:SER:HB2	1:B:342:GLU:HG2	1.98	0.46
1:B:26:ASN:O	1:B:30:LYS:HE2	2.17	0.45
1:A:94:PHE:O	1:A:95:LEU:HB2	2.17	0.44
1:B:94:PHE:O	1:B:95:LEU:HB2	2.18	0.44
1:B:167:PRO:HD3	1:B:346:ILE:HG21	1.98	0.44
1:A:30:LYS:HE2	1:A:51:GLU:OE2	2.18	0.43
1:A:339:LYS:O	1:A:342:GLU:HG3	2.18	0.43
1:B:191:ARG:NH2	2:B:401:UD2:O1B	2.45	0.43
1:B:301:ASP:OD2	1:B:337:ARG:NH1	2.51	0.42
1:A:61:ARG:NH1	1:A:63:ASP:OD2	2.52	0.42
1:A:30:LYS:HE2	1:A:51:GLU:CD	2.40	0.42
1:A:148:SER:O	1:A:152:VAL:HG13	2.20	0.42
1:A:217:LYS:HE2	1:A:217:LYS:HB3	1.78	0.41
1:A:115:ILE:HG13	1:A:143:LEU:HD13	2.02	0.41
1:A:199:VAL:HG22	1:A:203:LYS:HE3	2.02	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	358/373 (96%)	346 (97%)	11 (3%)	1 (0%)	44	55
1	B	358/373 (96%)	349 (98%)	9 (2%)	0	100	100
All	All	716/746 (96%)	695 (97%)	20 (3%)	1 (0%)	55	67

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	211	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	324/332 (98%)	322 (99%)	2 (1%)	89	95
1	B	324/332 (98%)	322 (99%)	2 (1%)	89	95
All	All	648/664 (98%)	644 (99%)	4 (1%)	89	95

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	236	SER
1	A	290	ASN
1	B	3	MSE
1	B	290	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	59	GLN
1	B	67	HIS
1	B	249	ASN
1	B	290	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 ⓘ

Of 13 ligands modelled in this entry, 9 are monoatomic - leaving 4 for Mogul analysis.

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	UD2	A	401	5	34,41,41	2.79	10 (29%)	43,62,62	2.83	14 (32%)
4	GOL	A	403	-	5,5,5	0.35	0	5,5,5	0.26	0
2	UD2	B	401	5	34,41,41	2.74	10 (29%)	43,62,62	2.80	12 (27%)
4	GOL	B	403	-	5,5,5	0.36	0	5,5,5	0.27	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
2	UD2	A	401	5	1/1/12/13	0/22/63/63	0/3/3/3
4	GOL	A	403	-	-	0/4/4/4	0/0/0/0
2	UD2	B	401	5	-	0/22/63/63	0/3/3/3
4	GOL	B	403	-	-	0/4/4/4	0/0/0/0

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	UD2	O4B-C1B	-7.97	1.30	1.41
2	B	401	UD2	O4B-C1B	-7.71	1.30	1.41
2	A	401	UD2	C2B-C1B	-7.64	1.41	1.53
2	B	401	UD2	C2B-C1B	-7.43	1.41	1.53
2	A	401	UD2	O4B-C4B	-6.60	1.30	1.45
2	B	401	UD2	O4B-C4B	-6.46	1.30	1.45
2	B	401	UD2	C3B-C4B	-4.64	1.40	1.53
2	A	401	UD2	C3B-C4B	-4.56	1.41	1.53
2	B	401	UD2	C3B-C2B	-3.93	1.43	1.53
2	A	401	UD2	C3B-C2B	-3.87	1.43	1.53
2	B	401	UD2	O5'-C1'	2.16	1.47	1.41
2	A	401	UD2	O5'-C1'	2.16	1.47	1.41
2	B	401	UD2	PB-O1'	2.34	1.66	1.60
2	A	401	UD2	C2'-N2'	2.41	1.49	1.45
2	A	401	UD2	PB-O1'	2.46	1.66	1.60
2	B	401	UD2	C2'-N2'	2.48	1.49	1.45
2	B	401	UD2	O5'-C5'	2.65	1.50	1.44
2	A	401	UD2	C7'-N2'	2.74	1.44	1.34
2	B	401	UD2	C7'-N2'	2.75	1.44	1.34
2	A	401	UD2	O5'-C5'	2.75	1.51	1.44

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	UD2	O5'-C1'-O1'	-5.42	104.28	111.36
2	A	401	UD2	O5'-C1'-O1'	-4.83	105.05	111.36
2	A	401	UD2	C4B-O4B-C1B	-3.48	106.06	109.77
2	B	401	UD2	C4B-O4B-C1B	-3.09	106.48	109.77
2	B	401	UD2	C1'-O5'-C5'	-2.41	109.17	113.72
2	A	401	UD2	C1'-O5'-C5'	-2.39	109.21	113.72
2	A	401	UD2	O1B-PB-O2B	-2.07	101.57	112.28
2	B	401	UD2	C8'-C7'-N2'	2.15	119.98	116.11
2	A	401	UD2	C8'-C7'-N2'	2.16	120.01	116.11
2	A	401	UD2	C5B-C4B-C3B	2.16	123.53	115.29
2	B	401	UD2	O3A-PB-O1'	3.42	108.16	102.05
2	B	401	UD2	O4B-C4B-C5B	3.49	121.19	109.40
2	A	401	UD2	O4B-C4B-C5B	3.59	121.51	109.40
2	A	401	UD2	O3A-PB-O1'	3.92	109.05	102.05
2	B	401	UD2	O3B-C3B-C2B	4.04	124.78	111.83
2	A	401	UD2	O3B-C3B-C2B	4.05	124.81	111.83
2	A	401	UD2	O2'-C2B-C3B	4.48	126.18	111.83
2	B	401	UD2	O2'-C2B-C3B	4.62	126.61	111.83
2	A	401	UD2	O2'-C2B-C1B	4.74	126.43	111.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	UD2	O2'-C2B-C1B	4.95	127.11	111.61
2	A	401	UD2	O3B-C3B-C4B	5.06	125.86	111.09
2	B	401	UD2	O3B-C3B-C4B	5.11	126.01	111.09
2	B	401	UD2	O4B-C1B-N1	7.68	123.45	108.08
2	A	401	UD2	O4B-C1B-N1	8.07	124.24	108.08
2	B	401	UD2	C4-N3-C2	8.83	121.72	114.13
2	A	401	UD2	C4-N3-C2	9.01	121.87	114.13

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	401	UD2	C1B

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	B	401	UD2	1	0
4	B	403	GOL	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

### 6.1 Protein, DNA and RNA chains

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	356/373 (95%)	-0.07	7 (1%) 65 72	24, 37, 57, 75	7 (1%)
1	B	356/373 (95%)	-0.01	17 (4%) 31 38	22, 36, 64, 81	2 (0%)
All	All	712/746 (95%)	-0.04	24 (3%) 46 53	22, 37, 59, 81	9 (1%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	65	LEU	4.1
1	B	177	ASP	3.6
1	B	108	GLY	3.2
1	B	63	ASP	3.1
1	B	66	TYR	2.8
1	B	232	TYR	2.8
1	B	69	ILE	2.8
1	A	63	ASP	2.7
1	A	260	LEU	2.7
1	B	348	GLU	2.7
1	B	260	LEU	2.6
1	B	214	GLU	2.6
1	B	215	ASN	2.5
1	B	59	GLN	2.5
1	B	225	GLU	2.5
1	B	272	VAL	2.4
1	A	30	LYS	2.3
1	A	189	ILE	2.3
1	A	62	PHE	2.3
1	B	58	GLU	2.2
1	B	259	VAL	2.1
1	A	232	TYR	2.1
1	B	139	PHE	2.1
1	A	248	GLU	2.0

## 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 [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
4	GOL	B	403	6/6	0.78	0.25	4.81	45,47,54,57	0
5	NA	B	404	1/1	0.97	0.19	4.70	44,44,44,44	0
2	UD2	B	401	39/39	0.97	0.16	0.07	23,30,34,37	0
2	UD2	A	401	39/39	0.96	0.13	-0.19	24,33,39,40	0
3	K	B	402	1/1	0.98	0.14	-0.63	34,34,34,34	0
5	NA	A	405	1/1	0.93	0.13	-1.17	43,43,43,43	0
5	NA	A	404	1/1	0.99	0.10	-1.49	40,40,40,40	0
3	K	A	402	1/1	0.96	0.07	-2.48	44,44,44,44	0
4	GOL	A	403	6/6	0.87	0.40	-	62,66,67,68	0
6	CL	A	407	1/1	0.94	0.08	-	79,79,79,79	0
5	NA	B	405	1/1	0.98	0.06	-	38,38,38,38	0
6	CL	B	406	1/1	0.94	0.18	-	72,72,72,72	0
6	CL	A	406	1/1	0.94	0.15	-	67,67,67,67	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.