



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jun 18, 2024 – 02:32 AM EDT

PDB ID : 3LAQ  
Title : Structure-based engineering of species selectivity in the uPA-uPAR interaction  
Authors : Huang, M.  
Deposited on : 2010-01-06  
Resolution : 3.20 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)

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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>.

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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	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
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)	:	2.37.1

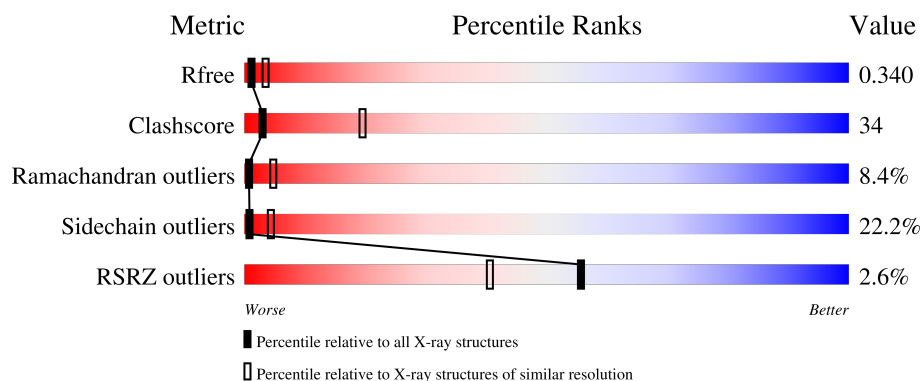
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.20 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	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

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  $\geq 3$ , 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	134	
1	B	134	
2	U	277	
2	V	277	

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
3	NAG	U	1160	-	-	-	X
3	NAG	V	1160	-	-	-	X
3	NAG	V	1259	-	-	X	-

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6035 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.

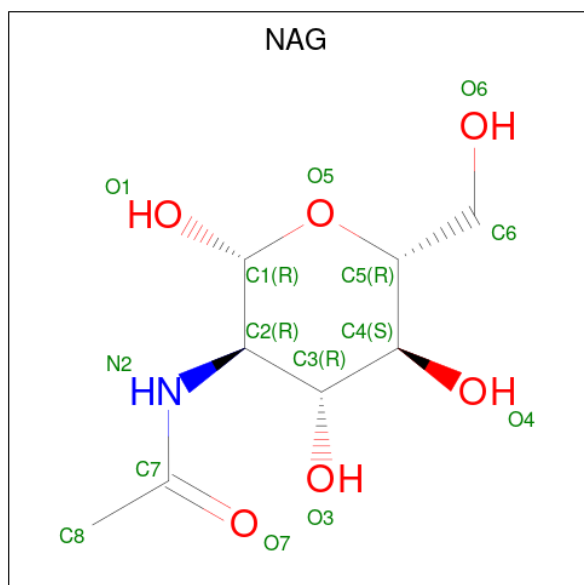
- Molecule 1 is a protein called Urokinase-type plasminogen activator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	124	Total	C	N	O	S	0	0	0
			981	601	190	177	13			
1	B	125	Total	C	N	O	S	0	0	0
			989	605	191	180	13			

- Molecule 2 is a protein called Urokinase plasminogen activator surface receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	U	258	Total	C	N	O	S	0	0	0
			1955	1173	355	394	33			
2	V	260	Total	C	N	O	S	0	0	0
			1970	1182	357	398	33			

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).

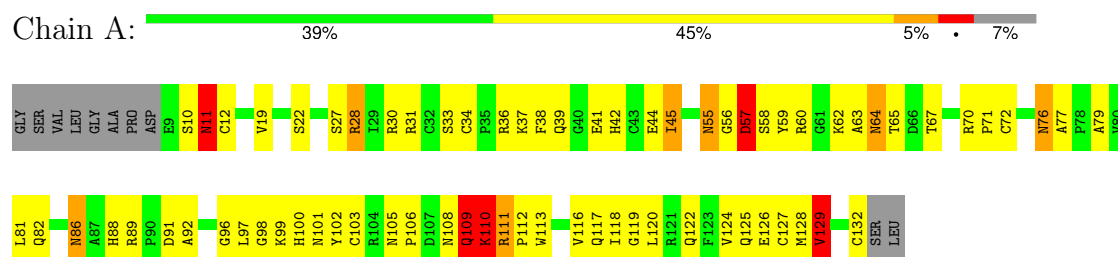


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	U	1	Total	C	N	O	0	0
			14	8	1	5		
3	U	1	Total	C	N	O	0	0
			14	8	1	5		
3	U	1	Total	C	N	O	0	0
			14	8	1	5		
3	U	1	Total	C	N	O	0	0
			14	8	1	5		
3	U	1	Total	C	N	O	0	0
			14	8	1	5		
3	V	1	Total	C	N	O	0	0
			14	8	1	5		
3	V	1	Total	C	N	O	0	0
			14	8	1	5		
3	V	1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

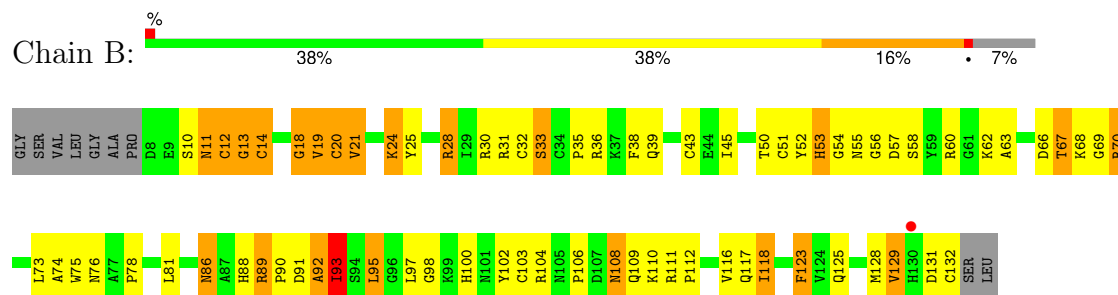
### 3 Residue-property plots [i](#)

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.

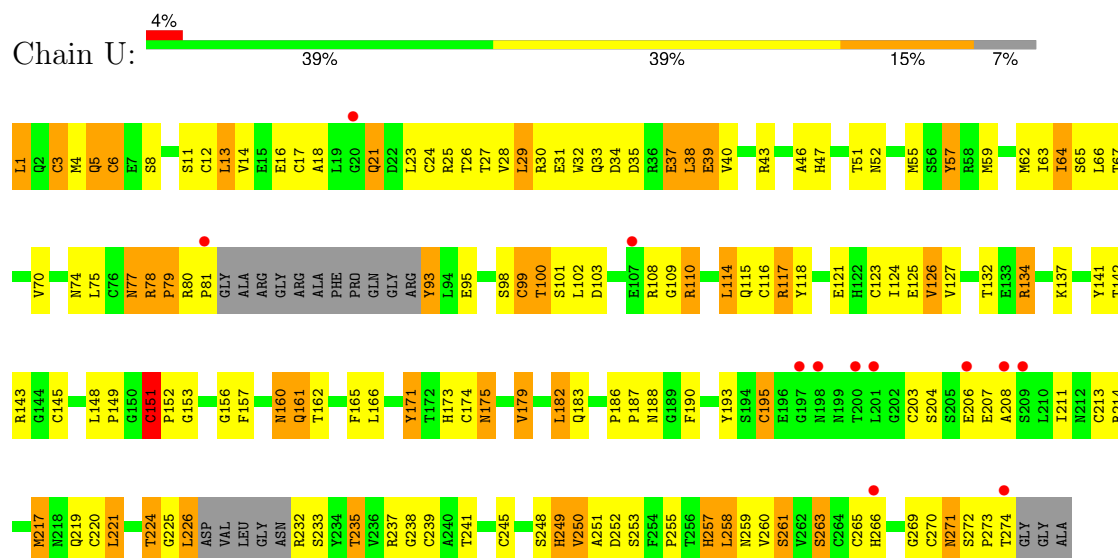
#### • Molecule 1: Urokinase-type plasminogen activator



#### • Molecule 1: Urokinase-type plasminogen activator

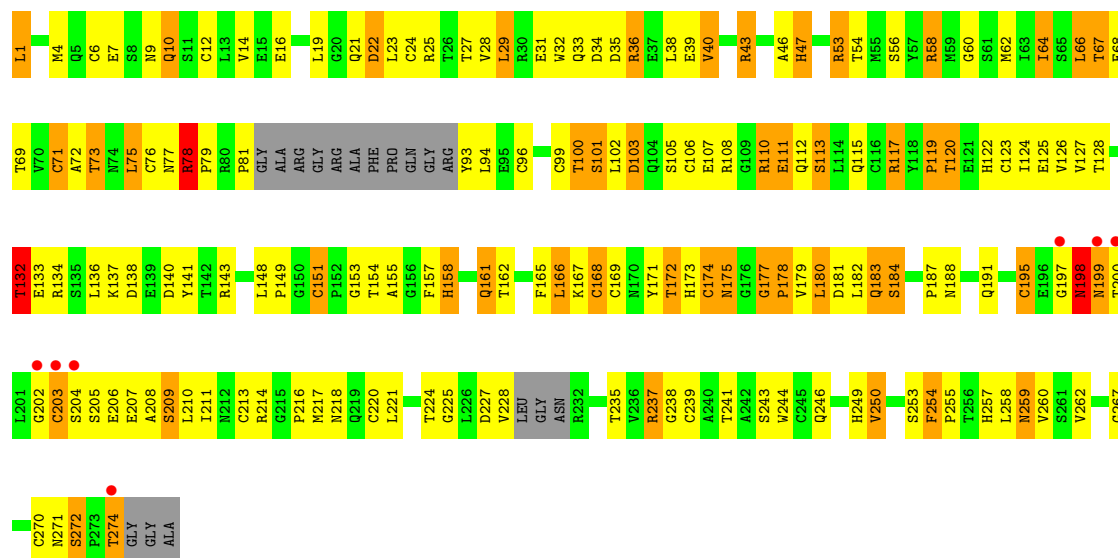


#### • Molecule 2: Urokinase plasminogen activator surface receptor



● Molecule 2: Urokinase plasminogen activator surface receptor

Chain V:  3% 34% 42% 17% 6%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	61.42Å 137.25Å 65.78Å 90.00° 106.17° 90.00°	Depositor
Resolution (Å)	44.74 – 3.20 44.73 – 3.20	Depositor EDS
% Data completeness (in resolution range)	93.3 (44.74-3.20) 93.3 (44.73-3.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.46 (at 3.19Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, $R_{free}$	0.231 , 0.341 0.234 , 0.340	Depositor DCC
$R_{free}$ test set	853 reflections (5.28%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	66.7	Xtriage
Anisotropy	0.329	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 52.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.88	EDS
Total number of atoms	6035	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.37% 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 [i](#)

### 5.1 Standard geometry [i](#)

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	0.72	1/1006 (0.1%)	0.82	1/1356 (0.1%)
1	B	0.65	0/1014	0.75	0/1367
2	U	0.68	2/1988 (0.1%)	0.85	2/2690 (0.1%)
2	V	0.69	0/2003	0.86	5/2711 (0.2%)
All	All	0.69	3/6011 (0.0%)	0.83	8/8124 (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	V	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	U	24	CYS	CB-SG	-5.88	1.72	1.81
2	U	245	CYS	CB-SG	-5.37	1.73	1.81
1	A	103	CYS	CB-SG	-5.10	1.73	1.81

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	103	CYS	CA-CB-SG	-6.94	101.50	114.00
2	U	24	CYS	CA-CB-SG	-6.50	102.29	114.00
2	V	158	HIS	O-C-N	6.44	133.00	122.70
2	U	114	LEU	CA-CB-CG	5.60	128.17	115.30
2	V	168	CYS	CA-CB-SG	5.53	123.96	114.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	V	78	ARG	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	981	0	919	62	0
1	B	989	0	923	80	0
2	U	1955	0	1806	116	0
2	V	1970	0	1821	153	0
3	U	70	0	64	1	0
3	V	70	0	64	9	0
All	All	6035	0	5597	400	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 34.

The worst 5 of 400 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)
2:V:259:ASN:OD1	3:V:1259:NAG:C7	1.89	1.20
2:V:259:ASN:OD1	3:V:1259:NAG:H2	1.42	1.19
2:V:259:ASN:OD1	3:V:1259:NAG:H83	1.45	1.16
2:V:259:ASN:OD1	3:V:1259:NAG:C8	1.93	1.14
2:V:259:ASN:OD1	3:V:1259:NAG:C2	1.96	1.12

There are no symmetry-related clashes.

## 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	122/134 (91%)	91 (75%)	23 (19%)	8 (7%)	1	9
1	B	123/134 (92%)	86 (70%)	23 (19%)	14 (11%)	0	2
2	U	252/277 (91%)	196 (78%)	41 (16%)	15 (6%)	1	12
2	V	254/277 (92%)	191 (75%)	37 (15%)	26 (10%)	0	3
All	All	751/822 (91%)	564 (75%)	124 (16%)	63 (8%)	1	5

5 of 63 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	109	GLN
1	A	129	VAL
2	U	78	ARG
2	U	134	ARG

### 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	106/113 (94%)	83 (78%)	23 (22%)	1	5
1	B	107/113 (95%)	86 (80%)	21 (20%)	1	7
2	U	228/238 (96%)	173 (76%)	55 (24%)	0	3
2	V	230/238 (97%)	180 (78%)	50 (22%)	1	5
All	All	671/702 (96%)	522 (78%)	149 (22%)	1	4

5 of 149 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	V	73	THR
2	V	241	THR
2	V	100	THR
2	V	166	LEU
2	U	124	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 29 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	55	ASN
2	V	249	HIS
1	B	122	GLN
2	V	199	ASN
1	B	100	HIS

### 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 [i](#)

10 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	U	1259	2	14,14,15	0.68	1 (7%)	17,19,21	2.56	5 (29%)
3	NAG	U	1053	-	14,14,15	1.06	1 (7%)	17,19,21	1.57	3 (17%)
3	NAG	U	1170	2	14,14,15	0.77	0	17,19,21	1.42	3 (17%)
3	NAG	U	1052	2	14,14,15	0.80	1 (7%)	17,19,21	1.72	4 (23%)
3	NAG	V	1160	2	14,14,15	0.96	1 (7%)	17,19,21	2.13	6 (35%)
3	NAG	V	1259	2	14,14,15	1.10	1 (7%)	17,19,21	1.85	5 (29%)
3	NAG	V	1170	2	14,14,15	0.46	0	17,19,21	1.28	2 (11%)
3	NAG	U	1160	-	14,14,15	0.79	1 (7%)	17,19,21	1.91	2 (11%)
3	NAG	V	1053	-	14,14,15	0.63	0	17,19,21	1.31	3 (17%)
3	NAG	V	1052	2	14,14,15	0.40	0	17,19,21	1.78	4 (23%)

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	U	1259	2	-	4/6/23/26	0/1/1/1
3	NAG	U	1053	-	-	0/6/23/26	0/1/1/1
3	NAG	U	1170	2	-	3/6/23/26	0/1/1/1
3	NAG	U	1052	2	-	2/6/23/26	0/1/1/1
3	NAG	V	1160	2	-	4/6/23/26	0/1/1/1
3	NAG	V	1259	2	-	4/6/23/26	0/1/1/1
3	NAG	V	1170	2	-	0/6/23/26	0/1/1/1
3	NAG	U	1160	-	-	2/6/23/26	0/1/1/1
3	NAG	V	1053	-	-	4/6/23/26	0/1/1/1
3	NAG	V	1052	2	-	2/6/23/26	0/1/1/1

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	V	1160	NAG	C1-C2	3.08	1.56	1.52
3	U	1053	NAG	C1-C2	2.98	1.56	1.52
3	U	1052	NAG	C1-C2	2.46	1.55	1.52
3	V	1259	NAG	C1-C2	2.45	1.55	1.52
3	U	1160	NAG	C1-C2	2.24	1.55	1.52

The worst 5 of 37 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	U	1259	NAG	C1-O5-C5	7.51	122.25	112.19
3	U	1160	NAG	C1-O5-C5	6.37	120.73	112.19
3	U	1052	NAG	C3-C4-C5	-4.81	101.52	110.23
3	U	1259	NAG	O5-C1-C2	4.59	118.40	111.29
3	V	1052	NAG	C4-C3-C2	-4.58	104.30	111.02

There are no chirality outliers.

5 of 25 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	U	1160	NAG	O5-C5-C6-O6
3	U	1052	NAG	O5-C5-C6-O6
3	V	1160	NAG	O5-C5-C6-O6
3	V	1053	NAG	O5-C5-C6-O6
3	U	1160	NAG	C4-C5-C6-O6

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	U	1053	NAG	1	0
3	U	1052	NAG	1	0
3	V	1160	NAG	1	0
3	V	1259	NAG	8	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	124/134 (92%)	-0.23	0 100 100	24, 34, 49, 72	0
1	B	125/134 (93%)	-0.16	1 (0%) 86 78	18, 31, 51, 77	0
2	U	258/277 (93%)	0.19	12 (4%) 31 19	11, 35, 48, 56	0
2	V	260/277 (93%)	0.08	7 (2%) 54 39	2, 34, 43, 50	0
All	All	767/822 (93%)	0.03	20 (2%) 56 40	2, 34, 47, 77	0

The worst 5 of 20 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	U	201	LEU	6.0
2	V	199	ASN	5.6
2	U	274	THR	4.8
2	V	200	THR	4.8
2	V	202	GLY	4.2

### 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](#)

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, 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	B-factors( $\text{\AA}^2$ )	Q<0.9
3	NAG	U	1052	14/15	0.62	0.28	60,66,67,72	0
3	NAG	U	1053	14/15	0.62	0.28	78,81,82,83	0
3	NAG	V	1259	14/15	0.68	0.28	71,78,85,85	0
3	NAG	V	1053	14/15	0.69	0.23	78,80,83,84	0
3	NAG	V	1160	14/15	0.69	0.44	62,67,73,74	0
3	NAG	U	1160	14/15	0.69	0.46	70,77,79,80	0
3	NAG	U	1259	14/15	0.85	0.30	69,77,78,78	0
3	NAG	U	1170	14/15	0.86	0.19	61,67,69,71	0
3	NAG	V	1052	14/15	0.90	0.14	59,66,68,73	0
3	NAG	V	1170	14/15	0.93	0.15	58,63,65,65	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.