



# wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 7, 2020 – 02:28 PM BST

PDB ID : 4R6U  
Title : IL-18 receptor complex  
Authors : Wei, H.; Wang, X.  
Deposited on : 2014-08-26  
Resolution : 2.80 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)

A user guide is available at

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.13.1  
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) : 2.13.1

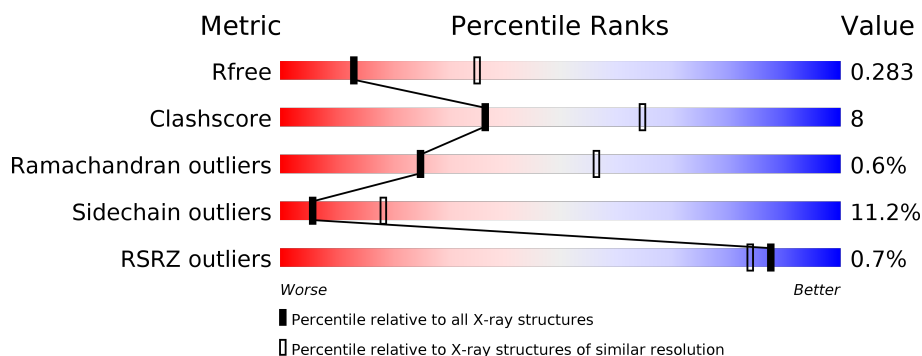
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 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	317	<div> <div></div> <div>58% 23% 5% 14%</div> </div>
1	C	317	<div> <div></div> <div>59% 23% • 13%</div> </div>
2	B	157	<div> <div></div> <div>83% 14% ••</div> </div>
2	D	157	<div> <div></div> <div>86% 11% ••</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7101 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 Interleukin-18 receptor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	275	Total	C	N	O	S	0	0	0
			2241	1425	388	414	14			
1	A	274	Total	C	N	O	S	0	0	0
			2234	1421	387	412	14			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	330	HIS	-	expression tag	UNP Q13478
C	331	HIS	-	expression tag	UNP Q13478
C	332	HIS	-	expression tag	UNP Q13478
C	333	HIS	-	expression tag	UNP Q13478
C	334	HIS	-	expression tag	UNP Q13478
C	335	HIS	-	expression tag	UNP Q13478
A	330	HIS	-	expression tag	UNP Q13478
A	331	HIS	-	expression tag	UNP Q13478
A	332	HIS	-	expression tag	UNP Q13478
A	333	HIS	-	expression tag	UNP Q13478
A	334	HIS	-	expression tag	UNP Q13478
A	335	HIS	-	expression tag	UNP Q13478

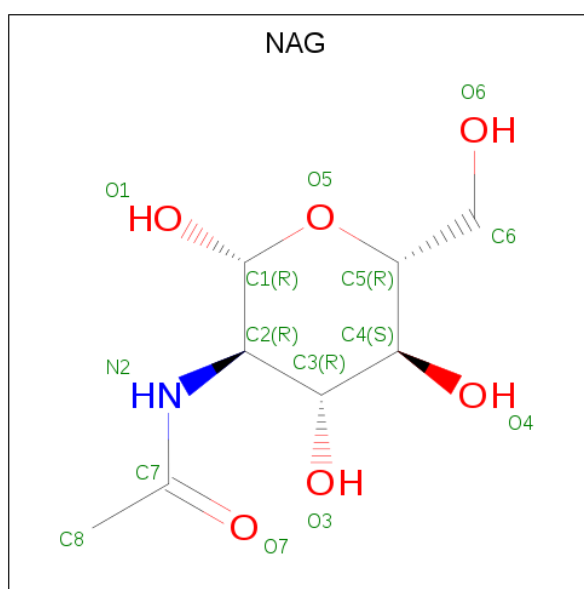
- Molecule 2 is a protein called Interleukin-18.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	155	Total	C	N	O	S	0	0	0
			1257	792	210	249	6			
2	D	155	Total	C	N	O	S	0	0	0
			1257	792	210	249	6			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	38	SER	CYS	engineered mutation	UNP Q14116
B	68	SER	CYS	engineered mutation	UNP Q14116
B	76	SER	CYS	engineered mutation	UNP Q14116
B	127	SER	CYS	engineered mutation	UNP Q14116
D	38	SER	CYS	engineered mutation	UNP Q14116
D	68	SER	CYS	engineered mutation	UNP Q14116
D	76	SER	CYS	engineered mutation	UNP Q14116
D	127	SER	CYS	engineered mutation	UNP Q14116

- 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	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		

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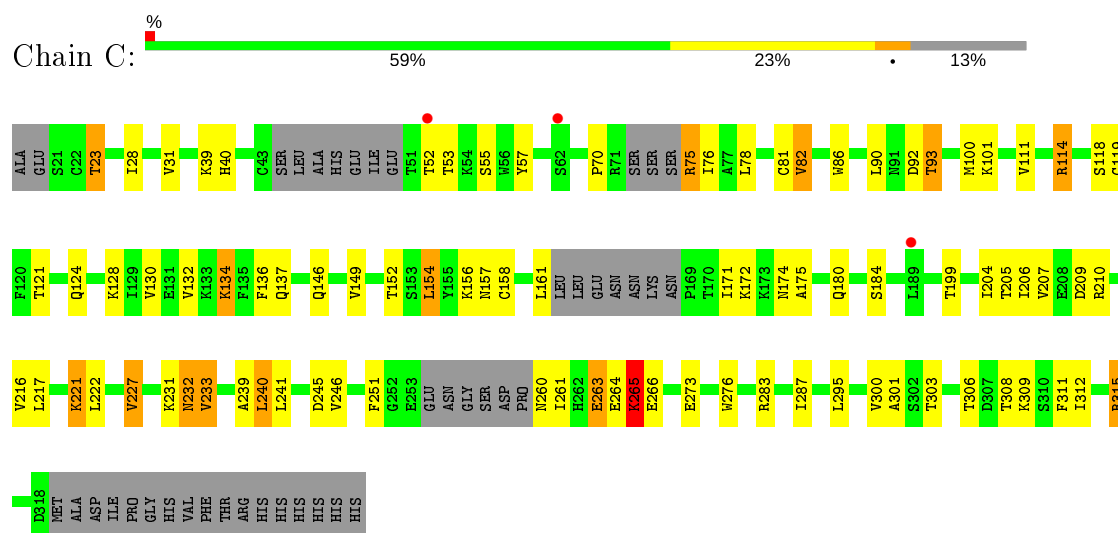
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	14	8	1	5	0	0

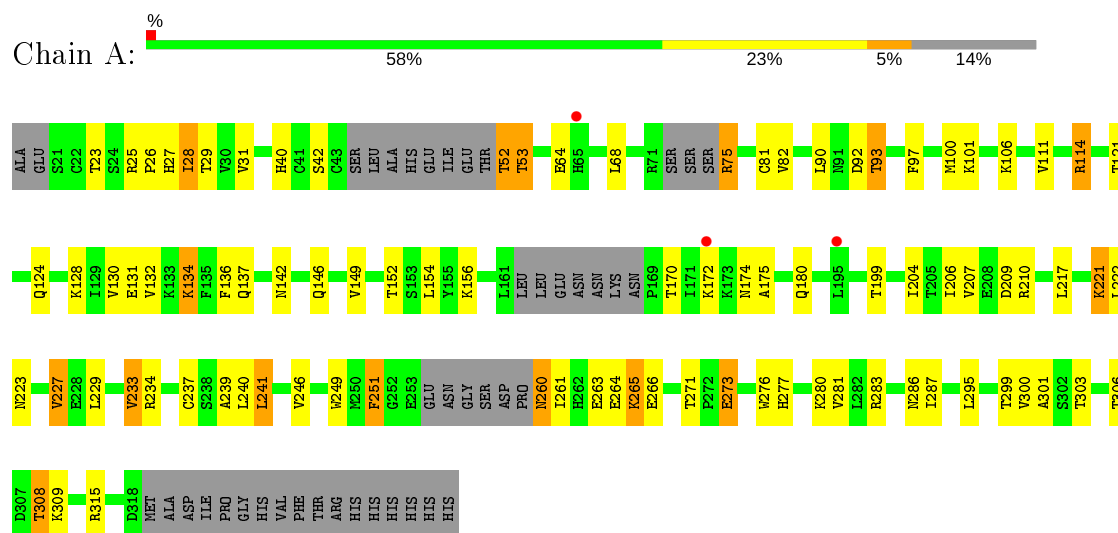
### 3 Residue-property plots [i](#)

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.

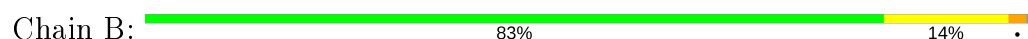
- Molecule 1: Interleukin-18 receptor 1



- Molecule 1: Interleukin-18 receptor 1

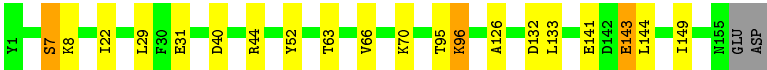
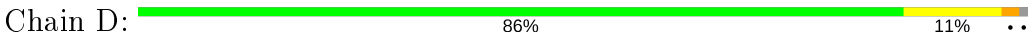


- Molecule 2: Interleukin-18





● Molecule 2: Interleukin-18



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	71.39Å 85.60Å 88.12Å 90.00° 97.60° 90.00°	Depositor
Resolution (Å)	38.90 – 2.80 38.90 – 2.60	Depositor EDS
% Data completeness (in resolution range)	96.9 (38.90-2.80) 96.1 (38.90-2.60)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.81 (at 2.61Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, $R_{free}$	0.233 , 0.282 0.236 , 0.283	Depositor DCC
$R_{free}$ test set	1591 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	57.3	Xtriage
Anisotropy	0.364	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 35.9	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	0.92	EDS
Total number of atoms	7101	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.02% 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: 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	0.63	0/2282	0.82	1/3078 (0.0%)
1	C	0.61	0/2289	0.82	0/3088
2	B	0.63	0/1278	0.86	4/1713 (0.2%)
2	D	0.61	1/1278 (0.1%)	0.81	2/1713 (0.1%)
All	All	0.62	1/7127 (0.0%)	0.83	7/9592 (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	C	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	141	GLU	CG-CD	5.12	1.59	1.51

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	44	ARG	NE-CZ-NH1	-9.63	115.49	120.30
2	B	44	ARG	NE-CZ-NH2	8.29	124.45	120.30
1	A	265	LYS	N-CA-C	7.78	132.00	111.00
2	D	143	GLU	OE1-CD-OE2	5.65	130.08	123.30
2	D	144	LEU	CB-CG-CD2	-5.53	101.60	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	264	GLU	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	2234	0	2197	47	0
1	C	2241	0	2203	49	0
2	B	1257	0	1249	10	0
2	D	1257	0	1249	7	0
3	A	56	0	52	3	0
3	C	56	0	52	1	0
All	All	7101	0	7002	112	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 8.

The worst 5 of 112 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:C:40:HIS:CE1	1:C:81:CYS:HB2	2.06	0.90
1:A:40:HIS:NE2	1:A:81:CYS:HB2	1.87	0.90
1:C:40:HIS:HE1	1:C:81:CYS:HB2	1.33	0.89
1:A:40:HIS:HE2	1:A:81:CYS:HB2	1.37	0.89
1:C:233:VAL:HG22	1:C:287:ILE:HD11	1.60	0.84

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	264/317 (83%)	240 (91%)	22 (8%)	2 (1%)	19	49
1	C	265/317 (84%)	240 (91%)	23 (9%)	2 (1%)	19	49
2	B	153/157 (98%)	146 (95%)	7 (5%)	0	100	100
2	D	153/157 (98%)	147 (96%)	5 (3%)	1 (1%)	22	53
All	All	835/948 (88%)	773 (93%)	57 (7%)	5 (1%)	25	56

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	143	GLU
1	A	265	LYS
1	A	64	GLU
1	C	265	LYS
1	C	315	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	255/293 (87%)	218 (86%)	37 (14%)	3	9
1	C	256/293 (87%)	218 (85%)	38 (15%)	3	9
2	B	146/148 (99%)	137 (94%)	9 (6%)	18	47
2	D	146/148 (99%)	140 (96%)	6 (4%)	30	64
All	All	803/882 (91%)	713 (89%)	90 (11%)	6	18

5 of 90 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	THR
1	A	134	LYS
2	B	154	GLN
1	A	53	THR

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Mol	Chain	Res	Type
1	A	93	THR

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

Mol	Chain	Res	Type
1	C	40	HIS
1	C	180	GLN
1	A	180	GLN
2	B	26	ASN

### 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](#)

8 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	A	402	1	14,14,15	1.59	1 (7%)	17,19,21	0.85	1 (5%)
3	NAG	C	801	1	14,14,15	1.77	3 (21%)	17,19,21	1.54	3 (17%)
3	NAG	A	401	1	14,14,15	0.93	1 (7%)	17,19,21	1.18	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	NAG	C	804	1	14,14,15	1.11	1 (7%)	17,19,21	1.17	1 (5%)
3	NAG	A	404	1	14,14,15	1.14	1 (7%)	17,19,21	1.12	1 (5%)
3	NAG	C	803	1	14,14,15	0.89	1 (7%)	17,19,21	0.80	0
3	NAG	A	403	1	14,14,15	0.35	0	17,19,21	1.10	0
3	NAG	C	802	1	14,14,15	0.95	2 (14%)	17,19,21	0.78	1 (5%)

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	A	402	1	-	2/6/23/26	0/1/1/1
3	NAG	C	801	1	-	2/6/23/26	0/1/1/1
3	NAG	A	401	1	-	2/6/23/26	0/1/1/1
3	NAG	C	804	1	-	2/6/23/26	0/1/1/1
3	NAG	A	404	1	-	0/6/23/26	0/1/1/1
3	NAG	C	803	1	-	0/6/23/26	0/1/1/1
3	NAG	A	403	1	-	1/6/23/26	0/1/1/1
3	NAG	C	802	1	-	0/6/23/26	0/1/1/1

The worst 5 of 10 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	NAG	O5-C1	5.59	1.52	1.43
3	C	801	NAG	O5-C1	-4.16	1.37	1.43
3	C	804	NAG	C1-C2	3.95	1.58	1.52
3	C	801	NAG	C2-N2	3.83	1.52	1.46
3	A	404	NAG	O5-C1	3.76	1.49	1.43

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	801	NAG	O5-C1-C2	-4.63	103.97	111.29
3	A	404	NAG	C1-O5-C5	3.55	117.01	112.19
3	C	804	NAG	O3-C3-C2	3.15	115.98	109.47
3	A	401	NAG	O5-C1-C2	-2.92	106.68	111.29
3	C	801	NAG	C1-O5-C5	-2.76	108.46	112.19

There are no chirality outliers.

5 of 9 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	804	NAG	O5-C5-C6-O6
3	C	804	NAG	C4-C5-C6-O6
3	C	801	NAG	O5-C5-C6-O6
3	A	402	NAG	C4-C5-C6-O6
3	A	401	NAG	O5-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	804	NAG	1	0
3	A	404	NAG	1	0
3	A	403	NAG	2	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	274/317 (86%)	0.01	3 (1%) 80 75	46, 61, 95, 118	0
1	C	275/317 (86%)	0.04	3 (1%) 80 75	44, 62, 91, 129	0
2	B	155/157 (98%)	-0.24	0 100 100	29, 45, 81, 92	0
2	D	155/157 (98%)	-0.24	0 100 100	32, 48, 85, 118	0
All	All	859/948 (90%)	-0.07	6 (0%) 87 84	29, 57, 91, 129	0

The worst 5 of 6 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	62	SER	3.9
1	A	172	LYS	3.2
1	A	65	HIS	3.0
1	C	52	THR	3.0
1	C	189	LEU	2.4

### 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	A	402	14/15	0.76	0.28	54,54,54,54	0
3	NAG	C	804	14/15	0.78	0.20	54,54,54,54	0
3	NAG	C	801	14/15	0.81	0.24	54,54,54,54	0
3	NAG	A	404	14/15	0.82	0.14	54,54,54,54	0
3	NAG	A	403	14/15	0.82	0.26	54,54,54,54	0
3	NAG	A	401	14/15	0.86	0.22	54,54,54,54	0
3	NAG	C	802	14/15	0.88	0.19	54,54,54,54	0
3	NAG	C	803	14/15	0.90	0.28	54,54,54,54	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.