



# wwPDB X-ray Structure Validation Summary Report

Feb 26, 2014 – 04:35 PM GMT

PDB ID : 1RKY  
Title : PPLO + Xe  
Authors : Guss, J.M.; Duff, A.P.  
Deposited on : 2003-11-24  
Resolution : 1.68 Å(reported)

This is a wwPDB 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 <http://wwpdb.org/ValidationPDFNotes.html>

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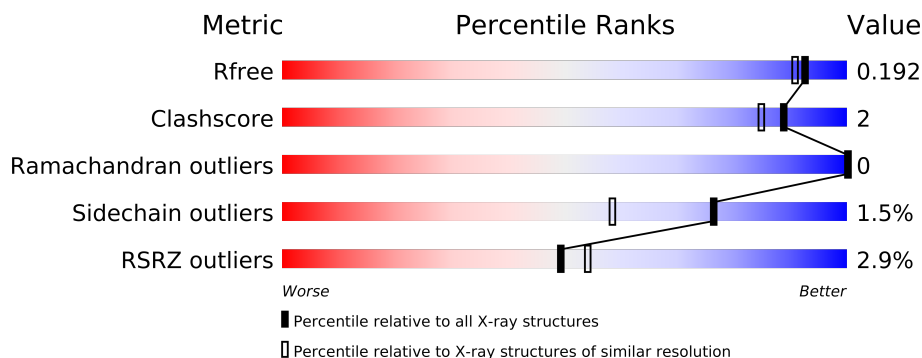
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance


The reported resolution of this entry is 1.68 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))
$R_{free}$	66092	3587 (1.70-1.66)
Clashscore	79885	4225 (1.70-1.66)
Ramachandran outliers	78287	4144 (1.70-1.66)
Sidechain outliers	78261	4143 (1.70-1.66)
RSRZ outliers	66119	3587 (1.70-1.66)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	747	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	NAG	A	1081[A]	-	X
2	NAG	A	1081[B]	-	X
2	NAG	A	1191	-	X
2	NAG	A	1309	-	X
2	NAG	A	1434	-	X
7	CL	A	812	-	X
7	CL	A	813	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
7	CL	A	814	-	X
7	CL	A	816	-	X
8	IMD	A	821	-	X

## 2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 6853 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

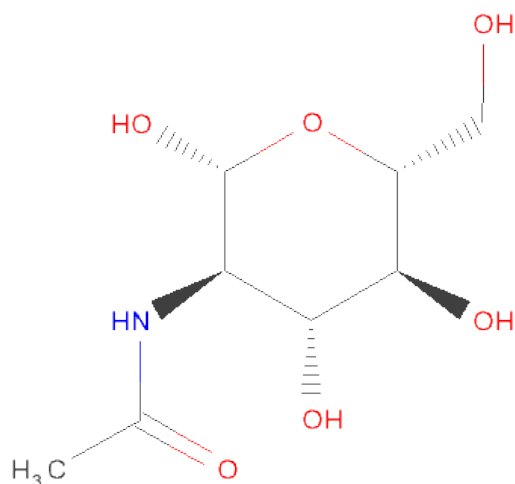
- Molecule 1 is a protein called lysyl oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	735	Total	C	N	O	S	2	49	0
			6087	3861	974	1236	16			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	314	VAL	ILE	CONFLICT	GB 13936870
A	338	ASP	GLU	CONFLICT	GB 13936870
A	417	SER	ALA	CONFLICT	GB 13936870
A	478	TPQ	TYR	MODIFIED RESIDUE	GB 13936870
A	549	GLN	LYS	CONFLICT	GB 13936870
A	577	PHE	LEU	CONFLICT	GB 13936870
A	579	ASN	LYS	CONFLICT	GB 13936870
A	650	THR	GLU	CONFLICT	GB 13936870
A	758	VAL	LEU	CONFLICT	GB 13936870
A	761	ALA	PRO	CONFLICT	GB 13936870

- Molecule 2 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	1
			28	16	2	10		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	3	Total	C	N	O	0	0
			39	22	2	15		

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	314	VAL	ILE	CONFLICT	GB 13936870
A	338	ASP	GLU	CONFLICT	GB 13936870
A	417	SER	ALA	CONFLICT	GB 13936870
A	478	TPQ	TYR	MODIFIED RESIDUE	GB 13936870
A	549	GLN	LYS	CONFLICT	GB 13936870
A	577	PHE	LEU	CONFLICT	GB 13936870
A	579	ASN	LYS	CONFLICT	GB 13936870
A	650	THR	GLU	CONFLICT	GB 13936870
A	758	VAL	LEU	CONFLICT	GB 13936870

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Chain	Residue	Modelled	Actual	Comment	Reference
A	761	ALA	PRO	CONFLICT	GB 13936870

- Molecule 4 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Cu 1 1	0	0

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	2	Total Ca 2 2	0	0

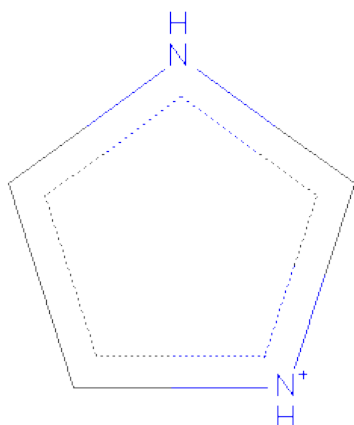
- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	3	Total Mg 3 3	0	0

- Molecule 7 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	6	Total Cl 6 6	0	0

- Molecule 8 is IMIDAZOLE (three-letter code: IMD) (formula: C<sub>3</sub>H<sub>5</sub>N<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	N	0	0
			5	3	2		

- Molecule 9 is XENON (three-letter code: XE) (formula: Xe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	8	Total	Xe	0	0
			8	8		

- Molecule 10 is water.

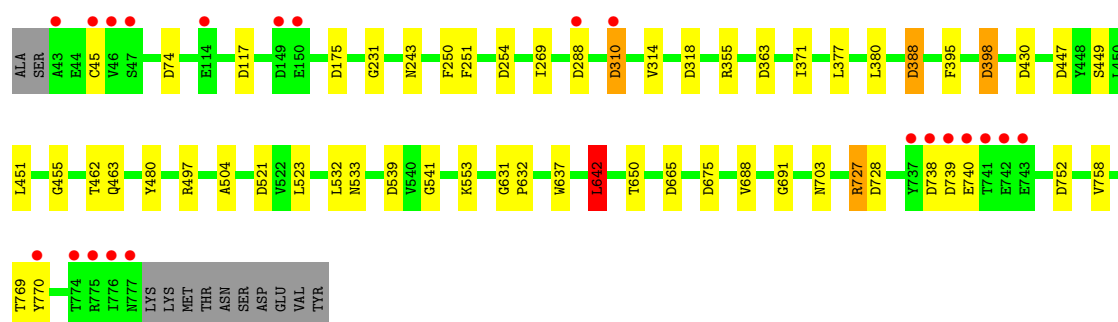
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	632	Total	O	0	0
			632	632		

### 3 Residue-property plots

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: lysyl oxidase

Chain A: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	140.14Å 66.70Å 108.62Å 90.00° 119.24° 90.00°	Depositor
Resolution (Å)	31.47 – 1.68 31.59 – 1.68	Depositor EDS
% Data completeness (in resolution range)	87.7 (31.47-1.68) 87.7 (31.59-1.68)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.63 (at 1.68Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, $R_{free}$	0.151 , 0.182 0.165 , 0.192	Depositor DCC
$R_{free}$ test set	1705 reflections (1.99%)	DCC
Wilson B-factor (Å <sup>2</sup> )	23.9	Xtriage
Anisotropy	0.076	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 49.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 182513 reflections	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	6853	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.72% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 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: MG, BMA, NAG, CL, CA, XE, TPQ, CU, IMD

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	0.59	0/6412	0.83	22/8732 (0.3%)

There are no bond length outliers.

The worst 5 of 22 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	642	LEU	CA-CB-CG	7.82	133.29	115.30
1	A	288	ASP	CB-CG-OD2	6.56	124.20	118.30
1	A	521[A]	ASP	CB-CG-OD2	6.43	124.09	118.30
1	A	521[B]	ASP	CB-CG-OD2	6.43	124.09	118.30
1	A	430	ASP	CB-CG-OD2	6.29	123.96	118.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6087	0	5599	28	1
2	A	70	0	65	0	0
3	A	39	0	34	0	0
4	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	2	0	0	0	0
6	A	3	0	0	0	0
7	A	6	0	0	0	0
8	A	5	0	5	0	0
9	A	8	0	0	9	0
10	A	632	0	0	4	0
All	All	6853	0	5703	28	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 2.

The worst 5 of 28 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:371[A]:ILE:HD12	9:A:901:XE:XE	1.76	1.63
1:A:532:LEU:CD1	9:A:907:XE:XE	2.31	1.56
1:A:371[A]:ILE:CD1	9:A:901:XE:XE	2.30	1.56
1:A:371[A]:ILE:HD11	9:A:901:XE:XE	2.07	1.23
1:A:532:LEU:HD11	9:A:907:XE:XE	2.12	1.19

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	Distance(Å)	Clash(Å)
1:A:455[B]:GLY:O	1:A:770[B]:TYR:OH[2.655]	2.06	0.14

## 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	776/747 (104%)	754 (97%)	22 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	686/658 (104%)	675 (98%)	11 (2%)	75 56

5 of 11 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	388	ASP
1	A	395	PHE
1	A	739	ASP
1	A	355	ARG
1	A	703	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	173	HIS
1	A	608	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is 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
1	TPQ	A	478	1,4	14,14,15	5.27	5 (35%)	17,19,21	1.77	2 (11%)

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
1	TPQ	A	478	1,4	-	0/4/22/24	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	478	TPQ	O-C	18.70	1.24	1.11
1	A	478	TPQ	O4-C4	-3.34	1.25	1.34
1	A	478	TPQ	C3-C4	3.18	1.40	1.35
1	A	478	TPQ	C6-C1	2.08	1.40	1.34
1	A	478	TPQ	CA-C	2.01	1.52	1.48

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	478	TPQ	CA-CB-C1	-5.03	104.21	114.41
1	A	478	TPQ	C-CA-N	-3.84	110.00	113.83

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.5 Carbohydrates ⓘ

3 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
3	NAG	A	1104	1,3	12,14,15	0.86	1 (8%)	15,19,21	0.83	0
3	NAG	A	1105	3	12,14,15	0.65	0	15,19,21	1.14	2 (13%)
3	BMA	A	1106	3	10,11,12	0.73	0	11,15,17	0.82	1 (9%)

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	1104	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	1105	3	-	0/6/23/26	0/1/1/1
3	BMA	A	1106	3	-	0/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1104	NAG	O5-C5	-2.68	1.40	1.45

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1105	NAG	O5-C5-C6	2.64	109.75	106.98
3	A	1105	NAG	C3-C2-N2	-2.29	108.28	111.76
3	A	1106	BMA	O5-C5-C6	2.18	109.27	106.98

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.6 Ligand geometry ⓘ

Of 26 ligands modelled in this entry, 20 are monoatomic - leaving 6 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	NAG	A	1081[A]	1	12,14,15	0.55	0	15,19,21	0.95	1 (6%)
2	NAG	A	1081[B]	1	12,14,15	0.60	0	15,19,21	0.86	1 (6%)
2	NAG	A	1191	1	12,14,15	0.50	0	15,19,21	1.27	1 (6%)
2	NAG	A	1309	1	12,14,15	0.66	0	15,19,21	0.91	0
2	NAG	A	1434	1	12,14,15	0.61	0	15,19,21	0.81	0
8	IMD	A	821	-	5,5,5	1.37	1 (20%)	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	NAG	A	1081[A]	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1081[B]	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1191	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1309	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1434	1	-	0/6/23/26	0/1/1/1
8	IMD	A	821	-	-	0/0/0/0	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	A	821	IMD	C2-N3	2.02	1.35	1.31

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1191	NAG	O5-C5-C6	3.18	110.31	106.98
2	A	1081[A]	NAG	C3-C2-N2	-2.11	108.55	111.76
2	A	1081[B]	NAG	C3-C2-N2	-2.11	108.55	111.76

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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 ⓘ

### 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	735/747 (98%)	-0.23	21 (2%) 49 54	21, 30, 52, 105	8 (1%)

The worst 5 of 21 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	776[A]	ILE	9.4
1	A	740	GLU	7.1
1	A	739	ASP	5.6
1	A	777[A]	ASN	5.0
1	A	47	SER	4.2

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
1	TPQ	A	478	14/15	0.14	3.07	25,40,73,108	9

### 6.3 Carbohydrates ⓘ

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	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	NAG	A	1105	14/15	0.13	2.22	39,52,78,88	0
3	NAG	A	1104	14/15	0.08	-0.19	31,48,62,63	0
3	BMA	A	1106	11/12	0.24	-	70,85,97,105	0

## 6.4 Ligands ⓘ

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	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
7	CL	A	816	1/1	0.30	56.03	49,49,49,49	1
2	NAG	A	1434	14/15	0.24	9.55	44,60,78,89	0
2	NAG	A	1191	14/15	0.28	8.02	39,53,71,75	0
8	IMD	A	821	5/5	0.15	6.58	38,44,48,60	5
2	NAG	A	1309	14/15	0.39	6.31	66,86,102,103	0
2	NAG	A	1081[B]	14/15	0.33	6.21	30,35,50,57	14
2	NAG	A	1081[A]	14/15	0.33	4.78	35,56,70,80	14
7	CL	A	814	1/1	0.07	4.08	48,48,48,48	1
7	CL	A	812	1/1	0.10	2.32	41,41,41,41	0
7	CL	A	813	1/1	0.14	2.19	43,43,43,43	0
7	CL	A	815	1/1	0.10	1.72	46,46,46,46	1
9	XE	A	908	1/1	0.08	0.99	39,39,39,39	1
9	XE	A	907	1/1	0.07	0.98	39,39,39,39	1
9	XE	A	902	1/1	0.08	0.46	41,41,41,41	1
9	XE	A	905	1/1	0.05	-0.46	38,38,38,38	1
7	CL	A	811	1/1	0.06	-0.56	40,40,40,40	0
6	MG	A	805	1/1	0.09	-0.75	37,37,37,37	1
5	CA	A	802	1/1	0.05	-1.05	23,23,23,23	0
9	XE	A	903	1/1	0.04	-1.10	34,34,34,34	1
6	MG	A	806	1/1	0.06	-1.69	59,59,59,59	1
9	XE	A	906	1/1	0.05	-1.72	43,43,43,43	1
9	XE	A	904	1/1	0.04	-2.46	38,38,38,38	1
9	XE	A	901	1/1	0.04	-2.50	27,27,27,27	1
4	CU	A	801	1/1	0.02	-2.51	32,32,32,32	0
5	CA	A	803	1/1	0.02	-3.51	25,25,25,25	0
6	MG	A	804	1/1	0.02	-10.31	31,31,31,31	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.