



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:10 PM GMT

PDB ID : 3T3N  
Title : Molecular basis for the recognition and cleavage of RNA (UUCCGU) by the bifunctional 5'-3' exo/endoribonuclease RNase J  
Authors : Dorleans, A.; Li de la Sierra-Gallay, I.; Piton, J.; Zig, L.; Gilet, L.; Putzer, H.; Condon, C.  
Deposited on : 2011-07-25  
Resolution : 3.09 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

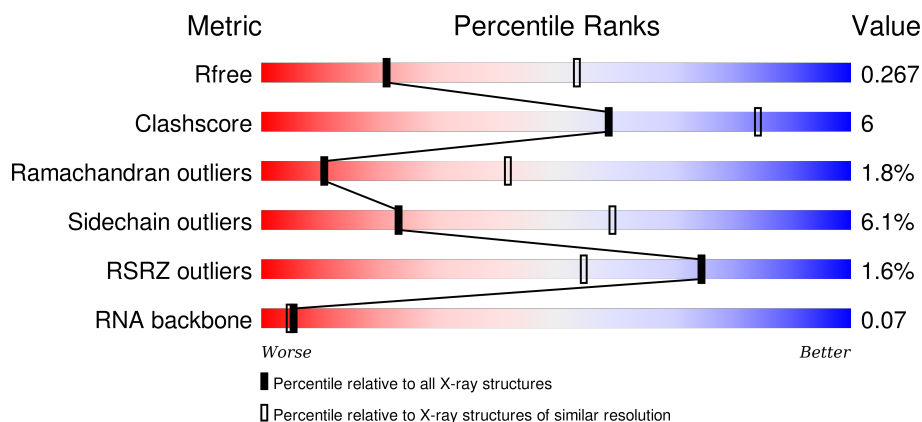
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.09 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))
$R_{free}$	91344	1114 (3.14-3.06)
Clashscore	102246	1222 (3.14-3.06)
Ramachandran outliers	100387	1174 (3.14-3.06)
Sidechain outliers	100360	1174 (3.14-3.06)
RSRZ outliers	91569	1119 (3.14-3.06)
RNA backbone	2183	1010 (3.52-2.68)

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	562	 2% 78% 18% ..
2	B	6	 33% 67%

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4455 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 Metal dependent hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	550	4311	2768	757	772	14	0	0	0

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	EXPRESSION TAG	UNP Q72JJ7
A	-6	SER	-	EXPRESSION TAG	UNP Q72JJ7
A	-5	HIS	-	EXPRESSION TAG	UNP Q72JJ7
A	-4	HIS	-	EXPRESSION TAG	UNP Q72JJ7
A	-3	HIS	-	EXPRESSION TAG	UNP Q72JJ7
A	-2	HIS	-	EXPRESSION TAG	UNP Q72JJ7
A	-1	HIS	-	EXPRESSION TAG	UNP Q72JJ7
A	0	HIS	-	EXPRESSION TAG	UNP Q72JJ7
A	1	SER	-	EXPRESSION TAG	UNP Q72JJ7
A	77	ALA	HIS	ENGINEERED MUTATION	UNP Q72JJ7

- Molecule 2 is a RNA chain called O2'methyl-RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	B	6	130	61	17	46	6	0	0	0

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		

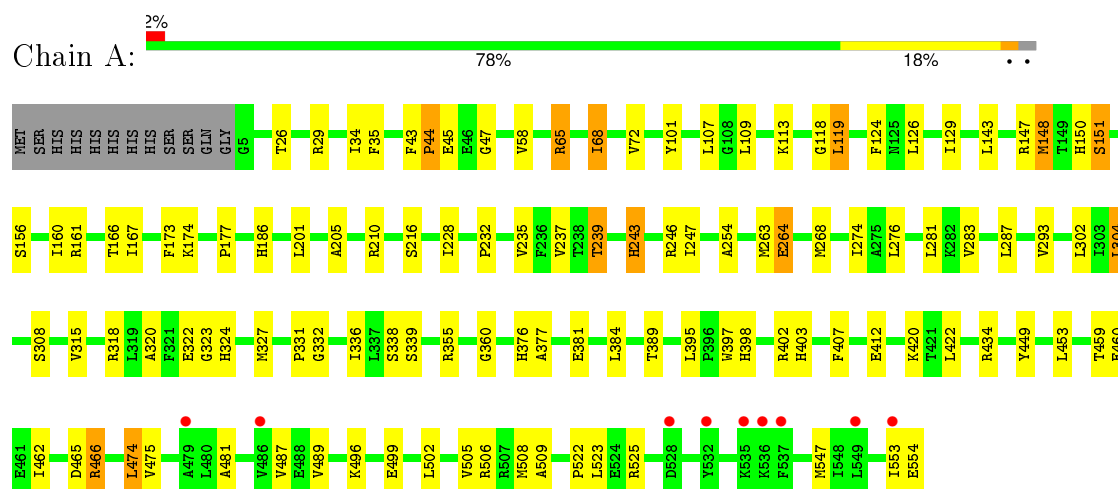
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	13	Total	O	0	0
			13	13		

### 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 ( $\text{RSRZ} > 2$ ). 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: Metal dependent hydrolase



- Molecule 2: O2'methyl-RNA



## 4 Data and refinement statistics

Property	Value	Source
Space group	F 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	101.95Å 116.78Å 231.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	54.41 – 3.09 54.41 – 3.09	Depositor EDS
% Data completeness (in resolution range)	(Not available) (54.41-3.09) 95.7 (54.41-3.09)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.06 (at 3.07Å)	Xtriage
Refinement program	BUSTER 2.8.0	Depositor
R, $R_{free}$	0.202 , 0.260 0.217 , 0.267	Depositor DCC
$R_{free}$ test set	1004 reflections (8.89%)	DCC
Wilson B-factor (Å <sup>2</sup> )	71.1	Xtriage
Anisotropy	0.264	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 53.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 12301 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	4455	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.38% 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 [i](#)

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

Bond lengths and bond angles in the following residue types are not validated in this section: OMC, ZN, OMG, OMU

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	0.41	0/4408	0.66	0/5968

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-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 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	4311	0	4412	49	0
2	B	130	0	76	4	0
3	A	1	0	0	0	0
4	A	13	0	0	0	0
All	All	4455	0	4488	52	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 6.

All (52) 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:147:ARG:HB3	1:A:186:HIS:HD2	1.56	0.71
1:A:268:MET:HE3	1:A:304:LEU:HB3	1.73	0.69
1:A:243:HIS:O	1:A:247:ILE:HG12	1.99	0.62
2:B:903:OMC:H5'	2:B:903:OMC:H6	1.69	0.58
1:A:173:PHE:O	1:A:376:HIS:HA	2.05	0.56
1:A:505:VAL:HA	1:A:508:MET:HB2	1.89	0.55
1:A:459:THR:H	1:A:462:ILE:HD12	1.72	0.54
1:A:148:MET:HG3	1:A:173:PHE:HB2	1.90	0.54
1:A:320:ALA:HB1	1:A:355:ARG:HB3	1.89	0.53
1:A:72:VAL:HG21	1:A:143:LEU:HD22	1.91	0.52
1:A:465:ASP:HB2	1:A:547:MET:HE1	1.92	0.51
1:A:324:HIS:HB2	1:A:327:MET:O	2.11	0.51
1:A:237:VAL:HG22	1:A:336:ILE:HB	1.94	0.49
2:B:905:OMG:HM21	2:B:906:OMU:C6	2.42	0.49
1:A:318:ARG:O	1:A:323:GLY:HA2	2.13	0.49
1:A:109:LEU:HA	1:A:274:ILE:HD13	1.95	0.49
1:A:239:THR:HG21	1:A:246:ARG:HG2	1.96	0.48
1:A:107:LEU:HD22	1:A:126:LEU:HB3	1.96	0.48
1:A:150:HIS:CG	1:A:151:SER:H	2.32	0.47
1:A:489:VAL:H	1:A:506:ARG:HD3	1.80	0.47
1:A:65:ARG:HA	1:A:68:ILE:HD12	1.97	0.47
1:A:420:LYS:HE2	1:A:422:LEU:HD21	1.97	0.47
1:A:205:ALA:O	1:A:403:HIS:HB3	2.15	0.46
1:A:119:LEU:HD11	1:A:124:PHE:HE2	1.80	0.46
1:A:283:VAL:HG21	1:A:287:LEU:HG	1.97	0.46
1:A:101:TYR:HB3	1:A:129:ILE:HG12	1.98	0.46
1:A:239:THR:HG23	1:A:338:SER:O	2.16	0.45
1:A:481:ALA:HB3	1:A:554:GLU:HA	1.99	0.45
2:B:905:OMG:HM23	2:B:905:OMG:N3	2.33	0.44
1:A:160:ILE:HB	1:A:167:ILE:HB	2.00	0.44
1:A:177:PRO:HD2	1:A:216:SER:HB2	1.99	0.44
1:A:29:ARG:HG3	1:A:34:ILE:HG13	1.98	0.44
1:A:377:ALA:HB1	1:A:381:GLU:HB2	2.00	0.44
1:A:264:GLU:HB3	1:A:315:VAL:HG21	1.99	0.43
1:A:487:VAL:HG11	1:A:509:ALA:HB3	2.00	0.43
1:A:174:LYS:HA	1:A:377:ALA:HB2	2.00	0.43
1:A:201:LEU:HD13	1:A:389:THR:HG21	2.00	0.43
1:A:324:HIS:HB2	1:A:327:MET:H	1.84	0.43
1:A:228:ILE:HG12	1:A:235:VAL:HG11	2.00	0.42
1:A:43:PHE:CE1	2:B:902:OMU:HM22	2.55	0.42
1:A:254:ALA:HB2	1:A:302:LEU:HD13	2.01	0.42
1:A:489:VAL:HG21	1:A:505:VAL:HG23	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:276:LEU:HD12	1:A:281:LEU:HB3	2.01	0.41
1:A:522:PRO:HD2	1:A:525:ARG:HB2	2.02	0.41
1:A:474:LEU:HG	1:A:475:VAL:N	2.35	0.41
1:A:499:GLU:HA	1:A:502:LEU:HG	2.02	0.41
1:A:308:SER:HB2	1:A:339:SER:OG	2.20	0.41
1:A:449:TYR:HE1	1:A:460:GLU:HG3	1.86	0.41
1:A:263:MET:HG3	1:A:287:LEU:HD22	2.02	0.40
1:A:462:ILE:O	1:A:466:ARG:HB2	2.21	0.40
1:A:397:TRP:CG	1:A:398:HIS:N	2.89	0.40
1:A:331:PRO:HB3	1:A:360:GLY:HA3	2.04	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	548/562 (98%)	507 (92%)	31 (6%)	10 (2%)	11	42

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	44	PRO
1	A	232	PRO
1	A	45	GLU
1	A	47	GLY
1	A	65	ARG
1	A	322	GLU
1	A	118	GLY
1	A	58	VAL
1	A	68	ILE
1	A	332	GLY

### 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	459/470 (98%)	431 (94%)	28 (6%)	23 59

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	THR
1	A	35	PHE
1	A	44	PRO
1	A	113	LYS
1	A	119	LEU
1	A	148	MET
1	A	151	SER
1	A	156	SER
1	A	161	ARG
1	A	166	THR
1	A	210	ARG
1	A	239	THR
1	A	243	HIS
1	A	264	GLU
1	A	293	VAL
1	A	304	LEU
1	A	384	LEU
1	A	395	LEU
1	A	402	ARG
1	A	407	PHE
1	A	412	GLU
1	A	434	ARG
1	A	453	LEU
1	A	466	ARG
1	A	474	LEU
1	A	496	LYS
1	A	523	LEU
1	A	553	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	HIS
1	A	136	GLN
1	A	186	HIS
1	A	345	ASN
1	A	379	GLN

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	5/6 (83%)	4 (80%)	0

All (4) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	902	OMU
2	B	903	OMC
2	B	905	OMG
2	B	906	OMU

There are no RNA pucker outliers to report.

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

6 non-standard protein/DNA/RNA residues 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	OMU	B	901	2	17,23,23	1.22	3 (17%)	25,34,34	1.70	3 (12%)
2	OMU	B	902	2	12,22,23	1.25	2 (16%)	19,31,34	1.97	3 (15%)
2	OMC	B	903	2	13,22,23	1.21	1 (7%)	20,31,34	1.60	4 (20%)
2	OMC	B	904	2	13,22,23	1.17	1 (7%)	20,31,34	1.75	4 (20%)
2	OMG	B	905	2	17,26,27	1.42	3 (17%)	21,38,41	1.89	5 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	OMU	B	906	2	12,22,23	1.55	2 (16%)	19,31,34	1.78	2 (10%)

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	OMU	B	901	2	-	0/8/28/28	0/2/2/2
2	OMU	B	902	2	-	0/5/27/28	0/2/2/2
2	OMC	B	903	2	-	0/5/27/28	0/2/2/2
2	OMC	B	904	2	-	0/5/27/28	0/2/2/2
2	OMG	B	905	2	-	0/5/27/28	0/3/3/3
2	OMU	B	906	2	-	0/5/27/28	0/2/2/2

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	905	OMG	C8-N7	-2.36	1.30	1.34
2	B	902	OMU	C4-N3	2.04	1.36	1.33
2	B	901	OMU	P-OP3	2.12	1.62	1.54
2	B	905	OMG	C2-N1	2.25	1.39	1.35
2	B	901	OMU	C4-N3	2.28	1.37	1.33
2	B	901	OMU	C6-N1	3.13	1.40	1.35
2	B	906	OMU	C4-N3	3.27	1.39	1.33
2	B	905	OMG	C6-N1	3.34	1.39	1.33
2	B	902	OMU	C6-N1	3.55	1.40	1.35
2	B	904	OMC	C6-N1	3.65	1.40	1.35
2	B	903	OMC	C6-N1	3.81	1.41	1.35
2	B	906	OMU	C6-N1	4.03	1.41	1.35

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	905	OMG	N3-C2-N1	-4.41	120.72	127.44
2	B	905	OMG	C5-C6-N1	-3.47	118.84	123.59
2	B	904	OMC	CM2-O2'-C2'	-2.65	107.10	114.59
2	B	903	OMC	CM2-O2'-C2'	-2.52	107.47	114.59
2	B	906	OMU	CM2-O2'-C2'	-2.51	107.50	114.59
2	B	902	OMU	CM2-O2'-C2'	-2.49	107.56	114.59
2	B	905	OMG	CM2-O2'-C2'	-2.42	107.74	114.59
2	B	904	OMC	N4-C4-N3	2.02	120.18	116.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	901	OMU	O4'-C1'-N1	2.19	112.70	108.08
2	B	901	OMU	OP2-P-OP1	2.22	117.73	110.58
2	B	903	OMC	N4-C4-N3	2.39	120.86	116.50
2	B	905	OMG	C1'-N9-C4	2.54	130.77	126.94
2	B	902	OMU	O4'-C1'-N1	2.62	113.60	108.08
2	B	903	OMC	O4'-C1'-N1	3.25	114.94	108.08
2	B	905	OMG	C6-N1-C2	3.31	120.53	115.94
2	B	903	OMC	C2-N3-C4	4.17	121.49	115.61
2	B	904	OMC	C2-N3-C4	4.29	121.66	115.61
2	B	904	OMC	O4'-C1'-N1	4.38	117.32	108.08
2	B	901	OMU	C4-N3-C2	6.79	120.87	114.14
2	B	906	OMU	C4-N3-C2	6.92	121.00	114.14
2	B	902	OMU	C4-N3-C2	7.32	121.39	114.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	902	OMU	1	0
2	B	903	OMC	1	0
2	B	905	OMG	2	0
2	B	906	OMU	1	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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	550/562 (97%)	-0.04	9 (1%) 74 55	37, 71, 112, 134	0
2	B	0/6	-	-	-	-
All	All	550/568 (96%)	-0.04	9 (1%) 74 55	37, 71, 112, 134	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	537	PHE	3.4
1	A	479	ALA	2.9
1	A	536	LYS	2.7
1	A	549	LEU	2.2
1	A	528	ASP	2.1
1	A	486	VAL	2.1
1	A	535	LYS	2.0
1	A	553	ILE	2.0
1	A	532	TYR	2.0

### 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	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	OMC	B	904	21/22	0.90	0.20	-	118,125,127,127	0
2	OMU	B	901	22/22	0.85	0.28	-	116,119,119,120	0
2	OMC	B	903	21/22	0.94	0.20	-	115,115,116,118	0
2	OMU	B	902	21/22	0.91	0.22	-	104,114,118,118	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	OMU	B	906	21/22	0.72	0.36	-	142,145,150,150	0
2	OMG	B	905	24/25	0.78	0.28	-	129,138,144,144	0

### 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( $\text{\AA}^2$ )	Q<0.9
3	ZN	A	600	1/1	0.86	0.14	-3.02	169,169,169,169	0

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

There are no such residues in this entry.