



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 15, 2017 – 03:08 am GMT

PDB ID : 3PU4  
Title : Crystal Structure of a vesicular stomatitis virus nucleocapsid-polyU complex  
Authors : Luo, M.; Green, T.J.; Rowse, M.  
Deposited on : 2010-12-03  
Resolution : 3.00 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

<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
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
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)	:	recalc28949

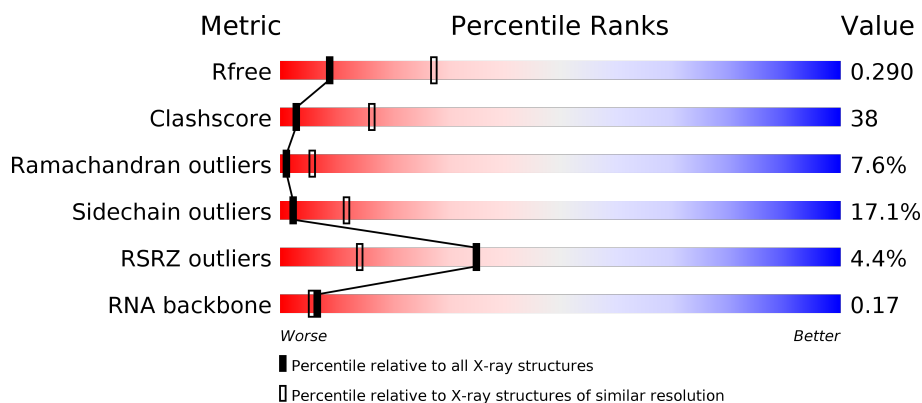
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.00 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	1692 (3.00-3.00)
Clashscore	112137	2037 (3.00-3.00)
Ramachandran outliers	110173	1973 (3.00-3.00)
Sidechain outliers	110143	1976 (3.00-3.00)
RSRZ outliers	101464	1716 (3.00-3.00)
RNA backbone	2435	1007 (3.34-2.66)

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. 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	421	<div> <div>4%</div> <div>48% 38% 12%</div> </div>
1	B	421	<div> <div>2%</div> <div>47% 38% 11%</div> </div>
1	C	421	<div> <div>3%</div> <div>48% 37% 11%</div> </div>
1	D	421	<div> <div>5%</div> <div>48% 37% 12%</div> </div>

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Mol	Chain	Length	Quality of chain
1	E	421	<div><div></div><div>4%</div><div>46%</div><div>39%</div><div>12%</div><div></div></div>
2	R	45	<div><div></div><div>42%</div><div>7%</div><div>24%</div><div>33%</div><div>36%</div><div></div></div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 17437 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.

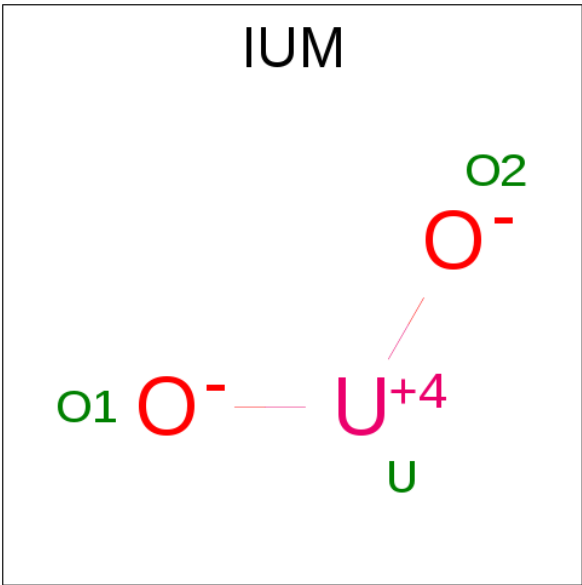
- Molecule 1 is a protein called Nucleoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	421	Total	C	N	O	S	0	0	0
			3327	2118	558	633	18			
1	B	415	Total	C	N	O	S	0	0	0
			3290	2097	552	623	18			
1	C	413	Total	C	N	O	S	0	0	0
			3275	2089	550	618	18			
1	D	416	Total	C	N	O	S	0	0	0
			3298	2103	553	624	18			
1	E	421	Total	C	N	O	S	0	0	0
			3327	2118	558	633	18			

- Molecule 2 is a RNA chain called RNA (45-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	R	45	Total	C	N	O	P	0	0	0
			900	405	90	360	45			

- Molecule 3 is URANYL (VI) ION (three-letter code: IUM) (formula: O<sub>2</sub>U).



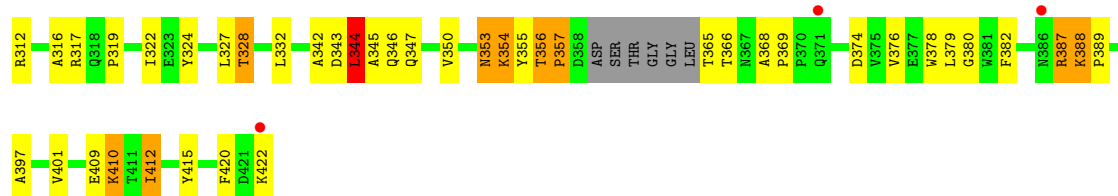
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total U 1 1	0	0
3	A	1	Total U 1 1	0	0
3	A	1	Total U 1 1	0	0
3	A	1	Total U 1 1	0	0
3	B	1	Total U 1 1	0	0
3	B	1	Total U 1 1	0	0
3	B	1	Total U 1 1	0	0
3	C	1	Total U 1 1	0	0
3	C	1	Total U 1 1	0	0
3	C	1	Total U 1 1	0	0
3	D	1	Total U 1 1	0	0
3	D	1	Total U 1 1	0	0
3	E	1	Total U 1 1	0	0
3	E	1	Total U 1 1	0	0

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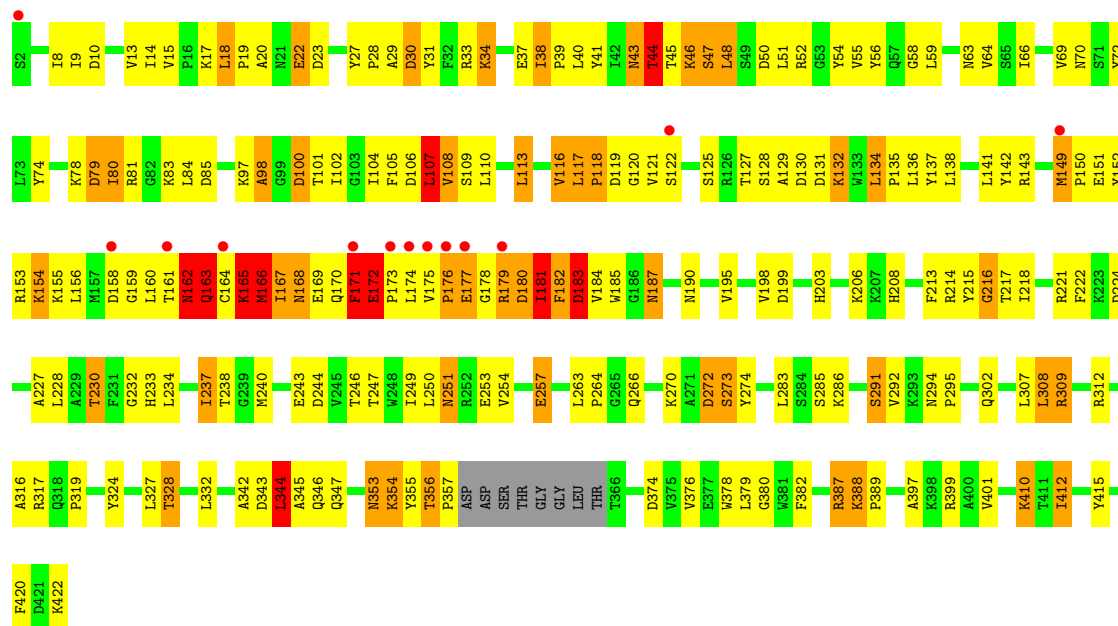
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	1	Total 1	U 1	0	0
3	R	1	Total 1	U 1	0	0
3	R	1	Total 1	U 1	0	0
3	R	1	Total 1	U 1	0	0
3	R	1	Total 1	U 1	0	0
3	R	1	Total 1	U 1	0	0

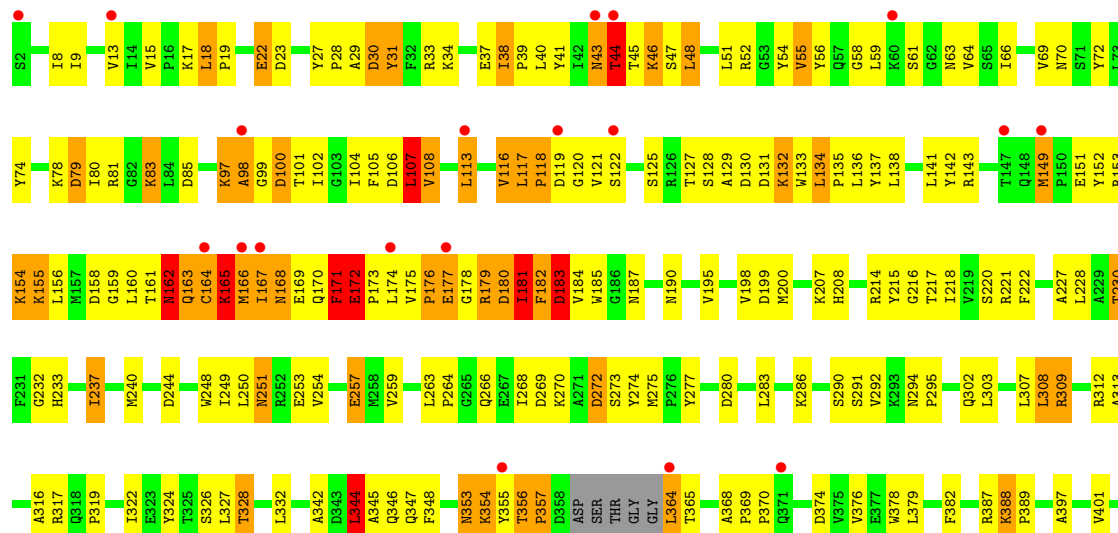




### • Molecule 1: Nucleoprotein



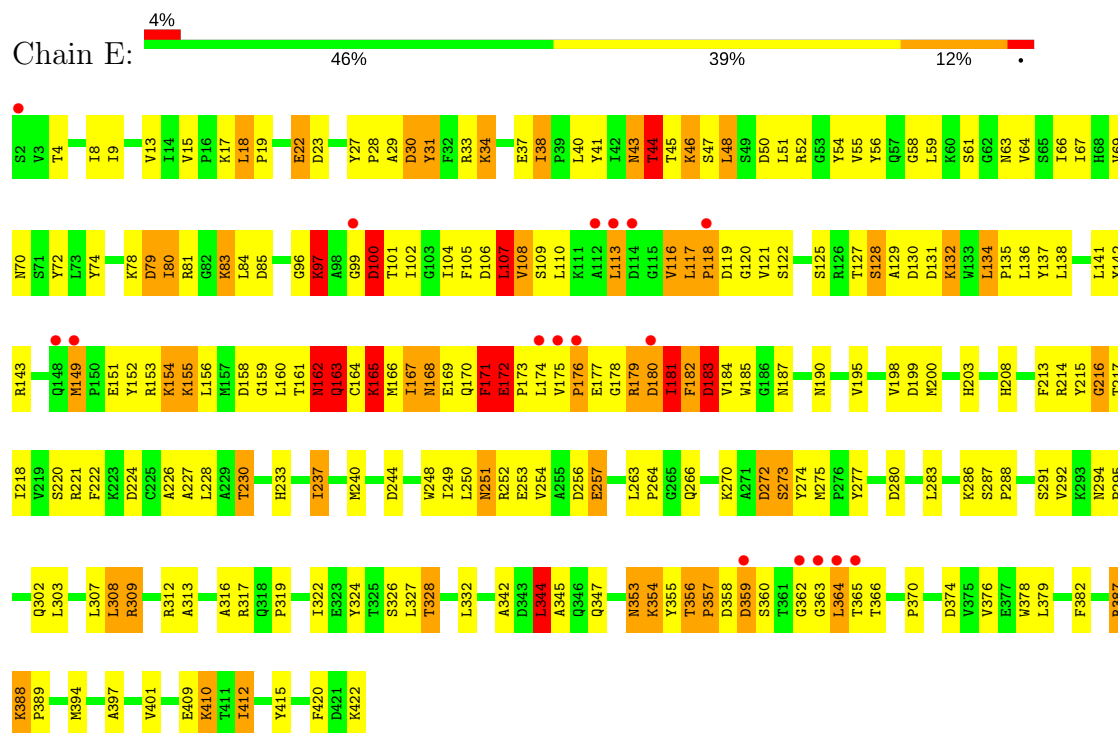
### • Molecule 1: Nucleoprotein



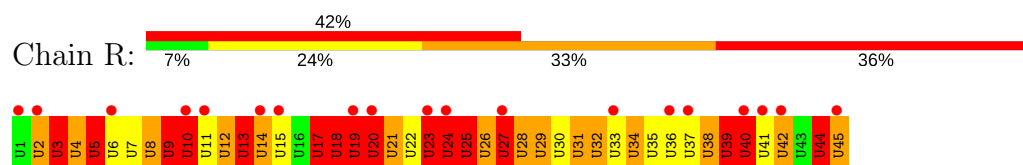




● Molecule 1: Nucleoprotein



● Molecule 2: RNA (45-MER)



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	165.48Å 235.06Å 75.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.77 – 3.00 40.88 – 2.85	Depositor EDS
% Data completeness (in resolution range)	83.0 (34.77-3.00) 74.2 (40.88-2.85)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.85 (at 2.86Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6_289)	Depositor
R, $R_{free}$	0.251 , 0.288 0.246 , 0.290	Depositor DCC
$R_{free}$ test set	1846 reflections (3.69%)	DCC
Wilson B-factor (Å <sup>2</sup> )	73.9	Xtriage
Anisotropy	0.545	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 47.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	17437	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.51% 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: IUM

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	0.56	2/3403 (0.1%)	0.67	0/4607
1	B	0.54	1/3365 (0.0%)	0.67	1/4554 (0.0%)
1	C	0.55	0/3350	0.67	1/4533 (0.0%)
1	D	0.54	0/3373	0.67	0/4565
1	E	0.52	0/3403	0.65	0/4607
2	R	0.75	0/989	1.71	30/1526 (2.0%)
All	All	0.56	3/17883 (0.0%)	0.77	32/24392 (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	A	0	3
1	B	0	4
1	C	0	3
1	D	0	3
1	E	0	4
All	All	0	17

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	225	CYS	CB-SG	-5.97	1.72	1.81
1	B	179	ARG	C-N	5.42	1.46	1.34
1	A	30	ASP	CB-CG	5.20	1.62	1.51

The worst 5 of 32 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	R	13	U	O4'-C1'-N1	12.47	118.17	108.20
2	R	9	U	O4'-C1'-N1	11.70	117.56	108.20
2	R	27	U	O4'-C1'-N1	11.37	117.30	108.20
2	R	39	U	O4'-C1'-N1	-11.05	99.36	108.20
2	R	18	U	O4'-C1'-N1	9.81	116.05	108.20

There are no chirality outliers.

5 of 17 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	107	LEU	Peptide
1	A	98	ALA	Peptide
1	A	99	GLY	Peptide
1	B	107	LEU	Peptide
1	B	97	LYS	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	3327	0	3287	269	0
1	B	3290	0	3253	260	0
1	C	3275	0	3242	279	0
1	D	3298	0	3264	266	0
1	E	3327	0	3287	247	0
2	R	900	0	451	85	0
3	A	4	0	0	0	0
3	B	3	0	0	0	0
3	C	3	0	0	0	0
3	D	2	0	0	0	0
3	E	3	0	0	0	0
3	R	5	0	0	0	0
All	All	17437	0	16784	1310	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 38.

The worst 5 of 1310 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:B:165:LYS:HB3	1:B:166:MET:SD	1.48	1.53
1:C:165:LYS:HB3	1:C:166:MET:SD	1.47	1.51
1:B:167:ILE:HD13	1:B:168:ASN:N	1.30	1.37
1:A:165:LYS:HB3	1:A:166:MET:SD	1.67	1.33
1:B:165:LYS:CB	1:B:166:MET:SD	2.28	1.22

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	419/421 (100%)	316 (75%)	70 (17%)	33 (8%)	1	5
1	B	411/421 (98%)	315 (77%)	67 (16%)	29 (7%)	1	6
1	C	409/421 (97%)	318 (78%)	62 (15%)	29 (7%)	1	6
1	D	412/421 (98%)	319 (77%)	61 (15%)	32 (8%)	1	5
1	E	419/421 (100%)	317 (76%)	67 (16%)	35 (8%)	1	4
All	All	2070/2105 (98%)	1585 (77%)	327 (16%)	158 (8%)	1	6

5 of 158 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	34	LYS
1	A	99	GLY
1	A	117	LEU
1	A	120	GLY
1	A	132	LYS

### 5.3.2 Protein sidechains [i](#)

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	362/362 (100%)	301 (83%)	61 (17%)	2	12
1	B	358/362 (99%)	296 (83%)	62 (17%)	2	12
1	C	356/362 (98%)	298 (84%)	58 (16%)	3	13
1	D	359/362 (99%)	296 (82%)	63 (18%)	2	11
1	E	362/362 (100%)	298 (82%)	64 (18%)	2	11
All	All	1797/1810 (99%)	1489 (83%)	308 (17%)	2	12

5 of 308 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	100	ASP
1	C	328	THR
1	E	187	ASN
1	C	134	LEU
1	C	181	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 45 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	57	GLN
1	C	203	HIS
1	E	203	HIS
1	C	162	ASN
1	C	266	GLN

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	R	44/45 (97%)	27 (61%)	0

5 of 27 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	R	2	U
2	R	3	U

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Mol	Chain	Res	Type
2	R	5	U
2	R	9	U
2	R	10	U

There are no RNA pucker outliers to report.

## 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 20 ligands modelled in this entry, 20 are modelled with single atom - 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 [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	421/421 (100%)	-0.08	16 (3%) 41 17	50, 81, 131, 148	0
1	B	415/421 (98%)	-0.15	9 (2%) 62 33	50, 81, 130, 140	0
1	C	413/421 (98%)	-0.08	13 (3%) 49 22	50, 81, 129, 141	0
1	D	416/421 (98%)	0.06	20 (4%) 31 12	50, 81, 129, 141	0
1	E	421/421 (100%)	-0.07	17 (4%) 39 16	49, 82, 131, 145	0
2	R	45/45 (100%)	1.86	19 (42%) 0 0	106, 124, 140, 151	0
All	All	2131/2150 (99%)	-0.02	94 (4%) 35 14	49, 82, 131, 151	0

The worst 5 of 94 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	361	THR	7.9
1	C	2	SER	7.7
1	E	2	SER	6.9
1	D	166	MET	6.8
1	B	2	SER	6.5

### 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
3	IUM	E	534	1/3	0.99	0.12	-0.94	105,105,105,105	0
3	IUM	B	526	1/3	0.99	0.12	-1.16	104,104,104,104	0
3	IUM	D	532	1/3	1.00	0.11	-1.60	100,100,100,100	0
3	IUM	A	522	1/3	0.99	0.10	-1.64	109,109,109,109	0
3	IUM	C	529	1/3	0.99	0.09	-2.67	105,105,105,105	0
3	IUM	A	521	1/3	0.96	0.04	-	162,162,162,162	0
3	IUM	E	535	1/3	0.98	0.06	-	186,186,186,186	0
3	IUM	R	540	1/3	0.69	0.20	-	257,257,257,257	0
3	IUM	B	527	1/3	0.94	0.05	-	199,199,199,199	0
3	IUM	A	523	1/3	0.95	0.04	-	174,174,174,174	0
3	IUM	C	528	1/3	0.97	0.03	-	138,138,138,138	0
3	IUM	C	530	1/3	0.99	0.04	-	151,151,151,151	0
3	IUM	R	538	1/3	0.56	0.07	-	272,272,272,272	0
3	IUM	R	536	1/3	0.91	0.04	-	245,245,245,245	0
3	IUM	A	524	1/3	0.99	0.04	-	158,158,158,158	0
3	IUM	R	537	1/3	0.90	0.12	-	258,258,258,258	0
3	IUM	B	525	1/3	0.90	0.09	-	187,187,187,187	0
3	IUM	D	531	1/3	0.89	0.08	-	188,188,188,188	0
3	IUM	E	533	1/3	0.91	0.03	-	173,173,173,173	0
3	IUM	R	539	1/3	0.88	0.14	-	265,265,265,265	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.