



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 04:53 pm GMT

PDB ID : 3LWR  
Title : Structure of H/ACA RNP bound to a substrate RNA containing 4SU  
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Deposited on : 2010-02-24  
Resolution : 2.20 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
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

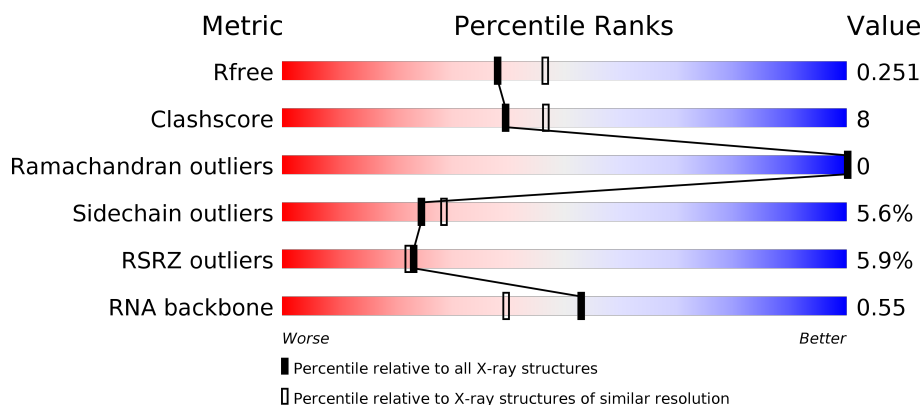
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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))
$R_{free}$	100719	4002 (2.20-2.20)
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)
RSRZ outliers	101464	4033 (2.20-2.20)
RNA backbone	2435	1007 (2.74-1.66)

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. 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	340	<div> <div>74%</div> <div>18%</div> <div>• 7%</div> </div>
2	B	60	<div> <div>3%</div> <div>62%</div> <div>20%</div> <div>7%</div> <div>12%</div> </div>
3	C	123	<div> <div>8%</div> <div>79%</div> <div>17%</div> <div>• •</div> </div>
4	D	58	<div> <div>10%</div> <div>64%</div> <div>21%</div> <div>12%</div> <div>•</div> </div>

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Mol	Chain	Length	Quality of chain
5	E	13	<div> <div>85%</div> <div> <div></div> <div>46%</div> <div>31%</div> <div>8%</div> <div>8%</div> <div>8%</div> </div> </div>

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
6	ZN	B	61	-	-	-	X

## 2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 5432 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 Pseudouridine synthase Cbf5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	317	Total	C	N	O	S	0	0	0
			2521	1629	432	450	10			

- Molecule 2 is a protein called Ribosome biogenesis protein Nop10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	53	Total	C	N	O	S	0	0	0
			444	283	86	71	4			

- Molecule 3 is a protein called 50S ribosomal protein L7Ae.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	120	Total	C	N	O	S	0	0	0
			916	585	151	177	3			

- Molecule 4 is a RNA chain called H/ACA RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	58	Total	C	N	O	P	0	0	0
			1236	552	227	400	57			

- Molecule 5 is a RNA chain called 5'-R(\*GP\*AP\*GP\*CP\*GP\*(4SU)P\*GP\*CP\*GP\*GP\*UP\*UP\*U)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	12	Total	C	N	O	P	0	0	0
			257	115	47	84	11			

- Molecule 6 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total 1	Zn 1	0	0

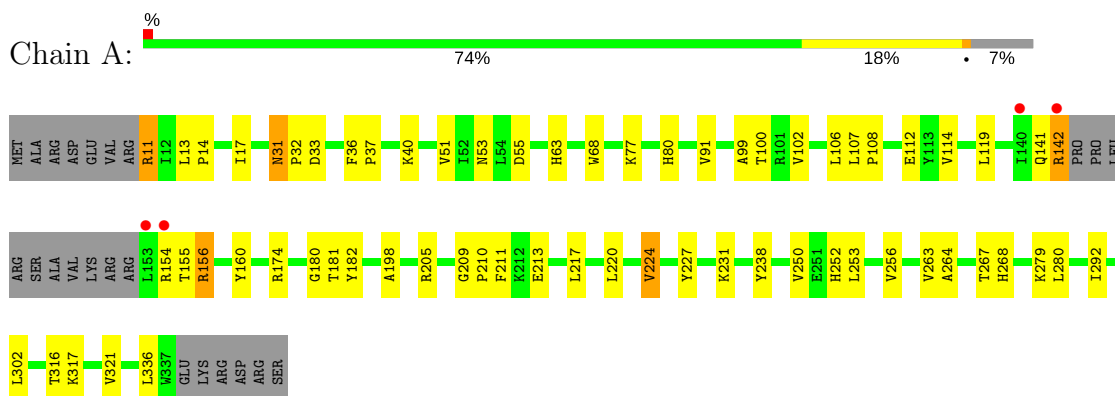
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	32	Total 32	O 32	0	0
7	B	8	Total 8	O 8	0	0
7	C	2	Total 2	O 2	0	0
7	D	15	Total 15	O 15	0	0

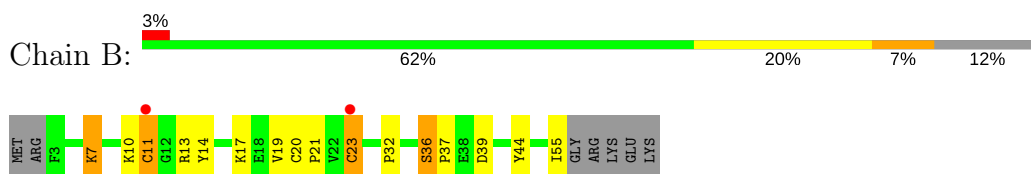
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

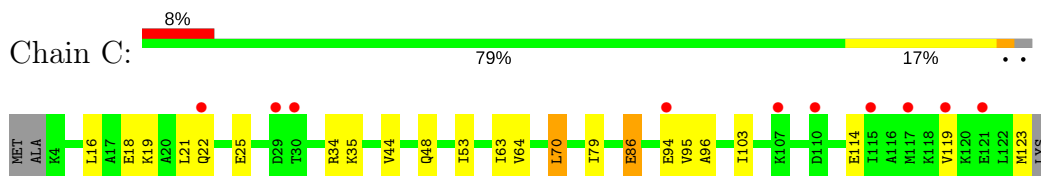
#### • Molecule 1: Pseudouridine synthase Cbf5



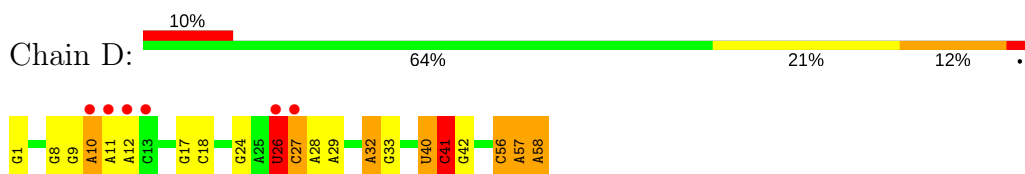
#### • Molecule 2: Ribosome biogenesis protein Nop10



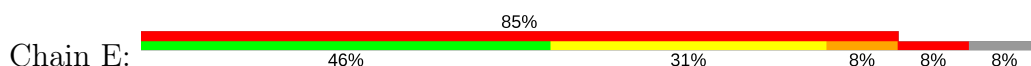
#### • Molecule 3: 50S ribosomal protein L7Ae



#### • Molecule 4: H/ACA RNA



#### • Molecule 5: 5'-R(\*GP\*AP\*GP\*CP\*GP\*(4SU)P\*GP\*CP\*GP\*GP\*UP\*UP\*U)-3'





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	187.35Å 64.43Å 83.39Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.69 – 2.20 41.69 – 2.20	Depositor EDS
% Data completeness (in resolution range)	52.4 (41.69-2.20) 99.4 (41.69-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.11	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.10 (at 2.20Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, $R_{free}$	0.222 , 0.259 0.214 , 0.251	Depositor DCC
$R_{free}$ test set	2631 reflections (5.10%)	DCC
Wilson B-factor (Å <sup>2</sup> )	32.6	Xtriage
Anisotropy	0.255	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 45.5	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	0.94	EDS
Total number of atoms	5432	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.36% 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: ZN

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	0/2578	0.57	0/3488
2	B	0.75	2/457 (0.4%)	0.59	0/613
3	C	0.34	0/928	0.50	0/1253
4	D	0.52	0/1381	1.15	8/2152 (0.4%)
5	E	0.48	0/286	1.09	1/443 (0.2%)
All	All	0.47	2/5630 (0.0%)	0.80	9/7949 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	23	CYS	CB-SG	8.68	1.97	1.82
2	B	11	CYS	CB-SG	7.63	1.95	1.82

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	26	U	O4'-C1'-N1	-8.77	101.18	108.20
4	D	24	G	P-O3'-C3'	7.94	129.22	119.70
4	D	41	C	N1-C1'-C2'	-7.55	103.70	112.00
4	D	1	G	O5'-C5'-C4'	6.78	124.58	111.70
4	D	40	U	P-O3'-C3'	5.78	126.63	119.70
5	E	13	G	C3'-C2'-C1'	5.77	106.11	101.50
4	D	56	C	P-O3'-C3'	5.50	126.30	119.70
4	D	32	A	C3'-C2'-C1'	5.35	105.78	101.50
4	D	41	C	C3'-C2'-C1'	5.06	105.55	101.50

There are no chirality outliers.

There are no planarity outliers.

## 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	2521	0	2588	42	0
2	B	444	0	451	13	0
3	C	916	0	964	15	0
4	D	1236	0	633	17	0
5	E	257	0	132	10	0
6	B	1	0	0	0	0
7	A	32	0	0	0	0
7	B	8	0	0	0	0
7	C	2	0	0	0	0
7	D	15	0	0	0	0
All	All	5432	0	4768	85	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.

All (85) 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)
4:D:32:A:H2'	4:D:33:G:C8	2.09	0.88
4:D:41:C:H42	5:E:9:G:H1	1.22	0.84
1:A:141:GLN:HE22	1:A:182:TYR:H	1.32	0.77
1:A:68:TRP:CE2	2:B:37:PRO:HG3	2.22	0.75
2:B:11:CYS:SG	2:B:13:ARG:HB3	2.28	0.74
3:C:18:GLU:O	3:C:22:GLN:HG2	1.97	0.64
1:A:31:ASN:C	1:A:31:ASN:HD22	2.02	0.63
1:A:11:ARG:N	1:A:11:ARG:HD2	2.13	0.63
1:A:63:HIS:HD2	1:A:80:HIS:NE2	1.97	0.62
1:A:220:LEU:O	1:A:224:VAL:HG13	2.01	0.61
3:C:44:VAL:HG11	3:C:70:LEU:HD13	1.82	0.61
1:A:114:VAL:HG22	1:A:174:ARG:HG2	1.84	0.60
1:A:63:HIS:HE1	5:E:12:C:OP2	1.84	0.60
2:B:44:TYR:CZ	3:C:70:LEU:HD23	2.36	0.60
3:C:44:VAL:CG1	3:C:70:LEU:HD13	2.33	0.58
1:A:31:ASN:ND2	1:A:33:ASP:H	2.01	0.57
4:D:41:C:N3	5:E:9:G:N2	2.48	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:ASN:HD22	1:A:32:PRO:N	2.03	0.56
4:D:40:U:C4	5:E:12:C:N4	2.73	0.56
4:D:32:A:H2'	4:D:33:G:H8	1.66	0.55
4:D:26:U:H5''	4:D:26:U:C6	2.41	0.55
4:D:10:A:H2'	4:D:11:A:C8	2.43	0.54
1:A:99:ALA:O	1:A:102:VAL:HG22	2.08	0.54
4:D:26:U:H6	4:D:26:U:H5''	1.72	0.53
4:D:40:U:H5''	4:D:41:C:OP1	2.08	0.53
1:A:112:GLU:OE1	1:A:160:TYR:OH	2.27	0.53
5:E:12:C:HO2'	5:E:13:G:H8	1.55	0.53
1:A:156:ARG:HG2	1:A:181:THR:OG1	2.08	0.52
3:C:119:VAL:O	3:C:123:MET:HG2	2.09	0.52
5:E:12:C:O2'	5:E:13:G:H8	1.92	0.52
1:A:264:ALA:O	1:A:268:HIS:HD2	1.92	0.52
1:A:142:ARG:HA	1:A:142:ARG:HH11	1.75	0.52
1:A:107:LEU:N	1:A:108:PRO:CD	2.72	0.51
1:A:119:LEU:HD23	1:A:198:ALA:HB2	1.93	0.51
4:D:58:A:H2'	4:D:58:A:N3	2.26	0.51
1:A:107:LEU:HB2	1:A:108:PRO:HD3	1.93	0.51
1:A:77:LYS:HE2	1:A:100:THR:HG21	1.93	0.49
1:A:141:GLN:NE2	1:A:182:TYR:CD1	2.80	0.49
3:C:53:ILE:HD12	3:C:53:ILE:N	2.27	0.49
3:C:34:ARG:HG3	3:C:103:ILE:HD11	1.94	0.49
3:C:16:LEU:HG	3:C:79:ILE:HD11	1.95	0.49
3:C:95:VAL:CG1	3:C:96:ALA:N	2.76	0.49
4:D:10:A:H2'	4:D:11:A:H8	1.75	0.48
4:D:17:G:H2'	4:D:18:C:O4'	2.13	0.48
2:B:7:LYS:HE3	2:B:14:TYR:CZ	2.49	0.48
1:A:31:ASN:C	1:A:31:ASN:ND2	2.68	0.47
1:A:91:VAL:O	1:A:205:ARG:HD2	2.15	0.47
4:D:41:C:H2'	4:D:42:G:O4'	2.14	0.47
1:A:263:VAL:O	1:A:267:THR:HG23	2.15	0.46
3:C:63:ILE:HG22	3:C:64:VAL:HG13	1.98	0.46
1:A:238:TYR:CE1	2:B:55:ILE:HD11	2.51	0.46
1:A:55:ASP:CG	2:B:32:PRO:HG3	2.36	0.46
1:A:68:TRP:NE1	2:B:37:PRO:HG3	2.31	0.46
1:A:213:GLU:HG2	1:A:217:LEU:HD22	1.98	0.45
4:D:57:A:N3	4:D:57:A:H2'	2.32	0.45
2:B:17:LYS:HD2	2:B:21:PRO:HG3	1.98	0.45
5:E:11:G:H4'	5:E:12:C:OP1	2.16	0.45
1:A:14:PRO:O	1:A:17:ILE:HG12	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:ASN:HD22	1:A:32:PRO:CD	2.29	0.45
3:C:95:VAL:HG12	3:C:96:ALA:O	2.16	0.45
2:B:36:SER:HB2	2:B:39:ASP:H	1.81	0.44
2:B:7:LYS:HE2	2:B:7:LYS:HB2	1.51	0.44
1:A:292:ILE:HD11	1:A:321:VAL:HG11	2.00	0.44
3:C:19:LYS:HE3	3:C:114:GLU:OE2	2.17	0.44
3:C:79:ILE:HD12	3:C:119:VAL:CG1	2.47	0.44
1:A:210:PRO:HG3	1:A:252:HIS:ND1	2.33	0.44
1:A:250:VAL:HG22	1:A:253:LEU:HD12	1.99	0.44
4:D:8:G:O2'	4:D:9:G:H5'	2.18	0.44
1:A:40:LYS:O	1:A:40:LYS:HG2	2.18	0.44
2:B:20:CYS:HA	2:B:21:PRO:HD3	1.87	0.43
3:C:94:GLU:HB3	4:D:27:C:O2	2.18	0.43
1:A:114:VAL:HG21	2:B:14:TYR:CD1	2.54	0.43
1:A:227:TYR:CZ	1:A:231:LYS:HG3	2.54	0.42
1:A:142:ARG:HA	1:A:142:ARG:NH1	2.34	0.42
1:A:51:VAL:HG11	1:A:211:PHE:CZ	2.54	0.42
3:C:21:LEU:HD11	3:C:86:GLU:CD	2.40	0.42
2:B:10:LYS:HB2	2:B:10:LYS:HE2	1.69	0.42
5:E:12:C:O2'	5:E:13:G:C8	2.70	0.42
1:A:154:ARG:NH2	1:A:180:GLY:O	2.53	0.41
1:A:36:PHE:HA	1:A:37:PRO:HD2	1.84	0.41
1:A:141:GLN:NE2	1:A:182:TYR:HD1	2.19	0.41
1:A:209:GLY:HA3	1:A:210:PRO:HD3	1.88	0.40
1:A:316:THR:O	1:A:317:LYS:HG3	2.20	0.40
4:D:41:C:N4	5:E:9:G:H1	2.04	0.40
5:E:11:G:H4'	5:E:12:C:O5'	2.19	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	313/340 (92%)	309 (99%)	4 (1%)	0	100	100
2	B	51/60 (85%)	49 (96%)	2 (4%)	0	100	100
3	C	118/123 (96%)	117 (99%)	1 (1%)	0	100	100
All	All	482/523 (92%)	475 (98%)	7 (2%)	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	268/289 (93%)	254 (95%)	14 (5%)	27	32
2	B	48/54 (89%)	44 (92%)	4 (8%)	13	13
3	C	97/99 (98%)	92 (95%)	5 (5%)	27	32
All	All	413/442 (93%)	390 (94%)	23 (6%)	25	29

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ARG
1	A	13	LEU
1	A	31	ASN
1	A	53	ASN
1	A	106	LEU
1	A	142	ARG
1	A	155	THR
1	A	156	ARG
1	A	224	VAL
1	A	256	VAL
1	A	279	LYS
1	A	280	LEU
1	A	302	LEU
1	A	336	LEU
2	B	7	LYS

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Mol	Chain	Res	Type
2	B	19	VAL
2	B	23	CYS
2	B	36	SER
3	C	25	GLU
3	C	35	LYS
3	C	48	GLN
3	C	70	LEU
3	C	86	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	ASN
1	A	46	HIS
1	A	53	ASN
1	A	63	HIS
1	A	141	GLN
1	A	221	HIS
1	A	268	HIS
3	C	38	ASN
3	C	48	GLN

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
4	D	57/58 (98%)	10 (17%)	0
5	E	10/13 (76%)	4 (40%)	0
All	All	67/71 (94%)	14 (20%)	0

All (14) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
4	D	10	A
4	D	12	A
4	D	26	U
4	D	27	C
4	D	28	A
4	D	29	A
4	D	41	C
4	D	56	C

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Mol	Chain	Res	Type
4	D	57	A
4	D	58	A
5	E	12	C
5	E	13	G
5	E	14	G
5	E	15	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 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	317/340 (93%)	-0.08	4 (1%) 77 75	14, 27, 58, 98	0
2	B	53/60 (88%)	0.06	2 (3%) 41 39	20, 33, 53, 74	0
3	C	120/123 (97%)	0.38	10 (8%) 12 11	26, 49, 86, 99	0
4	D	58/58 (100%)	0.18	6 (10%) 7 6	29, 45, 126, 132	0
5	E	12/13 (92%)	3.38	11 (91%) 0 0	89, 120, 135, 156	0
All	All	560/594 (94%)	0.13	33 (5%) 23 22	14, 34, 89, 156	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
5	E	17	U	7.7
5	E	5	G	4.5
3	C	30	THR	4.5
5	E	9	G	4.5
4	D	11	A	4.3
4	D	12	A	4.1
3	C	117	MET	3.9
4	D	10	A	3.9
4	D	27	C	3.8
1	A	142	ARG	3.7
5	E	14	G	3.7
1	A	153	LEU	3.7
5	E	16	U	3.7
2	B	11	CYS	3.4
1	A	154	ARG	3.2
3	C	107	LYS	3.1
5	E	12	C	3.1
5	E	13	G	3.0
3	C	94	GLU	2.8
3	C	121	GLU	2.8

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Mol	Chain	Res	Type	RSRZ
4	D	13	C	2.8
1	A	140	ILE	2.7
4	D	26	U	2.6
3	C	29	ASP	2.4
3	C	119	VAL	2.3
5	E	11	G	2.3
3	C	110	ASP	2.3
5	E	6	A	2.3
5	E	15	U	2.3
2	B	23	CYS	2.2
3	C	22	GLN	2.2
3	C	115	ILE	2.1
5	E	7	G	2.0

## 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
6	ZN	B	61	1/1	0.38	1.82	24.13	236,236,236,236	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.