



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 21, 2020 – 04:14 PM BST

PDB ID : 1GTF
Title : The structure of the trp RNA-binding attenuation protein (TRAP) bound to a 53-nucleotide RNA molecule containing GAGUU repeats
Authors : Hopcroft, N.H.; Wendt, A.L.; Gollnick, P.; Antson, A.A.
Deposited on : 2002-01-15
Resolution : 1.75 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
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

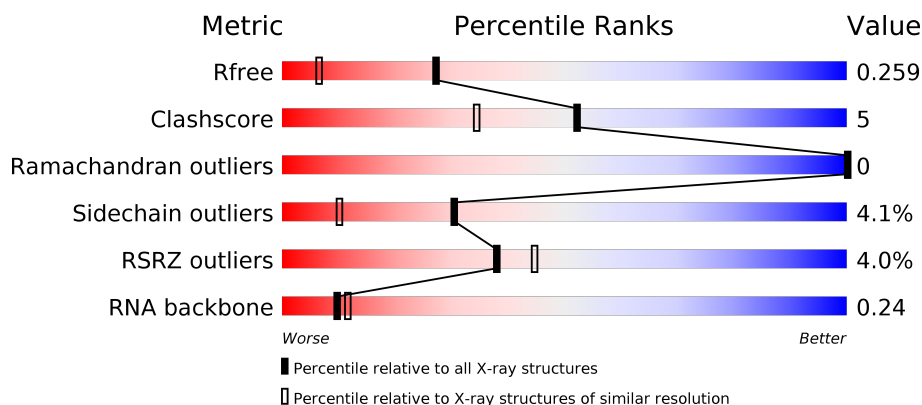
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)
RNA backbone	3102	1079 (2.40-1.12)

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 ≥ 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	74	<div> <div>5%</div> <div> <div></div> <div>85%</div> <div>7%</div> <div>• 7%</div> </div> </div>
1	B	74	<div> <div>80%</div> <div>12%</div> <div>8%</div> </div>
1	C	74	<div> <div>12%</div> <div>80%</div> <div>12%</div> <div>• 5%</div> </div>
1	D	74	<div> <div>11%</div> <div>77%</div> <div>15%</div> <div>• 7%</div> </div>

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Mol	Chain	Length	Quality of chain
1	E	74	
1	F	74	
1	G	74	
1	H	74	
1	I	74	
1	J	74	
1	K	74	
1	L	74	
1	M	74	
1	N	74	
1	O	74	
1	P	74	
1	Q	74	
1	R	74	
1	S	74	
1	T	74	
1	U	74	
1	V	74	
2	W	55	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14607 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 TRP RNA-BINDING ATTENUATION PROTEIN (TRAP).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	69	Total	C	N	O	15	0	0
			536	336	98	102			
1	B	68	Total	C	N	O	16	0	0
			527	330	96	101			
1	C	70	Total	C	N	O	13	0	0
			544	340	100	104			
1	D	69	Total	C	N	O	26	0	0
			536	336	98	102			
1	E	68	Total	C	N	O	21	0	0
			527	330	96	101			
1	F	69	Total	C	N	O	30	0	0
			536	336	98	102			
1	G	70	Total	C	N	O	26	0	0
			544	340	100	104			
1	H	68	Total	C	N	O	16	0	0
			527	330	96	101			
1	I	69	Total	C	N	O	23	0	0
			536	336	98	102			
1	J	67	Total	C	N	O	13	0	0
			523	328	95	100			
1	K	69	Total	C	N	O	21	0	0
			536	336	98	102			
1	L	70	Total	C	N	O	13	0	0
			542	338	99	105			
1	M	71	Total	C	N	O	14	0	0
			551	344	101	106			
1	N	70	Total	C	N	O	4	0	0
			542	338	99	105			
1	O	70	Total	C	N	O	0	0	0
			542	338	99	105			
1	P	70	Total	C	N	O	6	0	0
			542	338	99	105			

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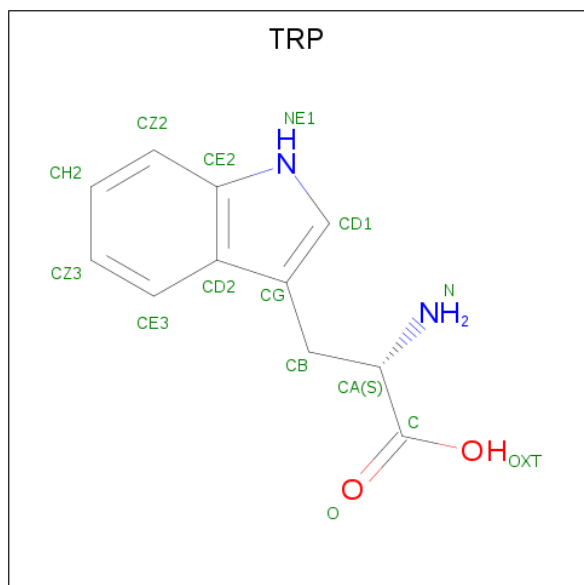
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	Q	70	Total	C	N	O	3	0	0
			542	338	99	105			
1	R	70	Total	C	N	O	17	0	0
			542	338	99	105			
1	S	70	Total	C	N	O	8	0	0
			542	338	99	105			
1	T	70	Total	C	N	O	7	0	0
			542	338	99	105			
1	U	70	Total	C	N	O	10	0	0
			542	338	99	105			
1	V	70	Total	C	N	O	7	0	0
			542	338	99	105			

- Molecule 2 is a RNA chain called (GAGUU)10GAG 53-NUCLEOTIDE RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	W	44	Total	C	N	O	P	0	0	0
			968	429	187	308	44			

- Molecule 3 is TRYPTOPHAN (three-letter code: TRP) (formula: $C_{11}H_{12}N_2O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			15	11	2	2		
3	B	1	Total	C	N	O	0	0
			15	11	2	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total 15	C 11	N 2	O 2	0	0
3	D	1	Total 15	C 11	N 2	O 2	0	0
3	E	1	Total 15	C 11	N 2	O 2	0	0
3	F	1	Total 15	C 11	N 2	O 2	0	0
3	G	1	Total 15	C 11	N 2	O 2	0	0
3	H	1	Total 15	C 11	N 2	O 2	0	0
3	I	1	Total 15	C 11	N 2	O 2	0	0
3	J	1	Total 15	C 11	N 2	O 2	0	0
3	K	1	Total 15	C 11	N 2	O 2	0	0
3	L	1	Total 15	C 11	N 2	O 2	0	0
3	M	1	Total 15	C 11	N 2	O 2	0	0
3	N	1	Total 15	C 11	N 2	O 2	0	0
3	O	1	Total 15	C 11	N 2	O 2	0	0
3	P	1	Total 15	C 11	N 2	O 2	0	0
3	Q	1	Total 15	C 11	N 2	O 2	0	0
3	R	1	Total 15	C 11	N 2	O 2	0	0
3	S	1	Total 15	C 11	N 2	O 2	0	0
3	T	1	Total 15	C 11	N 2	O 2	0	0
3	U	1	Total 15	C 11	N 2	O 2	0	0
3	V	1	Total 15	C 11	N 2	O 2	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	94	Total O 94 94	0	0
4	B	69	Total O 69 69	0	0
4	C	89	Total O 89 89	0	0
4	D	70	Total O 70 70	0	0
4	E	63	Total O 63 63	0	0
4	F	60	Total O 60 60	0	0
4	G	64	Total O 64 64	0	0
4	H	62	Total O 62 62	0	0
4	I	68	Total O 68 68	0	0
4	J	71	Total O 71 71	0	0
4	K	82	Total O 82 82	0	0
4	L	44	Total O 44 44	0	0
4	M	51	Total O 51 51	0	0
4	N	46	Total O 46 46	0	0
4	O	49	Total O 49 49	0	0
4	P	54	Total O 54 54	0	0
4	Q	59	Total O 59 59	0	0
4	R	75	Total O 75 75	0	0
4	S	51	Total O 51 51	0	0
4	T	63	Total O 63 63	0	0
4	U	51	Total O 51 51	0	0
4	V	54	Total O 54 54	0	0

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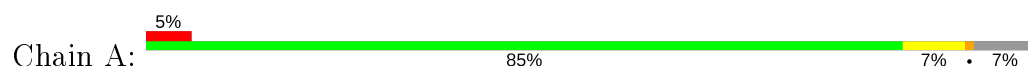
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	W	77	Total	O	0	0
			77	77		

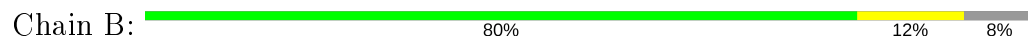
3 Residue-property plots [i](#)

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.

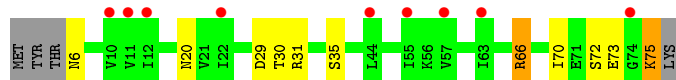
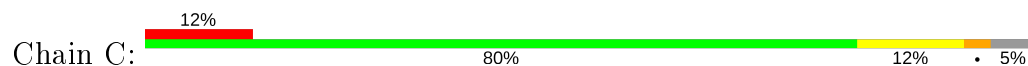
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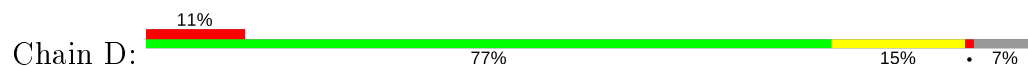
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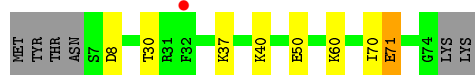
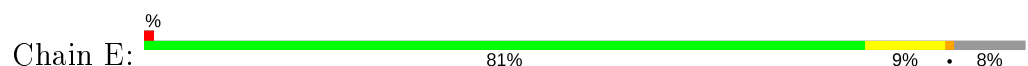
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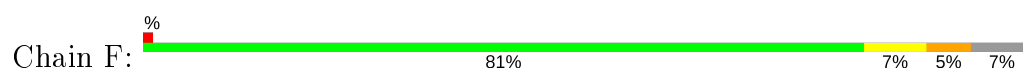
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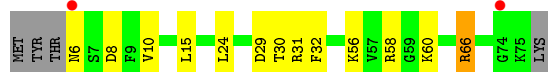
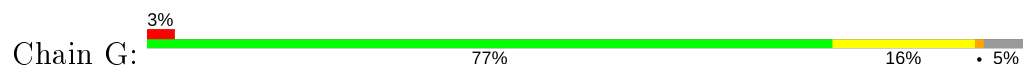
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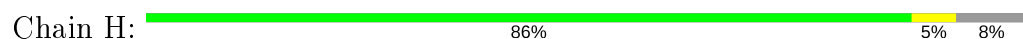
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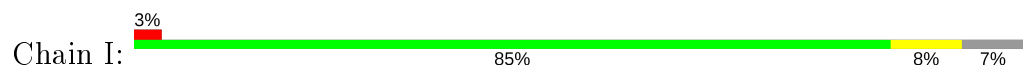
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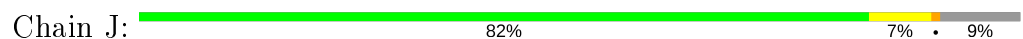
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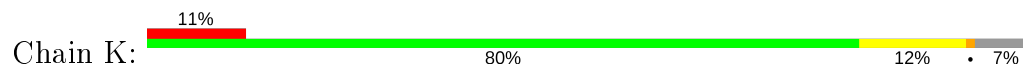
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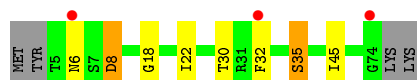
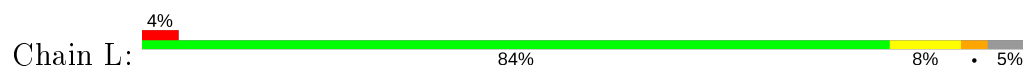
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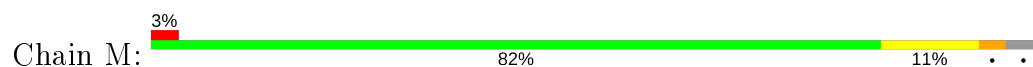
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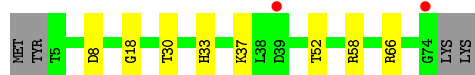
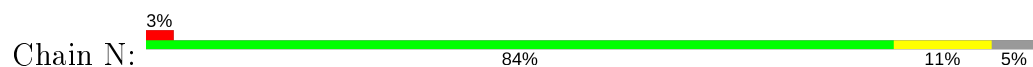
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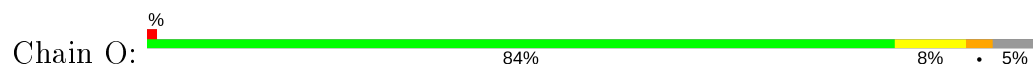
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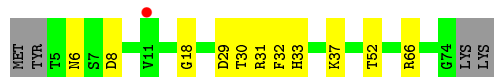
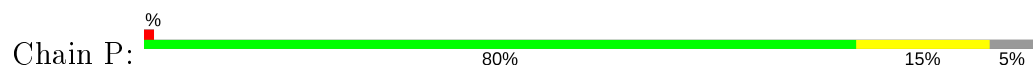
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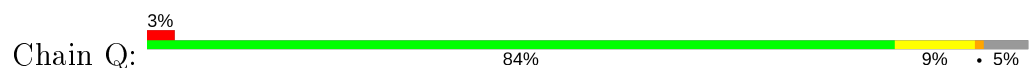
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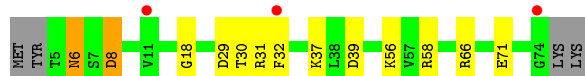
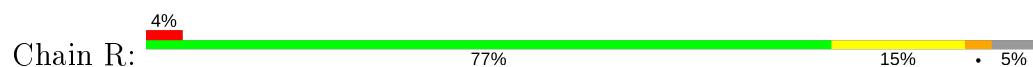
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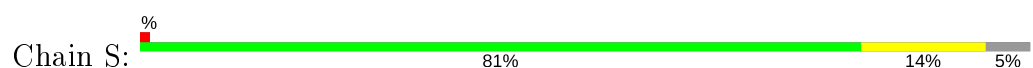
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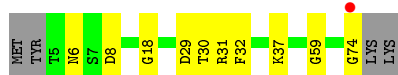


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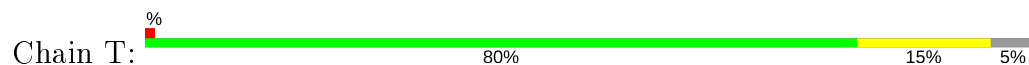


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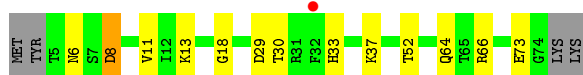
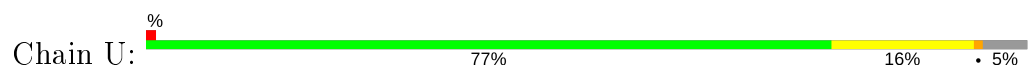




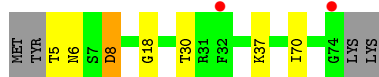
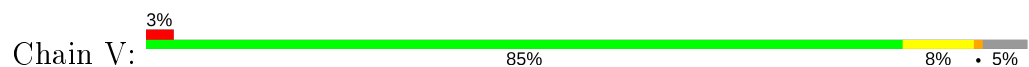
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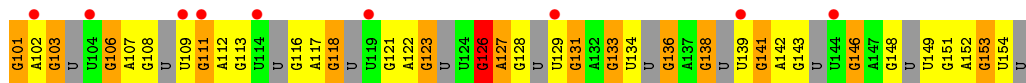
- Molecule 1: TRP RNA-BINDING ATTENUATION PROTEIN (TRAP)



- Molecule 1: TRP RNA-BINDING ATTENUATION PROTEIN (TRAP)



- Molecule 2: (GAGUU)10GAG 53-NUCLEOTIDE RNA



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	142.08Å 111.49Å 138.23Å 90.00° 117.28° 90.00°	Depositor
Resolution (Å)	47.67 – 1.75 47.93 – 1.75	Depositor EDS
% Data completeness (in resolution range)	100.0 (47.67-1.75) 95.9 (47.93-1.75)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.38 (at 1.75Å)	Xtriage
Refinement program	REFMAC 5.1.07	Depositor
R, R_{free}	0.194 , 0.242 0.234 , 0.259	Depositor DCC
R_{free} test set	1848 reflections (1.00%)	wwPDB-VP
Wilson B-factor (Å ²)	24.9	Xtriage
Anisotropy	0.281	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 50.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	14607	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.54% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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

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	1.16	3/543 (0.6%)	1.07	3/728 (0.4%)
1	B	1.15	4/534 (0.7%)	1.10	4/717 (0.6%)
1	C	1.09	3/551 (0.5%)	1.29	4/739 (0.5%)
1	D	1.00	2/543 (0.4%)	1.08	4/728 (0.5%)
1	E	1.19	3/534 (0.6%)	0.98	2/717 (0.3%)
1	F	1.32	5/543 (0.9%)	1.17	7/728 (1.0%)
1	G	1.03	2/551 (0.4%)	1.05	5/739 (0.7%)
1	H	1.02	1/534 (0.2%)	0.97	1/717 (0.1%)
1	I	0.95	1/543 (0.2%)	1.12	7/728 (1.0%)
1	J	1.32	2/530 (0.4%)	0.98	4/712 (0.6%)
1	K	1.15	3/543 (0.6%)	1.03	4/728 (0.5%)
1	L	0.74	0/549	0.90	1/738 (0.1%)
1	M	0.93	2/558 (0.4%)	0.91	1/749 (0.1%)
1	N	0.82	1/549 (0.2%)	0.84	0/738
1	O	0.76	0/549	0.96	0/738
1	P	0.82	0/549	0.90	0/738
1	Q	0.82	0/549	0.93	2/738 (0.3%)
1	R	0.88	1/549 (0.2%)	1.02	2/738 (0.3%)
1	S	0.85	1/549 (0.2%)	0.96	0/738
1	T	0.82	1/549 (0.2%)	0.93	0/738
1	U	0.84	1/549 (0.2%)	0.90	1/738 (0.1%)
1	V	0.75	1/549 (0.2%)	0.85	0/738
2	W	2.61	11/1078 (1.0%)	1.90	35/1661 (2.1%)
All	All	1.21	48/13075 (0.4%)	1.12	87/17771 (0.5%)

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	D	0	1

All (48) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	W	153	G	O3'-P	30.19	1.97	1.61
2	W	103	G	O3'-P	27.15	1.93	1.61
2	W	108	G	O3'-P	26.71	1.93	1.61
2	W	118	G	O3'-P	25.44	1.91	1.61
2	W	123	G	O3'-P	23.34	1.89	1.61
2	W	143	G	O3'-P	23.05	1.88	1.61
2	W	128	G	O3'-P	22.82	1.88	1.61
2	W	138	G	O3'-P	21.29	1.86	1.61
2	W	148	G	O3'-P	20.91	1.86	1.61
2	W	113	G	O3'-P	20.02	1.85	1.61
1	J	29	ASP	CB-CG	-19.90	1.09	1.51
2	W	133	G	O3'-P	17.78	1.82	1.61
1	F	66	ARG	NE-CZ	17.45	1.55	1.33
1	E	71	GLU	CG-CD	15.29	1.74	1.51
1	B	7	SER	CB-OG	-13.60	1.24	1.42
1	K	66	ARG	NE-CZ	12.41	1.49	1.33
1	H	73	GLU	CB-CG	12.10	1.75	1.52
1	A	66	ARG	NE-CZ	11.59	1.48	1.33
1	A	29	ASP	CB-CG	11.54	1.75	1.51
1	G	66	ARG	NE-CZ	11.33	1.47	1.33
1	A	39	ASP	CB-CG	11.12	1.75	1.51
1	F	73	GLU	CG-CD	11.11	1.68	1.51
1	M	73	GLU	CG-CD	-10.54	1.36	1.51
1	D	31	ARG	NE-CZ	10.15	1.46	1.33
1	K	31	ARG	CG-CD	10.05	1.77	1.51
1	J	31	ARG	CD-NE	9.95	1.63	1.46
1	E	60	LYS	CE-NZ	9.90	1.73	1.49
1	N	66	ARG	CD-NE	-9.81	1.29	1.46
1	F	39	ASP	CB-CG	9.77	1.72	1.51
1	G	8	ASP	CB-CG	8.92	1.70	1.51
1	I	73	GLU	CG-CD	8.45	1.64	1.51
1	B	39	ASP	CB-CG	8.43	1.69	1.51
1	C	66	ARG	CZ-NH1	8.11	1.43	1.33
1	U	29	ASP	CB-CG	-7.96	1.35	1.51
1	M	29	ASP	CB-CG	-7.64	1.35	1.51
1	F	29	ASP	CB-CG	-7.46	1.36	1.51
1	V	6	ASN	CB-CG	7.27	1.67	1.51
1	B	58	ARG	NE-CZ	6.62	1.41	1.33
1	S	6	ASN	CB-CG	6.59	1.66	1.51
1	T	66	ARG	CD-NE	-6.57	1.35	1.46
1	E	50	GLU	CG-CD	-6.08	1.42	1.51
1	K	29	ASP	CB-CG	5.81	1.64	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	60	LYS	CE-NZ	5.53	1.62	1.49
1	C	31	ARG	CD-NE	5.51	1.55	1.46
1	R	71	GLU	CG-CD	5.33	1.59	1.51
1	F	10	VAL	CB-CG2	-5.30	1.41	1.52
1	D	58	ARG	CB-CG	-5.15	1.38	1.52
1	C	66	ARG	CZ-NH2	-5.01	1.26	1.33

All (87) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	66	ARG	NE-CZ-NH2	15.20	127.90	120.30
1	F	66	ARG	CD-NE-CZ	-14.46	103.35	123.60
2	W	118	G	O4'-C1'-N9	12.22	117.98	108.20
2	W	126	G	P-O3'-C3'	11.70	133.74	119.70
1	C	66	ARG	NE-CZ-NH1	-11.62	114.49	120.30
2	W	103	G	O4'-C1'-N9	11.12	117.09	108.20
2	W	143	G	O4'-C1'-N9	11.02	117.01	108.20
2	W	113	G	O4'-C1'-N9	10.91	116.93	108.20
2	W	128	G	O4'-C1'-N9	10.03	116.22	108.20
2	W	133	G	O4'-C1'-N9	9.60	115.88	108.20
2	W	136	G	P-O3'-C3'	9.42	131.01	119.70
1	K	73	GLU	CB-CG-CD	9.32	139.37	114.20
2	W	153	G	O4'-C1'-N9	8.89	115.31	108.20
1	R	66	ARG	CG-CD-NE	-8.85	93.21	111.80
1	B	29	ASP	CB-CG-OD2	8.82	126.24	118.30
2	W	138	G	O4'-C1'-N9	8.77	115.22	108.20
1	A	39	ASP	CB-CG-OD1	8.71	126.14	118.30
2	W	131	G	P-O3'-C3'	8.56	129.97	119.70
1	D	31	ARG	NE-CZ-NH1	8.44	124.52	120.30
1	I	31	ARG	CD-NE-CZ	-8.29	112.00	123.60
1	B	58	ARG	NE-CZ-NH2	8.11	124.36	120.30
2	W	108	G	O4'-C1'-N9	7.82	114.45	108.20
1	H	29	ASP	CB-CG-OD1	7.72	125.25	118.30
2	W	143	G	P-O3'-C3'	-7.71	110.45	119.70
1	C	29	ASP	CB-CG-OD2	7.69	125.22	118.30
2	W	111	G	P-O3'-C3'	7.64	128.86	119.70
1	F	66	ARG	NE-CZ-NH2	-7.60	116.50	120.30
1	J	31	ARG	CG-CD-NE	-7.26	96.55	111.80
2	W	146	G	P-O3'-C3'	7.24	128.38	119.70
1	D	29	ASP	CB-CG-OD1	-7.05	111.96	118.30
1	G	29	ASP	CB-CG-OD2	7.04	124.64	118.30
2	W	123	G	O4'-C1'-N9	7.03	113.82	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	ARG	CD-NE-CZ	-7.02	113.77	123.60
1	E	60	LYS	CD-CE-NZ	-6.97	95.66	111.70
1	I	17	ASP	CB-CG-OD2	6.96	124.56	118.30
1	B	39	ASP	CB-CG-OD1	6.96	124.56	118.30
1	F	66	ARG	NE-CZ-NH1	6.66	123.63	120.30
1	I	31	ARG	NE-CZ-NH1	6.65	123.63	120.30
1	F	29	ASP	CB-CG-OD2	6.58	124.22	118.30
2	W	141	G	P-O3'-C3'	6.49	127.49	119.70
1	G	31	ARG	NE-CZ-NH1	6.33	123.47	120.30
1	D	31	ARG	NE-CZ-NH2	-6.32	117.14	120.30
1	F	39	ASP	CA-CB-CG	-6.19	99.77	113.40
1	B	58	ARG	NE-CZ-NH1	-6.15	117.22	120.30
1	G	31	ARG	NE-CZ-NH2	-6.08	117.26	120.30
1	F	31	ARG	CB-CG-CD	-6.06	95.85	111.60
1	I	31	ARG	NE-CZ-NH2	-6.00	117.30	120.30
2	W	106	G	P-O3'-C3'	5.93	126.82	119.70
1	K	66	ARG	CD-NE-CZ	-5.81	115.46	123.60
1	I	71	GLU	CG-CD-OE1	-5.79	106.72	118.30
1	M	66	ARG	CD-NE-CZ	-5.77	115.53	123.60
2	W	126	G	C5'-C4'-O4'	-5.76	102.18	109.10
1	J	29	ASP	CA-CB-CG	5.76	126.07	113.40
1	G	66	ARG	CD-NE-CZ	-5.67	115.66	123.60
2	W	148	G	O4'-C1'-N9	5.67	112.73	108.20
2	W	136	G	C5'-C4'-O4'	-5.66	102.31	109.10
2	W	128	G	O5'-P-OP2	-5.62	100.64	105.70
1	J	29	ASP	CB-CG-OD2	5.60	123.34	118.30
2	W	117	A	C2-N3-C4	5.55	113.37	110.60
2	W	126	G	N3-C2-N2	-5.55	116.02	119.90
1	F	39	ASP	CB-CG-OD1	5.54	123.29	118.30
1	C	31	ARG	CD-NE-CZ	5.53	131.35	123.60
2	W	141	G	OP1-P-OP2	5.53	127.90	119.60
1	Q	26	ARG	NE-CZ-NH2	-5.51	117.55	120.30
2	W	126	G	C5-C6-O6	-5.50	125.30	128.60
2	W	106	G	N3-C2-N2	-5.50	116.05	119.90
2	W	143	G	OP1-P-O3'	-5.49	93.12	105.20
1	Q	39	ASP	CB-CG-OD2	5.48	123.23	118.30
1	I	50	GLU	OE1-CD-OE2	-5.46	116.75	123.30
2	W	142	A	N1-C2-N3	-5.45	126.57	129.30
1	L	8	ASP	CB-CG-OD2	5.43	123.19	118.30
1	J	29	ASP	CB-CG-OD1	-5.42	113.42	118.30
2	W	101	G	N3-C2-N2	-5.42	116.11	119.90
1	A	39	ASP	CA-CB-CG	-5.39	101.53	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	10	VAL	CG1-CB-CG2	-5.38	102.30	110.90
2	W	113	G	N1-C6-O6	-5.37	116.68	119.90
2	W	131	G	N3-C2-N2	-5.31	116.18	119.90
1	K	17	ASP	CB-CG-OD1	5.30	123.07	118.30
2	W	117	A	N1-C2-N3	-5.29	126.66	129.30
1	I	50	GLU	CB-CG-CD	5.26	128.41	114.20
1	K	29	ASP	CB-CG-OD1	5.24	123.02	118.30
1	E	71	GLU	OE1-CD-OE2	-5.21	117.05	123.30
1	R	39	ASP	CB-CG-OD2	5.18	122.96	118.30
2	W	109	U	N1-C2-O2	5.16	126.41	122.80
1	U	66	ARG	CG-CD-NE	-5.12	101.06	111.80
1	D	31	ARG	CD-NE-CZ	-5.08	116.48	123.60
2	W	146	G	OP1-P-OP2	5.07	127.21	119.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	29	ASP	Sidechain

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	536	0	541	2	0
1	B	527	0	528	2	0
1	C	544	0	547	7	0
1	D	536	0	541	11	0
1	E	527	0	528	4	0
1	F	536	0	541	3	0
1	G	544	0	547	11	0
1	H	527	0	528	1	0
1	I	536	0	541	1	0
1	J	523	0	525	3	0
1	K	536	0	541	4	0
1	L	542	0	541	8	0
1	M	551	0	554	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	N	542	0	541	4	0
1	O	542	0	541	8	0
1	P	542	0	541	19	0
1	Q	542	0	541	14	0
1	R	542	0	541	9	0
1	S	542	0	541	6	0
1	T	542	0	541	5	0
1	U	542	0	541	7	0
1	V	542	0	541	6	0
2	W	968	0	484	26	0
3	A	15	0	9	1	0
3	B	15	0	9	1	0
3	C	15	0	9	1	0
3	D	15	0	9	1	0
3	E	15	0	9	1	0
3	F	15	0	9	1	0
3	G	15	0	9	1	0
3	H	15	0	9	1	0
3	I	15	0	9	1	0
3	J	15	0	9	1	0
3	K	15	0	9	1	0
3	L	15	0	9	1	0
3	M	15	0	9	1	0
3	N	15	0	9	1	0
3	O	15	0	9	1	0
3	P	15	0	9	1	0
3	Q	15	0	9	1	0
3	R	15	0	9	1	0
3	S	15	0	9	1	0
3	T	15	0	9	1	0
3	U	15	0	9	1	0
3	V	15	0	9	1	0
4	A	94	0	0	1	0
4	B	69	0	0	0	0
4	C	89	0	0	3	1
4	D	70	0	0	6	0
4	E	63	0	0	3	0
4	F	60	0	0	1	0
4	G	64	0	0	2	0
4	H	62	0	0	0	0
4	I	68	0	0	0	0
4	J	71	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	K	82	0	0	1	1
4	L	44	0	0	3	0
4	M	51	0	0	4	0
4	N	46	0	0	1	0
4	O	49	0	0	2	1
4	P	54	0	0	2	0
4	Q	59	0	0	2	0
4	R	75	0	0	2	0
4	S	51	0	0	1	1
4	T	63	0	0	0	0
4	U	51	0	0	2	0
4	V	54	0	0	3	0
4	W	77	0	0	4	0
All	All	14607	0	12554	138	2

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 5.

All (138) 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)
2:W:153:G:O3'	2:W:154:U:P	1.97	1.23
1:P:66:ARG:CZ	1:Q:66:ARG:HH21	1.62	1.12
1:R:8:ASP:OD1	4:R:2007:HOH:O	1.72	1.07
1:P:66:ARG:NH2	1:Q:66:ARG:HH21	1.53	1.04
1:P:66:ARG:HH12	1:Q:66:ARG:NH2	1.58	1.01
1:D:50:GLU:HG3	4:D:2052:HOH:O	1.61	1.01
1:P:66:ARG:HH22	1:Q:66:ARG:NH2	1.59	1.01
1:L:8:ASP:HB2	4:L:2002:HOH:O	1.60	0.99
1:P:66:ARG:NH1	1:Q:66:ARG:NH2	2.12	0.98
1:P:66:ARG:NH2	1:Q:66:ARG:NH2	2.13	0.97
1:G:24:LEU:CD2	1:G:32:PHE:HD1	1.80	0.94
1:G:24:LEU:HD21	1:G:32:PHE:CD1	2.04	0.91
1:G:24:LEU:HD21	1:G:32:PHE:CE1	2.06	0.91
1:G:24:LEU:CD2	1:G:32:PHE:CD1	2.54	0.90
1:V:8:ASP:OD1	4:V:2003:HOH:O	1.88	0.90
1:P:66:ARG:NH1	1:Q:66:ARG:HH21	1.67	0.89
2:W:134:U:H4'	4:W:2050:HOH:O	1.72	0.88
1:M:58:ARG:NH1	4:M:2037:HOH:O	2.08	0.86
1:C:73:GLU:HG2	4:C:2087:HOH:O	1.73	0.86
1:L:8:ASP:CB	4:L:2002:HOH:O	2.20	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:66:ARG:CZ	1:Q:66:ARG:NH2	2.38	0.85
1:P:6:ASN:O	1:P:6:ASN:OD1	2.00	0.80
1:V:8:ASP:OD2	4:V:2004:HOH:O	2.01	0.79
1:D:75:LYS:HG2	4:E:2040:HOH:O	1.84	0.77
1:D:18:GLY:HA2	1:D:37:LYS:HE3	1.66	0.75
1:N:18:GLY:HA2	2:W:116:G:O6	1.85	0.75
1:U:8:ASP:OD1	4:U:2004:HOH:O	2.03	0.74
1:P:66:ARG:HH12	1:Q:66:ARG:HH22	1.38	0.71
1:G:6:ASN:CB	4:G:2001:HOH:O	2.40	0.69
1:M:8:ASP:HB2	4:M:2003:HOH:O	1.92	0.69
1:E:40:LYS:NZ	4:E:2037:HOH:O	2.22	0.69
1:S:29:ASP:OD2	1:S:31:ARG:HD3	1.93	0.69
1:F:30:THR:HG1	3:G:81:TRP:N	1.92	0.68
1:C:75:LYS:H	1:C:75:LYS:CD	2.05	0.68
1:B:30:THR:HG1	3:C:81:TRP:N	1.95	0.65
1:G:30:THR:HG1	3:H:81:TRP:N	1.95	0.65
1:H:30:THR:HG1	3:I:81:TRP:N	1.94	0.64
1:D:50:GLU:OE2	4:D:2050:HOH:O	2.14	0.64
2:W:153:G:C3'	2:W:154:U:P	2.86	0.64
3:P:81:TRP:N	1:Q:30:THR:HG1	1.96	0.64
1:Q:6:ASN:ND2	1:Q:6:ASN:O	2.32	0.63
1:E:37:LYS:HE3	4:E:2034:HOH:O	1.99	0.63
1:Q:6:ASN:HA	4:Q:2005:HOH:O	1.99	0.62
1:C:72:SER:HB2	4:C:2073:HOH:O	1.99	0.62
1:C:75:LYS:H	1:C:75:LYS:HD3	1.62	0.62
1:S:59:GLY:HA2	1:S:74:GLY:HA3	1.81	0.61
1:P:18:GLY:HA2	2:W:126:G:O6	2.01	0.60
1:J:30:THR:HG1	3:K:81:TRP:N	1.99	0.60
3:L:81:TRP:N	1:M:30:THR:HG1	1.99	0.60
1:S:18:GLY:HA2	2:W:141:G:O6	2.02	0.60
3:U:81:TRP:N	1:V:30:THR:HG1	1.98	0.60
1:T:18:GLY:HA2	2:W:146:G:O6	2.02	0.60
1:M:75:LYS:HD3	1:M:75:LYS:H	1.66	0.59
1:C:75:LYS:HG3	4:D:2045:HOH:O	2.02	0.59
2:W:149:U:H4'	4:W:2070:HOH:O	2.01	0.59
1:A:30:THR:HG1	3:B:81:TRP:N	2.00	0.59
3:M:81:TRP:N	1:N:30:THR:HG1	2.01	0.58
1:U:73:GLU:HG2	4:U:2049:HOH:O	2.02	0.58
3:Q:81:TRP:N	1:R:30:THR:HG1	2.01	0.58
1:M:75:LYS:CD	1:M:75:LYS:H	2.15	0.58
1:L:22:ILE:HG12	1:L:35:SER:HB2	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:18:GLY:HA2	2:W:111:G:O6	2.02	0.57
1:N:58:ARG:NH1	4:N:2034:HOH:O	2.06	0.57
1:R:32:PHE:CZ	2:W:133:G:C6	2.93	0.57
3:N:81:TRP:N	1:O:30:THR:HG1	2.03	0.56
1:C:30:THR:HG1	3:D:81:TRP:N	2.04	0.56
1:J:66:ARG:HH22	1:K:66:ARG:HD3	1.70	0.56
1:V:18:GLY:HA2	2:W:101:G:O6	2.06	0.56
1:T:15:LEU:HB2	1:T:60:LYS:HG2	1.88	0.56
1:D:30:THR:HG1	3:E:81:TRP:N	2.05	0.55
1:A:31:ARG:HD3	4:A:2043:HOH:O	2.07	0.55
1:D:37:LYS:HG2	4:D:2042:HOH:O	2.07	0.55
1:S:8:ASP:HB2	4:S:2003:HOH:O	2.07	0.54
3:S:81:TRP:N	1:T:30:THR:HG1	2.05	0.54
1:L:30:THR:HG1	3:V:81:TRP:N	2.04	0.54
1:U:18:GLY:HA2	2:W:151:G:O6	2.06	0.54
1:L:18:GLY:HA2	2:W:106:G:O6	2.06	0.54
1:I:30:THR:HG1	3:J:81:TRP:N	2.05	0.54
1:Q:18:GLY:HA2	2:W:131:G:O6	2.08	0.54
1:O:18:GLY:HA2	2:W:121:G:O6	2.07	0.53
1:G:15:LEU:HB2	1:G:60:LYS:HG2	1.90	0.53
1:P:29:ASP:OD2	1:P:31:ARG:HD3	2.07	0.53
1:O:32:PHE:CZ	2:W:118:G:C6	2.96	0.53
3:O:81:TRP:N	1:P:30:THR:HG1	2.07	0.53
1:P:6:ASN:HA	4:Q:2007:HOH:O	2.07	0.53
3:R:81:TRP:N	1:S:30:THR:HG1	2.07	0.53
1:L:45:ILE:HD13	1:M:10:VAL:HG11	1.91	0.53
1:R:18:GLY:HA2	2:W:136:G:O6	2.08	0.53
1:R:56:LYS:NZ	1:R:58:ARG:HH11	2.07	0.53
1:E:30:THR:HG1	3:F:81:TRP:N	2.07	0.52
3:A:81:TRP:N	1:K:30:THR:HG1	2.07	0.52
1:C:75:LYS:HE2	4:C:2020:HOH:O	2.10	0.52
1:U:33:HIS:NE2	1:U:52:THR:OG1	2.41	0.51
1:L:8:ASP:OD2	4:L:2002:HOH:O	2.19	0.51
1:G:56:LYS:NZ	1:G:58:ARG:HH21	2.08	0.51
1:R:29:ASP:OD2	1:R:31:ARG:HD3	2.11	0.50
1:F:66:ARG:HD2	4:F:2048:HOH:O	2.12	0.50
3:T:81:TRP:N	1:U:30:THR:HG1	2.08	0.50
1:G:24:LEU:HD23	1:G:32:PHE:HD1	1.72	0.50
1:O:33:HIS:CE1	1:O:52:THR:OG1	2.65	0.49
1:V:5:THR:N	4:V:2001:HOH:O	2.46	0.49
1:P:32:PHE:CZ	2:W:123:G:C6	3.01	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:40:LYS:HE3	4:K:2013:HOH:O	2.12	0.48
2:W:101:G:H8	4:W:2014:HOH:O	1.95	0.48
1:G:24:LEU:HD21	1:G:32:PHE:HE1	1.68	0.48
1:L:32:PHE:CZ	2:W:103:G:C6	3.01	0.48
1:O:6:ASN:O	1:O:6:ASN:ND2	2.47	0.48
1:S:32:PHE:CZ	2:W:138:G:C6	3.01	0.47
1:O:8:ASP:OD2	4:O:2001:HOH:O	2.20	0.47
1:M:8:ASP:CB	4:M:2003:HOH:O	2.58	0.47
1:O:8:ASP:OD1	4:O:2002:HOH:O	2.20	0.47
1:R:6:ASN:HA	4:R:2005:HOH:O	2.13	0.47
1:D:37:LYS:HE2	4:D:2040:HOH:O	2.16	0.46
1:Q:6:ASN:HD22	1:Q:6:ASN:C	2.18	0.46
2:W:101:G:H2'	4:W:2014:HOH:O	2.15	0.46
1:M:8:ASP:OD2	4:M:2003:HOH:O	2.21	0.45
1:R:56:LYS:HZ3	1:R:58:ARG:HH11	1.63	0.45
1:T:33:HIS:NE2	1:T:52:THR:OG1	2.47	0.45
1:D:18:GLY:CA	1:D:37:LYS:HE3	2.40	0.45
1:U:13:LYS:HB2	1:V:70:ILE:HD11	1.99	0.45
1:B:66:ARG:HG3	1:B:66:ARG:O	2.18	0.44
1:J:66:ARG:HH22	1:K:66:ARG:CD	2.30	0.44
1:N:33:HIS:NE2	1:N:52:THR:OG1	2.44	0.44
2:W:126:G:H4'	2:W:127:A:O5'	2.17	0.44
1:D:59:GLY:CA	1:D:75:LYS:HE2	2.48	0.43
1:P:33:HIS:NE2	1:P:52:THR:OG1	2.37	0.43
1:D:40:LYS:HE3	4:D:2006:HOH:O	2.17	0.43
1:R:32:PHE:CZ	2:W:133:G:C5	3.07	0.43
1:G:6:ASN:HB3	4:G:2001:HOH:O	2.10	0.42
1:T:11:VAL:HB	1:T:64:GLN:HB2	2.01	0.42
1:U:11:VAL:HB	1:U:64:GLN:HB2	2.01	0.42
1:P:32:PHE:CZ	2:W:123:G:C5	3.08	0.42
1:O:6:ASN:HD22	1:O:6:ASN:C	2.21	0.42
1:P:66:ARG:HD3	4:P:2045:HOH:O	2.20	0.41
1:P:6:ASN:HB3	4:P:2003:HOH:O	2.19	0.41
1:D:59:GLY:N	1:D:75:LYS:HE2	2.36	0.41
1:E:71:GLU:O	1:F:13:LYS:HE3	2.21	0.40
2:W:152:A:C5	2:W:153:G:C6	3.10	0.40

All (2) 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	Interatomic distance (Å)	Clash overlap (Å)
4:C:2047:HOH:O	4:K:2061:HOH:O[4_555]	1.95	0.25
4:O:2005:HOH:O	4:S:2049:HOH:O[2_656]	1.98	0.22

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	67/74 (90%)	67 (100%)	0	0	100	100
1	B	66/74 (89%)	66 (100%)	0	0	100	100
1	C	68/74 (92%)	68 (100%)	0	0	100	100
1	D	67/74 (90%)	67 (100%)	0	0	100	100
1	E	66/74 (89%)	66 (100%)	0	0	100	100
1	F	67/74 (90%)	67 (100%)	0	0	100	100
1	G	68/74 (92%)	68 (100%)	0	0	100	100
1	H	66/74 (89%)	66 (100%)	0	0	100	100
1	I	67/74 (90%)	67 (100%)	0	0	100	100
1	J	65/74 (88%)	65 (100%)	0	0	100	100
1	K	67/74 (90%)	67 (100%)	0	0	100	100
1	L	68/74 (92%)	68 (100%)	0	0	100	100
1	M	69/74 (93%)	69 (100%)	0	0	100	100
1	N	68/74 (92%)	68 (100%)	0	0	100	100
1	O	68/74 (92%)	68 (100%)	0	0	100	100
1	P	68/74 (92%)	68 (100%)	0	0	100	100
1	Q	68/74 (92%)	68 (100%)	0	0	100	100
1	R	68/74 (92%)	68 (100%)	0	0	100	100
1	S	68/74 (92%)	68 (100%)	0	0	100	100
1	T	68/74 (92%)	68 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	U	68/74 (92%)	68 (100%)	0	0	100	100
1	V	68/74 (92%)	68 (100%)	0	0	100	100
All	All	1483/1628 (91%)	1483 (100%)	0	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	57/62 (92%)	55 (96%)	2 (4%)	36	13
1	B	56/62 (90%)	54 (96%)	2 (4%)	35	13
1	C	58/62 (94%)	52 (90%)	6 (10%)	7	1
1	D	57/62 (92%)	54 (95%)	3 (5%)	22	5
1	E	56/62 (90%)	54 (96%)	2 (4%)	35	13
1	F	57/62 (92%)	53 (93%)	4 (7%)	15	2
1	G	58/62 (94%)	57 (98%)	1 (2%)	60	42
1	H	56/62 (90%)	55 (98%)	1 (2%)	59	40
1	I	57/62 (92%)	57 (100%)	0	100	100
1	J	56/62 (90%)	53 (95%)	3 (5%)	22	5
1	K	57/62 (92%)	54 (95%)	3 (5%)	22	5
1	L	58/62 (94%)	56 (97%)	2 (3%)	37	14
1	M	59/62 (95%)	56 (95%)	3 (5%)	24	6
1	N	58/62 (94%)	56 (97%)	2 (3%)	37	14
1	O	58/62 (94%)	55 (95%)	3 (5%)	23	6
1	P	58/62 (94%)	56 (97%)	2 (3%)	37	14
1	Q	58/62 (94%)	55 (95%)	3 (5%)	23	6
1	R	58/62 (94%)	55 (95%)	3 (5%)	23	6
1	S	58/62 (94%)	57 (98%)	1 (2%)	60	42

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	T	58/62 (94%)	56 (97%)	2 (3%)	37	14
1	U	58/62 (94%)	55 (95%)	3 (5%)	23	6
1	V	58/62 (94%)	56 (97%)	2 (3%)	37	14
All	All	1264/1364 (93%)	1211 (96%)	53 (4%)	30	9

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	ASP
1	A	70	ILE
1	B	69	VAL
1	B	70	ILE
1	C	6	ASN
1	C	20	ASN
1	C	35	SER
1	C	66	ARG
1	C	70	ILE
1	C	75	LYS
1	D	8	ASP
1	D	29	ASP
1	D	70	ILE
1	E	8	ASP
1	E	70	ILE
1	F	29	ASP
1	F	31	ARG
1	F	70	ILE
1	F	73	GLU
1	G	66	ARG
1	H	70	ILE
1	J	8	ASP
1	J	29	ASP
1	J	70	ILE
1	K	20	ASN
1	K	35	SER
1	K	75	LYS
1	L	6	ASN
1	L	35	SER
1	M	6	ASN
1	M	8	ASP
1	M	75	LYS
1	N	8	ASP

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Mol	Chain	Res	Type
1	N	37	LYS
1	O	6	ASN
1	O	8	ASP
1	O	37	LYS
1	P	8	ASP
1	P	37	LYS
1	Q	6	ASN
1	Q	8	ASP
1	Q	37	LYS
1	R	6	ASN
1	R	8	ASP
1	R	37	LYS
1	S	37	LYS
1	T	35	SER
1	T	37	LYS
1	U	6	ASN
1	U	8	ASP
1	U	37	LYS
1	V	8	ASP
1	V	37	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	20	ASN
1	H	20	ASN
1	K	20	ASN
1	L	6	ASN
1	O	6	ASN
1	O	33	HIS
1	P	6	ASN
1	Q	6	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	W	34/55 (61%)	7 (20%)	1 (2%)

All (7) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	W	102	A
2	W	107	A
2	W	112	A
2	W	122	A
2	W	127	A
2	W	129	U
2	W	139	U

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	W	126	G

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](#)

22 ligands are modelled in this entry.

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

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	W	11

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	W	139:U	O3'	141:G	P	6.01
1	W	129:U	O3'	131:G	P	5.77
1	W	144:U	O3'	146:G	P	5.74
1	W	104:U	O3'	106:G	P	5.72
1	W	119:U	O3'	121:G	P	5.71
1	W	124:U	O3'	126:G	P	5.59
1	W	109:U	O3'	111:G	P	5.47
1	W	149:U	O3'	151:G	P	5.33
1	W	153:G	O3'	154:U	P	1.97
1	W	113:G	O3'	114:U	P	1.85
1	W	133:G	O3'	134:U	P	1.82

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, 95th 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 ²)	Q<0.9
1	A	69/74 (93%)	0.31	4 (5%) 23 28	8, 11, 20, 26	5 (7%)
1	B	68/74 (91%)	0.09	0 100 100	8, 11, 18, 21	7 (10%)
1	C	70/74 (94%)	0.77	9 (12%) 3 5	7, 11, 19, 26	5 (7%)
1	D	69/74 (93%)	0.46	8 (11%) 4 6	8, 11, 18, 29	8 (11%)
1	E	68/74 (91%)	0.05	1 (1%) 73 80	7, 11, 16, 22	8 (11%)
1	F	69/74 (93%)	-0.11	1 (1%) 75 82	8, 11, 16, 23	8 (11%)
1	G	70/74 (94%)	0.08	2 (2%) 51 57	9, 11, 20, 24	10 (14%)
1	H	68/74 (91%)	-0.17	0 100 100	9, 11, 17, 20	5 (7%)
1	I	69/74 (93%)	0.07	2 (2%) 51 57	7, 11, 17, 27	7 (10%)
1	J	67/74 (90%)	0.26	0 100 100	8, 11, 17, 21	4 (5%)
1	K	69/74 (93%)	0.60	8 (11%) 4 6	8, 11, 15, 26	6 (8%)
1	L	70/74 (94%)	0.19	3 (4%) 35 41	7, 10, 16, 23	3 (4%)
1	M	71/74 (95%)	0.17	2 (2%) 53 58	7, 11, 17, 24	5 (7%)
1	N	70/74 (94%)	0.06	2 (2%) 51 57	7, 10, 17, 23	1 (1%)
1	O	70/74 (94%)	0.12	1 (1%) 75 82	7, 11, 18, 20	0
1	P	70/74 (94%)	0.17	1 (1%) 75 82	7, 11, 18, 25	2 (2%)
1	Q	70/74 (94%)	0.18	2 (2%) 51 57	7, 11, 16, 23	1 (1%)
1	R	70/74 (94%)	0.32	3 (4%) 35 41	7, 11, 17, 19	5 (7%)
1	S	70/74 (94%)	0.11	1 (1%) 75 82	8, 11, 16, 21	2 (2%)
1	T	70/74 (94%)	0.13	1 (1%) 75 82	7, 11, 17, 23	2 (2%)
1	U	70/74 (94%)	-0.03	1 (1%) 75 82	7, 10, 17, 21	3 (4%)
1	V	70/74 (94%)	0.17	2 (2%) 51 57	7, 11, 17, 20	2 (2%)
2	W	44/55 (80%)	1.24	9 (20%) 1 1	11, 13, 24, 24	0
All	All	1571/1683 (93%)	0.21	63 (4%) 38 45	7, 11, 19, 29	99 (6%)

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	75	LYS	4.6
1	D	74	GLY	4.3
1	R	74	GLY	4.2
1	L	32	PHE	4.1
1	V	74	GLY	4.1
1	M	75	LYS	3.9
1	C	63	ILE	3.9
1	N	74	GLY	3.8
1	I	75	LYS	3.7
1	G	74	GLY	3.7
1	O	74	GLY	3.7
1	G	6	ASN	3.5
1	D	44	LEU	3.4
1	C	12	ILE	3.2
1	S	74	GLY	3.2
1	F	75	LYS	3.0
1	E	32	PHE	2.9
2	W	129	U	2.9
2	W	114	U	2.9
2	W	109	U	2.8
1	M	74	GLY	2.8
1	L	74	GLY	2.8
1	D	11	VAL	2.7
1	D	12	ILE	2.7
1	V	32	PHE	2.7
1	K	12	ILE	2.6
1	C	74	GLY	2.6
1	D	75	LYS	2.6
1	C	44	LEU	2.6
2	W	104	U	2.5
1	R	32	PHE	2.5
1	D	45	ILE	2.5
1	K	44	LEU	2.5
1	A	74	GLY	2.5
2	W	144	U	2.5
1	D	63	ILE	2.4
1	K	57	VAL	2.4
1	A	32	PHE	2.4
1	Q	10	VAL	2.4
1	U	32	PHE	2.4
1	C	22	ILE	2.3
1	K	46	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
2	W	119	U	2.3
1	K	55	ILE	2.2
1	N	39	ASP	2.2
2	W	139	U	2.2
1	C	10	VAL	2.2
1	K	10	VAL	2.2
1	C	55	ILE	2.2
1	C	57	VAL	2.1
1	D	43	VAL	2.1
1	P	11	VAL	2.1
1	Q	11	VAL	2.1
1	L	6	ASN	2.1
1	K	63	ILE	2.1
1	R	11	VAL	2.1
1	I	74	GLY	2.1
1	C	11	VAL	2.1
1	K	11	VAL	2.1
1	T	74	GLY	2.1
2	W	111	G	2.0
1	A	44	LEU	2.0
2	W	102	A	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 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, 95th 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(\AA^2)	Q<0.9
3	TRP	B	81	15/15	0.93	0.08	5,9,13,13	0
3	TRP	A	81	15/15	0.93	0.11	8,9,11,12	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	TRP	C	81	15/15	0.94	0.10	7,9,11,12	0
3	TRP	L	81	15/15	0.94	0.07	7,9,11,12	0
3	TRP	R	81	15/15	0.94	0.10	9,10,11,11	0
3	TRP	V	81	15/15	0.95	0.07	6,8,11,11	0
3	TRP	P	81	15/15	0.95	0.10	6,8,11,11	0
3	TRP	J	81	15/15	0.95	0.09	6,9,12,12	0
3	TRP	Q	81	15/15	0.95	0.09	7,8,10,10	0
3	TRP	E	81	15/15	0.95	0.08	6,9,12,12	0
3	TRP	F	81	15/15	0.95	0.08	7,9,11,12	0
3	TRP	M	81	15/15	0.95	0.06	6,8,11,12	0
3	TRP	S	81	15/15	0.95	0.08	8,9,12,13	0
3	TRP	K	81	15/15	0.95	0.12	7,10,12,12	0
3	TRP	D	81	15/15	0.95	0.10	7,9,12,13	0
3	TRP	G	81	15/15	0.96	0.07	7,9,11,12	0
3	TRP	H	81	15/15	0.96	0.08	7,9,10,11	0
3	TRP	U	81	15/15	0.96	0.09	7,9,10,12	0
3	TRP	I	81	15/15	0.96	0.07	7,10,12,13	0
3	TRP	O	81	15/15	0.96	0.07	7,9,10,10	0
3	TRP	N	81	15/15	0.96	0.07	7,8,11,12	0
3	TRP	T	81	15/15	0.97	0.07	7,8,10,11	0

6.5 Other polymers

There are no such residues in this entry.