



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:52 AM UTC

PDB ID : 4V9F / pdb_00004v9f
Title : The re-refined crystal structure of the Haloarcula marismortui large ribosomal subunit at 2.4 Angstrom resolution: more complete structure of the L7/L12 and L1 stalk, L5 and LX proteins
Authors : Gabdulkhakov, A.
Deposited on : 2012-11-02
Resolution : 2.40 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

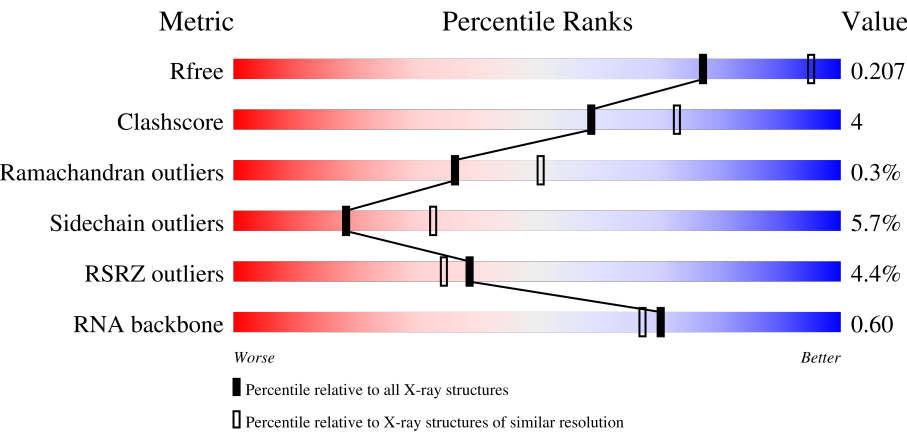
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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.



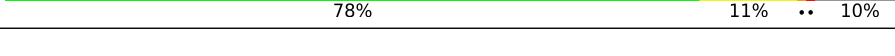
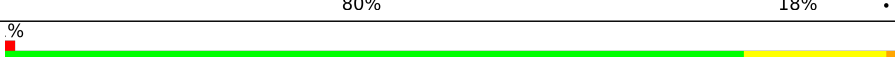
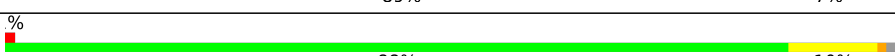
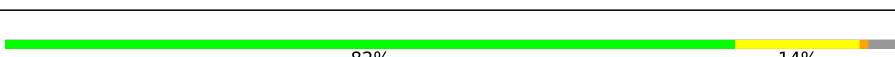
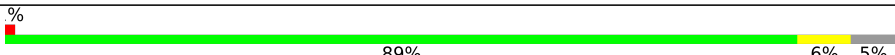
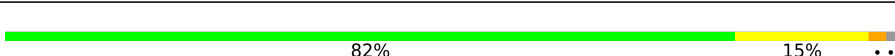
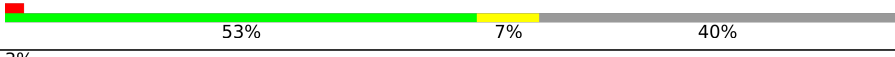
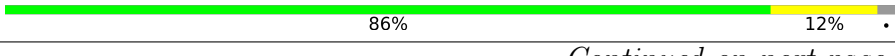
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)
RNA backbone	3983	1155 (2.70-2.10)

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 of 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	0	2910	% 76% 18% . .
2	9	122	2% 70% 25% 6%
3	A	240	% 85% 12% ..
4	B	338	85% 12% .

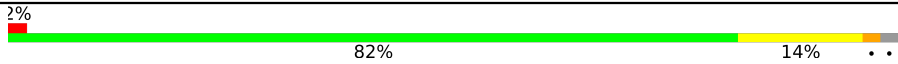
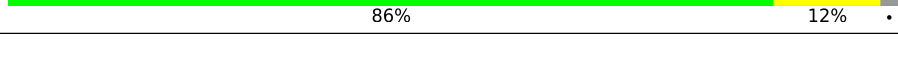
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Mol	Chain	Length	Quality of chain
5	C	246	
6	D	177	
7	E	178	
8	F	120	
9	G	255	
10	H	177	
11	I	162	
12	J	145	
13	K	132	
14	L	165	
15	M	196	
16	N	187	
17	O	116	
18	P	149	
19	Q	96	
20	R	155	
21	S	85	
22	T	120	
23	U	67	
24	V	71	
25	W	154	
26	X	92	
27	Y	241	
28	Z	92	
29	1	57	

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Mol	Chain	Length	Quality of chain
30	2	50	
31	3	92	
32	4	70	
32	5	70	
33	6	56	

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
37	CL	9	203	-	-	X	-
38	ACY	0	8213	-	-	X	-

2 Entry composition

There are 40 unique types of molecules in this entry. The entry contains 103831 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 RNA chain called 23S Ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	0	2808	Total	C	N	O	P	0	0	0
			60186	26859	11069	19451	2807			

- Molecule 2 is a RNA chain called 5S Ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	9	122	Total	C	N	O	P	0	0	0
			2599	1160	471	847	121			

- Molecule 3 is a protein called 50S ribosomal protein L2P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	237	Total	C	N	O	S	0	0	0
			1753	1072	352	324	5			

- Molecule 4 is a protein called 50S ribosomal protein L3P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	B	337	Total	C	N	O	S	0	0	0
			2625	1616	493	511	5			

- Molecule 5 is a protein called 50S ribosomal protein L4P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	C	246	Total	C	N	O	S	0	0	0
			1859	1130	345	383	1			

- Molecule 6 is a protein called 50S ribosomal protein L5P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	D	166	Total	C	N	O	S	0	0	0
			1299	812	227	256	4			

- Molecule 7 is a protein called 50S ribosomal protein L6P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	E	174	Total	C	N	O	S	0	0	0
			1376	850	230	292	4			

- Molecule 8 is a protein called 50S ribosomal protein L7Ae.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	F	119	Total	C	N	O	S	0	0	0
			889	551	141	196	1			

- Molecule 9 is a protein called 50S ribosomal protein L10E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	G	254	Total	C	N	O	S	0	0	0
			1805	1113	304	386	2			

- Molecule 10 is a protein called 50S ribosomal protein L10e.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	H	165	Total	C	N	O	S	0	0	0
			1320	822	246	246	6			

- Molecule 11 is a protein called 50S ribosomal protein L11P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	I	145	Total	C	N	O	S	0	0	0
			1088	681	170	236	1			

- Molecule 12 is a protein called 50S ribosomal protein L13P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	J	144	Total	C	N	O	S	0	0	0
			1132	704	201	224	3			

- Molecule 13 is a protein called 50S ribosomal protein L14P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	K	132	Total	C	N	O	S	0	0	0
			993	609	189	191	4			

- Molecule 14 is a protein called 50S ribosomal protein L15P.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
14	L	149	Total	C	N	O	0	0	0
			1138	682	226	230			

- Molecule 15 is a protein called 50S ribosomal protein L15e.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
15	M	195	Total	C	N	O	S	0	0	0
			1568	949	335	283	1			

- Molecule 16 is a protein called 50S ribosomal protein L18P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
16	N	186	Total	C	N	O	S	0	0	0
			1444	895	262	285	2			

- Molecule 17 is a protein called 50S ribosomal protein L18e.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
17	O	115	Total	C	N	O	0	0	0
			865	529	161	175			

- Molecule 18 is a protein called 50S ribosomal protein L19e.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
18	P	144	Total	C	N	O	0	0	0
			1144	687	231	226			

- Molecule 19 is a protein called 50S ribosomal protein L21e.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
19	Q	95	Total	C	N	O	0	0	0
			734	450	141	143			

- Molecule 20 is a protein called 50S ribosomal protein L22P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
20	R	150	Total	C	N	O	S	0	0	0
			1149	713	209	223	4			

- Molecule 21 is a protein called 50S ribosomal protein L23P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
21	S	81	Total	C	N	O	S	0	0	0
			641	389	111	138	3			

- Molecule 22 is a protein called 50S ribosomal protein L24P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
22	T	119	Total	C	N	O		0	0	0
			949	568	180	201				

- Molecule 23 is a protein called 50S ribosomal protein L24e.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
23	U	56	Total	C	N	O	S	0	0	0
			435	259	81	90	5			

- Molecule 24 is a protein called 50S ribosomal protein L29P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
24	V	65	Total	C	N	O	S	0	0	0
			499	304	94	100	1			

- Molecule 25 is a protein called 50S ribosomal protein L30P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
25	W	154	Total	C	N	O	S	0	0	0
			1196	737	209	244	6			

- Molecule 26 is a protein called 50S ribosomal protein L31e.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
26	X	90	Total	C	N	O	S	0	0	0
			714	438	137	138	1			

- Molecule 27 is a protein called 50S ribosomal protein L32e.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
27	Y	144	Total	C	N	O		0	0	0
			1146	696	230	220				

- Molecule 28 is a protein called 50S ribosomal protein L37Ae.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
28	Z	83	Total	C	N	O	S	0	0	0
			645	387	127	126	5			

- Molecule 29 is a protein called 50S ribosomal protein L37e.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
29	1	56	Total	C	N	O	S	0	0	0
			431	258	86	83	4			

- Molecule 30 is a protein called 50S ribosomal protein L39e.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
30	2	49	Total	C	N	O	S	0	0	0
			421	254	93	73	1			

- Molecule 31 is a protein called 50S ribosomal protein L44e.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
31	3	92	Total	C	N	O	S	0	0	0
			755	458	153	137	7			

- Molecule 32 is a protein called 50S ribosomal protein L12.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
32	4	58	Total	C	N	O	0	0	0
			290	174	58	58			
32	5	56	Total	C	N	O	0	0	0
			280	168	56	56			

- Molecule 33 is a protein called 50S ribosomal protein LX.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
33	6	55	Total	C	N	O	0	0	0
			441	271	78	92			

- Molecule 34 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
34	0	125	Total	Mg	0	0
			125	125		
34	9	1	Total	Mg	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
34	A	3	Total 3	Mg 3	0	0
34	K	1	Total 1	Mg 1	0	0
34	L	1	Total 1	Mg 1	0	0
34	N	1	Total 1	Mg 1	0	0
34	T	1	Total 1	Mg 1	0	0
34	Y	4	Total 4	Mg 4	0	0
34	3	1	Total 1	Mg 1	0	0

- Molecule 35 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
35	0	2	Total 2	K 2	0	0
35	O	1	Total 1	K 1	0	0

- Molecule 36 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
36	0	72	Total 72	Na 72	0	0
36	9	1	Total 1	Na 1	0	0
36	A	1	Total 1	Na 1	0	0
36	C	1	Total 1	Na 1	0	0
36	H	2	Total 2	Na 2	0	0
36	J	1	Total 1	Na 1	0	0
36	L	2	Total 2	Na 2	0	0
36	M	1	Total 1	Na 1	0	0

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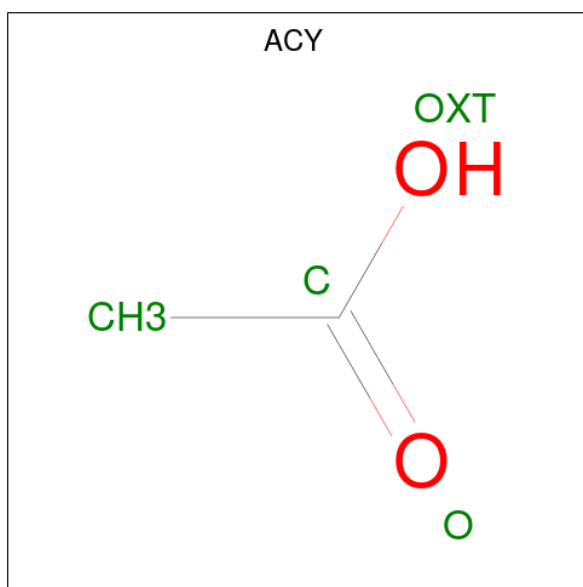
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
36	Q	1	Total 1	Na 1	0	0
36	R	2	Total 2	Na 2	0	0
36	S	1	Total 1	Na 1	0	0

- Molecule 37 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
37	0	12	Total 12	Cl 12	0	0
37	9	1	Total 1	Cl 1	0	0
37	A	1	Total 1	Cl 1	0	0
37	B	2	Total 2	Cl 2	0	0
37	F	1	Total 1	Cl 1	0	0
37	G	1	Total 1	Cl 1	0	0
37	J	3	Total 3	Cl 3	0	0
37	L	1	Total 1	Cl 1	0	0
37	M	1	Total 1	Cl 1	0	0
37	N	1	Total 1	Cl 1	0	0
37	O	2	Total 2	Cl 2	0	0
37	R	2	Total 2	Cl 2	0	0
37	Y	1	Total 1	Cl 1	0	0
37	3	1	Total 1	Cl 1	0	0

- Molecule 38 is ACETIC ACID (CCD ID: ACY) (formula: C₂H₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
38	0	1	Total	C	O	0	0
			4	2	2		
38	0	1	Total	C	O	0	0
			4	2	2		
38	0	1	Total	C	O	0	0
			4	2	2		
38	W	1	Total	C	O	0	0
			4	2	2		

- Molecule 39 is CADMIUM ION (CCD ID: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
39	O	1	Total	Cd	0	0
			1	1		
39	U	1	Total	Cd	0	0
			1	1		
39	Z	1	Total	Cd	0	0
			1	1		
39	1	1	Total	Cd	0	0
			1	1		
39	3	1	Total	Cd	0	0
			1	1		

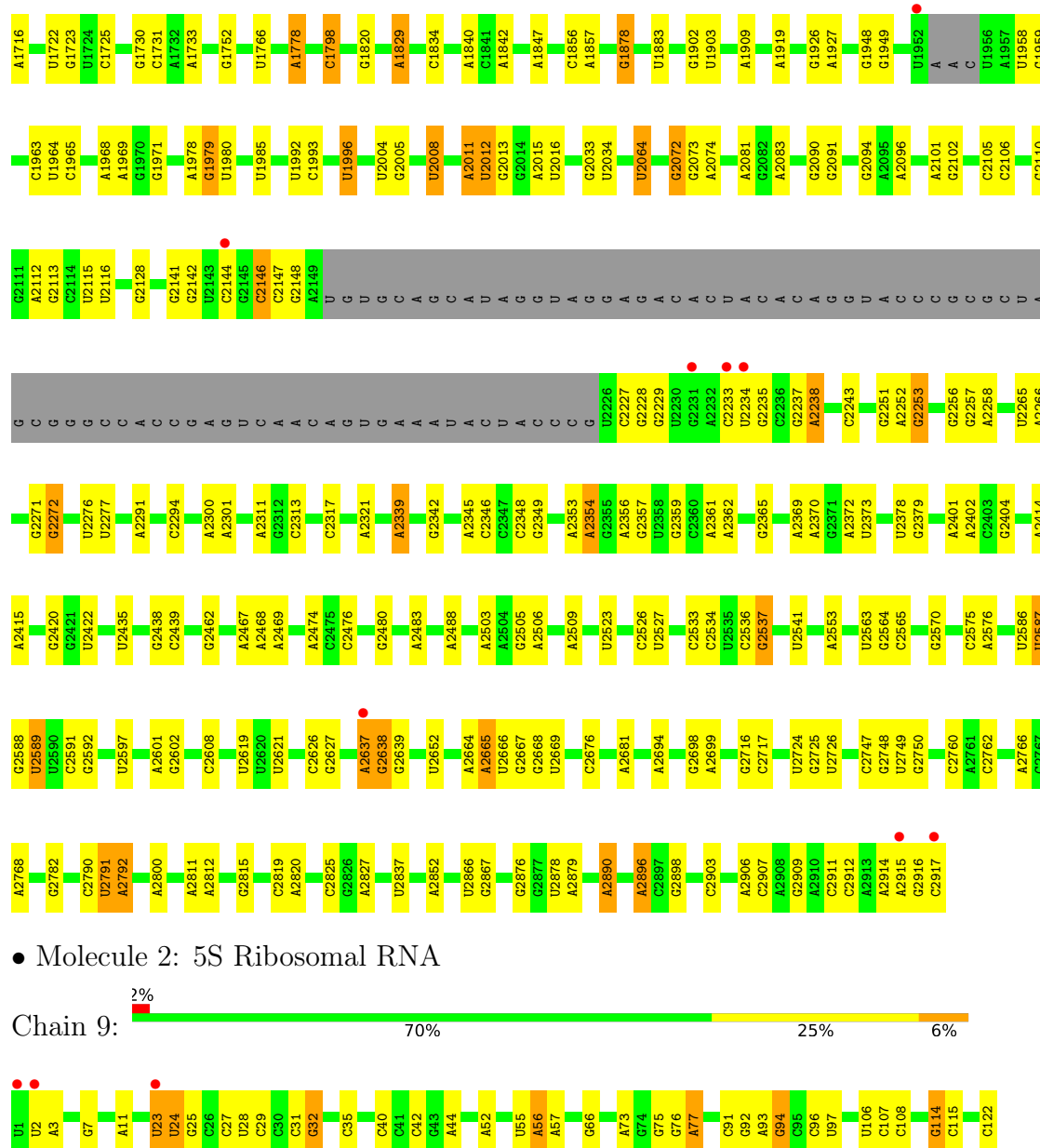
- Molecule 40 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
40	0	5924	Total 5924	O 5924	0	0
40	9	153	Total 153	O 153	0	0
40	A	126	Total 126	O 126	0	0
40	B	151	Total 151	O 151	0	0
40	C	171	Total 171	O 171	0	0
40	D	9	Total 9	O 9	0	0
40	E	44	Total 44	O 44	0	0
40	F	22	Total 22	O 22	0	0
40	G	3	Total 3	O 3	0	0
40	H	58	Total 58	O 58	0	0
40	I	1	Total 1	O 1	0	0
40	J	54	Total 54	O 54	0	0
40	K	59	Total 59	O 59	0	0
40	L	67	Total 67	O 67	0	0
40	M	133	Total 133	O 133	0	0
40	N	59	Total 59	O 59	0	0
40	O	45	Total 45	O 45	0	0
40	P	63	Total 63	O 63	0	0
40	Q	54	Total 54	O 54	0	0
40	R	72	Total 72	O 72	0	0
40	S	34	Total 34	O 34	0	0
40	T	39	Total 39	O 39	0	0

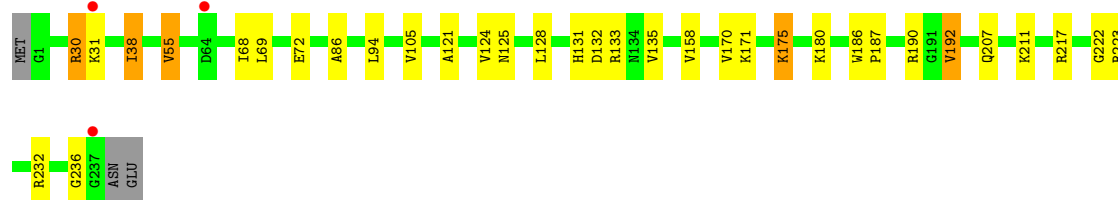
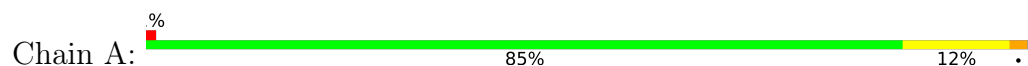
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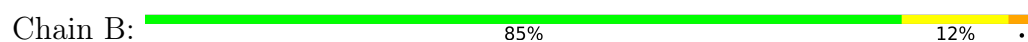
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40	U	23	Total 23	O 23	0	0
40	V	12	Total 12	O 12	0	0
40	W	64	Total 64	O 64	0	0
40	X	20	Total 20	O 20	0	0
40	Y	104	Total 104	O 104	0	0
40	Z	28	Total 28	O 28	0	0
40	1	50	Total 50	O 50	0	0
40	2	39	Total 39	O 39	0	0
40	3	63	Total 63	O 63	0	0
40	6	1	Total 1	O 1	0	0

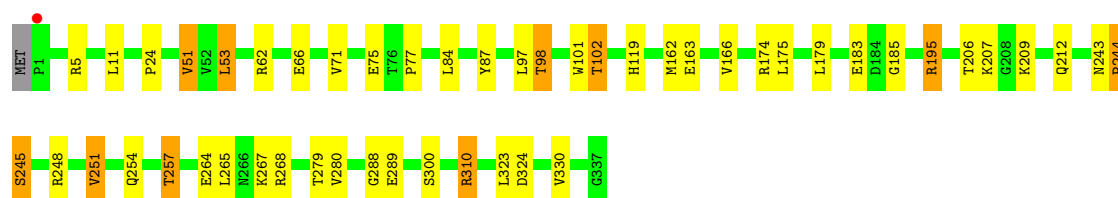


• Molecule 3: 50S ribosomal protein L2P



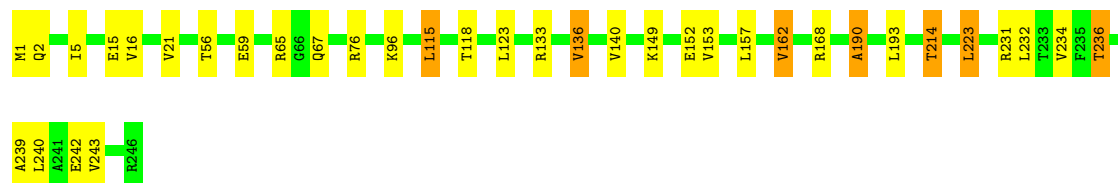
• Molecule 4: 50S ribosomal protein L3P





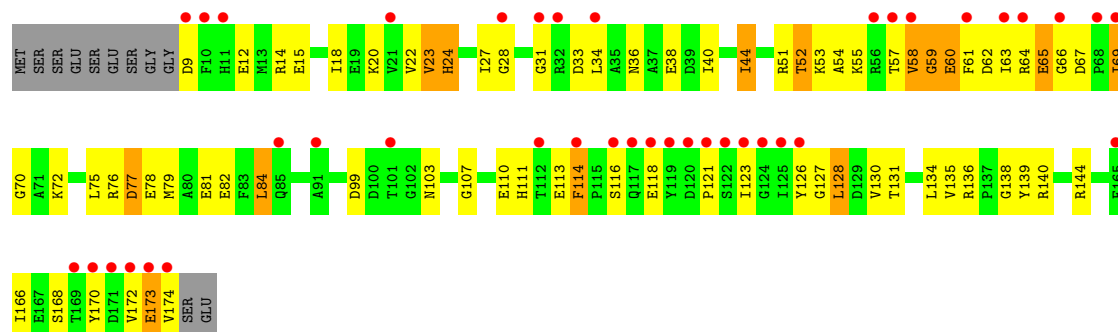
- Molecule 5: 50S ribosomal protein L4P

Chain C: 85% 12%



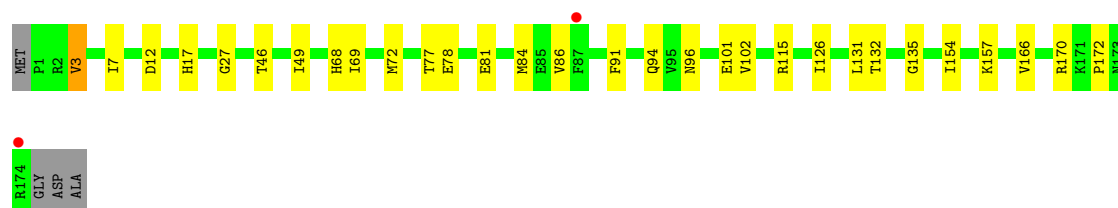
- Molecule 6: 50S ribosomal protein L5P

Chain D: 23% 52% 34% 8% 6%



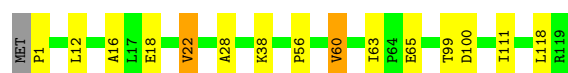
- Molecule 7: 50S ribosomal protein L6P

Chain E: % 81% 16%

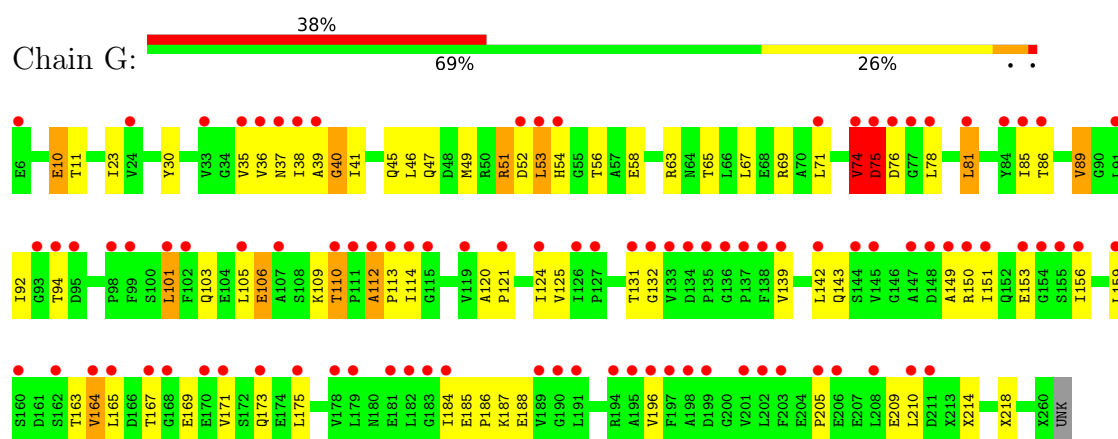


- Molecule 8: 50S ribosomal protein L7Ae

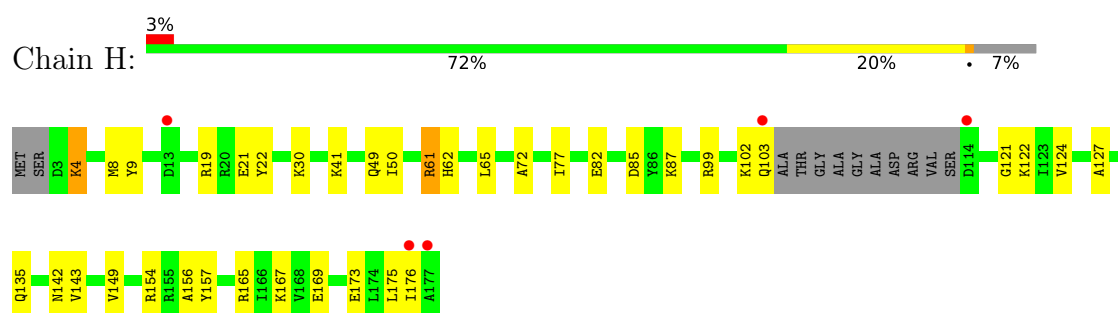
Chain F: 87% 11%



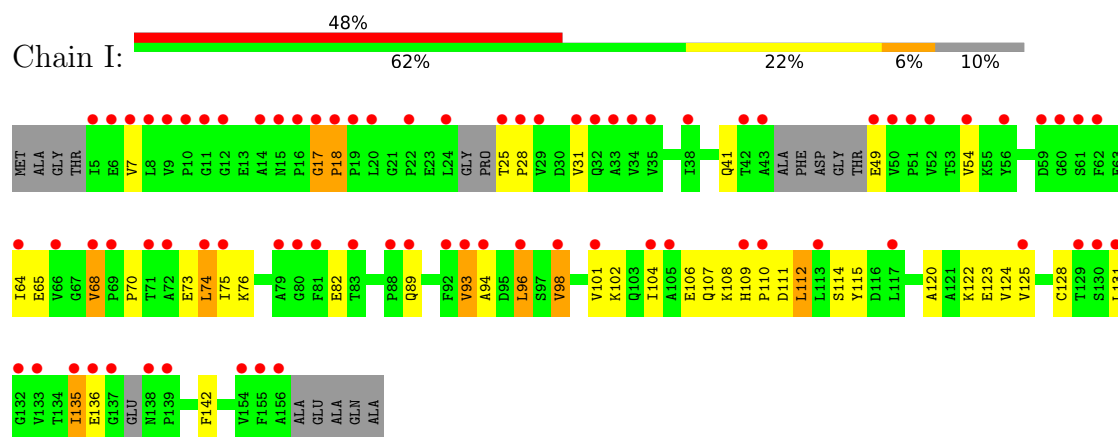
- Molecule 9: 50S ribosomal protein L10E



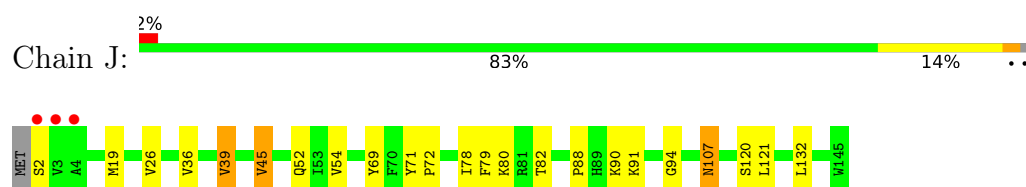
• Molecule 10: 50S ribosomal protein L10e



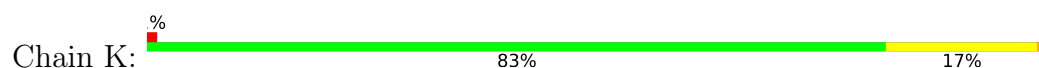
• Molecule 11: 50S ribosomal protein L11P



• Molecule 12: 50S ribosomal protein L13P

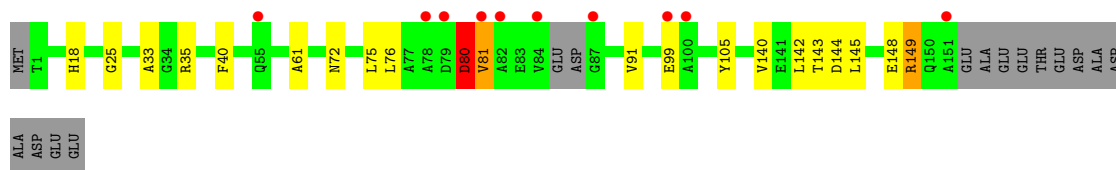
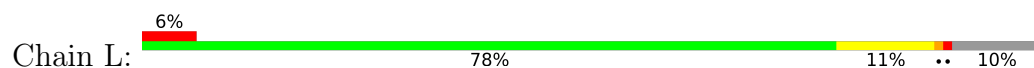


• Molecule 13: 50S ribosomal protein L14P





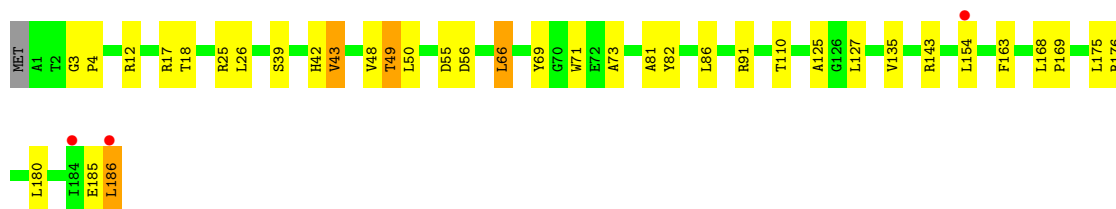
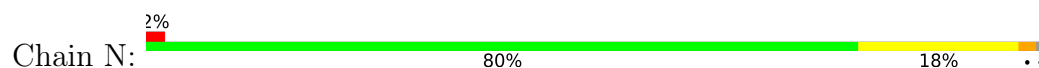
- Molecule 14: 50S ribosomal protein L15P



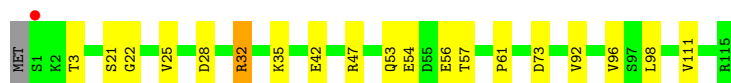
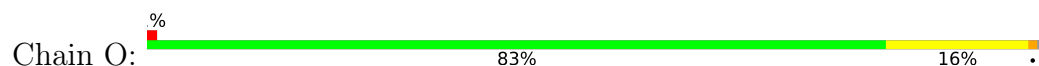
- Molecule 15: 50S ribosomal protein L15e



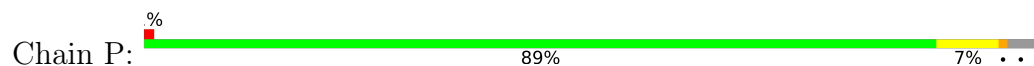
- Molecule 16: 50S ribosomal protein L18P



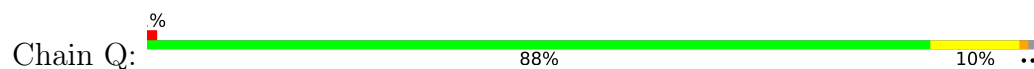
- Molecule 17: 50S ribosomal protein L18e

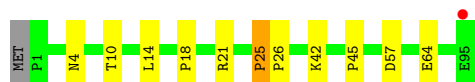


- Molecule 18: 50S ribosomal protein L19e



- Molecule 19: 50S ribosomal protein L21e





- Molecule 20: 50S ribosomal protein L22P

Chain R: 82% 14% ..



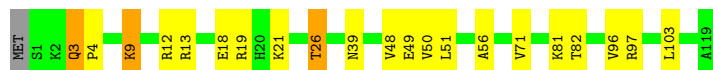
- Molecule 21: 50S ribosomal protein L23P

Chain S: 89% 6% 5%



- Molecule 22: 50S ribosomal protein L24P

Chain T: 82% 15% ..



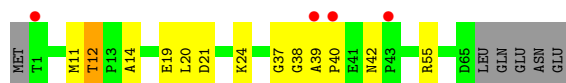
- Molecule 23: 50S ribosomal protein L24e

Chain U: 69% 13% 16%



- Molecule 24: 50S ribosomal protein L29P

Chain V: 73% 17% 8%



- Molecule 25: 50S ribosomal protein L30P

Chain W: 81% 19%

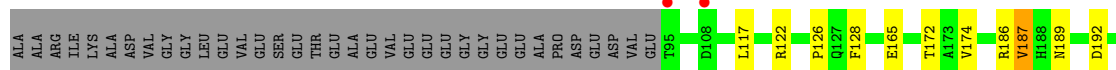
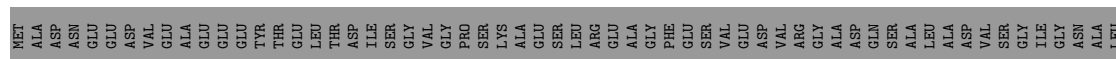


- Molecule 26: 50S ribosomal protein L31e

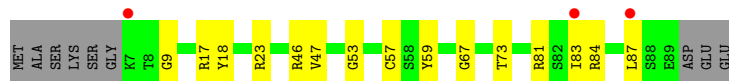
Chain X: 70% 26% ..



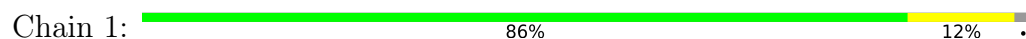
- Molecule 27: 50S ribosomal protein L32e



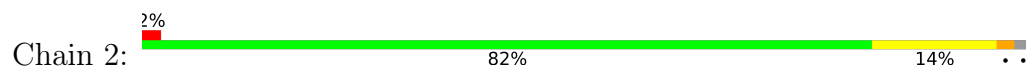
- Molecule 28: 50S ribosomal protein L37Ae



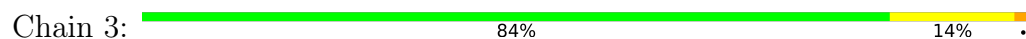
- Molecule 29: 50S ribosomal protein L37e



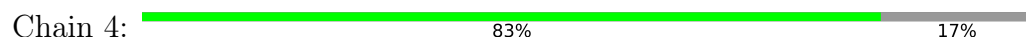
- Molecule 30: 50S ribosomal protein L39e



- Molecule 31: 50S ribosomal protein L44e

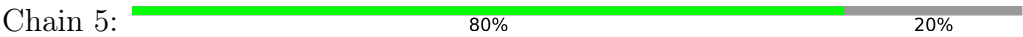


- Molecule 32: 50S ribosomal protein L12

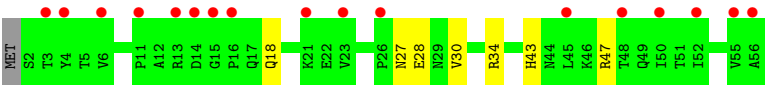
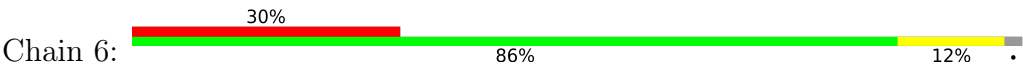




- Molecule 32: 50S ribosomal protein L12



- Molecule 33: 50S ribosomal protein LX



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	211.66Å 299.67Å 573.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	85.48 – 2.40 85.48 – 2.40	Depositor EDS
% Data completeness (in resolution range)	95.5 (85.48-2.40) 95.5 (85.48-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.40Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168), CNS	Depositor
R, R_{free}	0.166 , 0.206 0.167 , 0.207	Depositor DCC
R_{free} test set	6982 reflections (1.00%)	wwPDB-VP
Wilson B-factor (Å ²)	42.3	Xtriage
Anisotropy	0.271	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 44.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	103831	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.51% 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 ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: NA, MG, 1MA, OMG, PSU, CL, ACY, CD, OMU, UR3, K

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	0	0.30	1/67259 (0.0%)	0.50	3/104899 (0.0%)
2	9	0.26	0/2904	0.44	0/4526
3	A	0.50	0/1786	0.87	1/2408 (0.0%)
4	B	0.47	0/2690	0.87	2/3652 (0.1%)
5	C	0.55	0/1884	0.94	4/2552 (0.2%)
6	D	0.41	0/1324	0.94	5/1791 (0.3%)
7	E	0.44	0/1401	0.89	6/1905 (0.3%)
8	F	0.50	0/900	0.85	0/1224
9	G	0.43	0/1585	0.92	3/2156 (0.1%)
10	H	0.49	0/1340	0.87	2/1795 (0.1%)
11	I	0.44	0/1107	1.01	6/1512 (0.4%)
12	J	0.45	0/1148	0.93	2/1548 (0.1%)
13	K	0.53	0/1003	0.89	0/1351
14	L	0.53	0/1150	0.92	3/1536 (0.2%)
15	M	0.53	0/1592	0.87	3/2127 (0.1%)
16	N	0.46	0/1473	0.89	4/1999 (0.2%)
17	O	0.50	0/874	0.83	0/1181
18	P	0.45	0/1155	0.77	0/1539
19	Q	0.51	0/748	0.86	2/1005 (0.2%)
20	R	0.51	0/1172	0.88	1/1578 (0.1%)
21	S	0.45	0/648	0.80	0/875
22	T	0.49	0/957	0.88	2/1289 (0.2%)
23	U	0.43	0/443	0.75	0/597
24	V	0.44	0/502	0.79	0/675
25	W	0.50	0/1219	0.92	0/1655
26	X	0.47	0/725	0.96	3/978 (0.3%)
27	Y	0.51	0/1162	0.84	0/1558
28	Z	0.49	0/656	0.91	3/877 (0.3%)
29	1	0.62	0/438	1.03	2/578 (0.3%)
30	2	0.55	0/427	0.74	0/566
31	3	0.49	0/771	0.79	0/1024
33	6	0.45	0/449	0.83	1/607 (0.2%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
All	All	0.37	1/102892 (0.0%)	0.63	58/153563 (0.0%)

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
6	D	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	0	128	A	O3'-P	5.68	1.69	1.61

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	I	17	GLY	CA-C-N	9.63	130.30	120.38
11	I	17	GLY	C-N-CA	9.63	130.30	120.38
7	E	135	GLY	CA-C-N	8.96	128.61	118.85
7	E	135	GLY	C-N-CA	8.96	128.61	118.85
11	I	93	VAL	N-CA-C	8.63	119.42	110.36
6	D	114	PHE	CA-C-N	7.96	127.68	119.56
6	D	114	PHE	C-N-CA	7.96	127.68	119.56
9	G	40	GLY	N-CA-C	-7.85	97.38	110.80
26	X	80	GLU	CB-CA-C	-7.48	107.94	116.54
11	I	18	PRO	N-CA-C	7.19	119.47	110.70
26	X	65	ASN	N-CA-C	6.27	119.87	112.72
15	M	120	VAL	N-CA-C	6.22	117.44	111.91
1	0	128	A	OP2-P-O3'	6.18	126.53	108.00
6	D	15	GLU	CA-C-N	6.16	126.18	119.90
6	D	15	GLU	C-N-CA	6.16	126.18	119.90
19	Q	25	PRO	CA-C-N	6.10	125.72	119.56
19	Q	25	PRO	C-N-CA	6.10	125.72	119.56
9	G	112	ALA	CA-C-N	6.04	127.39	119.84
9	G	112	ALA	C-N-CA	6.04	127.39	119.84
14	L	81	VAL	N-CA-C	6.04	121.89	109.34
14	L	80	ASP	N-CA-C	6.03	118.66	107.99
14	L	72	ASN	N-CA-C	5.98	120.70	113.17
12	J	107	ASN	CA-C-N	5.93	125.55	119.56
12	J	107	ASN	C-N-CA	5.93	125.55	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	H	9	TYR	N-CA-C	5.87	121.36	113.72
11	I	68	VAL	CA-C-N	5.77	123.87	119.66
11	I	68	VAL	C-N-CA	5.77	123.87	119.66
1	0	518	G	O4'-C1'-N9	5.76	116.85	108.20
10	H	124	VAL	N-CA-C	5.63	119.57	113.43
7	E	7	ILE	CA-C-N	-5.60	114.98	120.52
7	E	7	ILE	C-N-CA	-5.60	114.98	120.52
16	N	3	GLY	CA-C-N	5.57	125.41	119.28
16	N	3	GLY	C-N-CA	5.57	125.41	119.28
3	A	222	GLY	N-CA-C	-5.56	107.02	114.25
28	Z	46	ARG	N-CA-C	5.53	119.65	113.02
6	D	77	ASP	CB-CA-C	-5.49	110.27	116.63
28	Z	47	VAL	N-CA-C	5.47	116.63	108.54
15	M	42	ARG	CA-C-N	5.46	125.38	119.76
15	M	42	ARG	C-N-CA	5.46	125.38	119.76
5	C	56	THR	CA-C-N	-5.42	113.07	119.84
5	C	56	THR	C-N-CA	-5.42	113.07	119.84
16	N	163	PHE	N-CA-C	5.36	118.04	111.82
7	E	17	HIS	CB-CA-C	-5.30	110.02	117.23
33	6	43	HIS	N-CA-C	5.21	120.28	113.30
1	0	128	A	O3'-P-O5'	-5.18	96.23	104.00
5	C	190	ALA	CA-C-N	-5.17	115.44	121.64
5	C	190	ALA	C-N-CA	-5.17	115.44	121.64
22	T	3	GLN	CA-C-N	5.17	125.21	119.32
22	T	3	GLN	C-N-CA	5.17	125.21	119.32
29	1	24	GLU	CA-C-N	-5.16	114.41	122.05
29	1	24	GLU	C-N-CA	-5.16	114.41	122.05
16	N	42	HIS	N-CA-C	5.13	116.47	109.18
28	Z	17	ARG	N-CA-C	5.12	117.47	110.55
26	X	24	LYS	N-CA-C	5.11	119.60	113.17
20	R	141	VAL	N-CA-C	5.10	115.19	107.75
7	E	27	GLY	N-CA-C	5.07	116.87	110.38
4	B	244	PRO	CA-C-N	-5.02	115.34	122.77
4	B	244	PRO	C-N-CA	-5.02	115.34	122.77

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
6	D	59	GLY	Peptide
6	D	65	GLU	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	0	60186	0	30379	291	0
2	9	2599	0	1325	24	0
3	A	1753	0	1766	18	0
4	B	2625	0	2533	29	0
5	C	1859	0	1813	23	0
6	D	1299	0	1263	51	0
7	E	1376	0	1285	17	0
8	F	889	0	843	6	0
9	G	1805	0	1560	48	0
10	H	1320	0	1331	23	0
11	I	1088	0	1029	27	0
12	J	1132	0	1112	15	0
13	K	993	0	1027	12	0
14	L	1138	0	1096	10	0
15	M	1568	0	1586	20	0
16	N	1444	0	1401	15	0
17	O	865	0	873	10	0
18	P	1144	0	1129	9	0
19	Q	734	0	729	9	0
20	R	1149	0	1122	15	0
21	S	641	0	605	3	0
22	T	949	0	924	14	0
23	U	435	0	394	6	0
24	V	499	0	511	7	0
25	W	1196	0	1137	21	0
26	X	714	0	700	15	0
27	Y	1146	0	1148	10	0
28	Z	645	0	610	7	0
29	1	431	0	426	5	0
30	2	421	0	437	8	0
31	3	755	0	728	7	0
32	4	290	0	60	0	0
32	5	280	0	59	0	0
33	6	441	0	411	4	0
34	0	125	0	0	0	0
34	3	1	0	0	0	0
34	9	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
34	A	3	0	0	0	0
34	K	1	0	0	0	0
34	L	1	0	0	0	0
34	N	1	0	0	0	0
34	T	1	0	0	0	0
34	Y	4	0	0	0	0
35	0	2	0	0	0	0
35	O	1	0	0	0	0
36	0	72	0	0	0	0
36	9	1	0	0	0	0
36	A	1	0	0	0	0
36	C	1	0	0	0	0
36	H	2	0	0	0	0
36	J	1	0	0	0	0
36	L	2	0	0	0	0
36	M	1	0	0	0	0
36	Q	1	0	0	0	0
36	R	2	0	0	0	0
36	S	1	0	0	0	0
37	0	12	0	0	1	0
37	3	1	0	0	0	0
37	9	1	0	0	3	0
37	A	1	0	0	0	0
37	B	2	0	0	0	0
37	F	1	0	0	1	0
37	G	1	0	0	0	0
37	J	3	0	0	1	0
37	L	1	0	0	0	0
37	M	1	0	0	0	0
37	N	1	0	0	0	0
37	O	2	0	0	1	0
37	R	2	0	0	0	0
37	Y	1	0	0	0	0
38	0	12	0	9	4	0
38	W	4	0	3	1	0
39	1	1	0	0	0	0
39	3	1	0	0	0	0
39	O	1	0	0	0	0
39	U	1	0	0	0	0
39	Z	1	0	0	0	0
40	0	5924	0	0	16	0
40	1	50	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
40	2	39	0	0	1	0
40	3	63	0	0	0	0
40	6	1	0	0	1	0
40	9	153	0	0	3	0
40	A	126	0	0	2	0
40	B	151	0	0	3	0
40	C	171	0	0	6	0
40	D	9	0	0	3	0
40	E	44	0	0	3	0
40	F	22	0	0	0	0
40	G	3	0	0	0	0
40	H	58	0	0	2	0
40	I	1	0	0	0	0
40	J	54	0	0	0	0
40	K	59	0	0	1	0
40	L	67	0	0	1	0
40	M	133	0	0	8	0
40	N	59	0	0	0	0
40	O	45	0	0	1	0
40	P	63	0	0	1	0
40	Q	54	0	0	2	0
40	R	72	0	0	1	0
40	S	34	0	0	0	0
40	T	39	0	0	0	0
40	U	23	0	0	0	0
40	V	12	0	0	0	0
40	W	64	0	0	1	0
40	X	20	0	0	2	0
40	Y	104	0	0	2	0
40	Z	28	0	0	1	0
All	All	103831	0	63364	698	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 4.

All (698) 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)
15:M:164:THR:HG22	15:M:167:GLY:H	1.28	0.97
5:C:236:THR:HG22	5:C:239:ALA:H	1.30	0.95
1:O:156:C:H5''	15:M:171:ARG:HD3	1.46	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:R:39:THR:HG22	20:R:42:GLU:H	1.33	0.94
1:0:1160:G:H5'	1:0:1161:A:H5'	1.58	0.85
12:J:19:MET:HE3	12:J:132:LEU:HD11	1.58	0.85
1:0:2586:U:H3	1:0:2592:G:H22	1.22	0.85
2:9:75:G:H1	2:9:106:U:H3	1.25	0.84
20:R:99:ALA:HB1	20:R:109:MET:HE1	1.61	0.82
1:0:2665:A:H62	1:0:2917:C:H41	1.29	0.81
12:J:19:MET:HE1	12:J:78:ILE:HG22	1.62	0.81
2:9:76:G:H3'	2:9:77:A:H5''	1.63	0.81
17:O:32:ARG:HH21	17:O:35:LYS:HG3	1.45	0.81
37:0:8210:CL:CL	40:B:637:HOH:O	2.36	0.80
23:U:52:THR:HG22	23:U:55:ALA:H	1.46	0.80
15:M:102:GLU:OE1	15:M:164:THR:HG21	1.81	0.80
1:0:1164:U:H3	1:0:1192:A:H2	1.29	0.79
25:W:5:VAL:HG11	25:W:153:MET:HE1	1.64	0.79
1:0:1242:A:H5'	12:J:82:THR:HG23	1.66	0.77
4:B:212:GLN:HB2	4:B:257:THR:HG21	1.69	0.74
7:E:101:GLU:OE1	7:E:115:ARG:NH1	2.20	0.73
4:B:195:ARG:NH1	4:B:324:ASP:OD1	2.21	0.73
1:0:929:A:N3	40:0:8652:HOH:O	2.21	0.73
1:0:2141:G:N2	1:0:2234:U:O2	2.21	0.72
25:W:149:LEU:HG	25:W:153:MET:HE2	1.70	0.72
1:0:1441:G:OP1	30:2:39:ARG:NH1	2.22	0.72
1:0:1165:G:N2	1:0:1173:A:OP1	2.22	0.72
9:G:110:THR:HG22	9:G:112:ALA:HB2	1.71	0.72
9:G:75:ASP:HA	9:G:78:LEU:H	1.55	0.71
1:0:2665:A:N6	1:0:2917:C:H41	1.88	0.71
1:0:967:U:H5''	38:0:8213:ACY:H3	1.72	0.71
37:9:203:CL:CL	40:9:447:HOH:O	2.45	0.71
10:H:102:LYS:HD2	10:H:122:LYS:HD3	1.71	0.70
11:I:108:LYS:HB2	11:I:111:ASP:HB2	1.73	0.70
26:X:8:ARG:NH1	40:X:104:HOH:O	2.25	0.69
1:0:1160:G:H8	1:0:1207:A:H62	1.36	0.69
1:0:2349:G:OP1	6:D:20:LYS:NZ	2.25	0.69
4:B:102:THR:HG21	4:B:185:GLY:HA2	1.74	0.69
16:N:49:THR:HG22	16:N:56:ASP:HB2	1.75	0.69
9:G:51:ARG:NH2	11:I:112:LEU:O	2.24	0.69
4:B:62:ARG:NH2	4:B:66:GLU:O	2.26	0.68
1:0:31:C:H4'	22:T:9:LYS:HD2	1.75	0.68
1:0:1351:G:OP1	5:C:96:LYS:NZ	2.24	0.68
9:G:150:ARG:HE	9:G:159:LEU:HD22	1.59	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:C:1:MET:HG2	5:C:2:GLN:H	1.60	0.66
2:9:42:C:O2	6:D:76:ARG:NH1	2.28	0.66
1:0:966:U:C2	38:0:8213:ACY:H2	2.32	0.65
26:X:18:ARG:NH1	40:X:109:HOH:O	2.26	0.64
4:B:310:ARG:NH2	40:B:548:HOH:O	2.30	0.64
7:E:77:THR:OG1	7:E:78:GLU:N	2.31	0.64
1:0:2345:A:H2	6:D:28:GLY:HA3	1.63	0.63
5:C:133:ARG:NH1	40:C:511:HOH:O	2.21	0.63
9:G:125:VAL:HG13	9:G:163:THR:HA	1.79	0.63
26:X:41:PHE:O	26:X:76:ARG:NH1	2.32	0.63
5:C:236:THR:HG21	40:C:482:HOH:O	1.99	0.63
5:C:162:VAL:HG22	5:C:232:LEU:HD21	1.81	0.63
1:0:73:U:OP1	24:V:55:ARG:NH2	2.31	0.63
19:Q:4:ASN:ND2	40:Q:250:HOH:O	2.32	0.62
20:R:122:GLN:HB3	20:R:138:SER:HB2	1.81	0.62
11:I:93:VAL:N	11:I:94:ALA:HA	2.13	0.62
2:9:73:A:H61	2:9:108:C:H42	1.47	0.62
11:I:98:VAL:HG13	11:I:135:ILE:HD11	1.81	0.62
1:0:1120:U:H5''	1:0:1120:U:H6	1.65	0.62
3:A:192:VAL:HG13	3:A:207:GLN:HB3	1.81	0.62
15:M:107:ARG:NH1	40:M:373:HOH:O	2.30	0.62
1:0:2488:A:H61	1:0:2534:C:H42	1.46	0.61
7:E:126:ILE:HG22	7:E:131:LEU:HG	1.81	0.61
22:T:12:ARG:O	22:T:19:ARG:NH2	2.33	0.61
1:0:2142:G:N2	1:0:2233:C:O2	2.33	0.61
6:D:134:LEU:HD11	6:D:166:ILE:HD11	1.83	0.61
15:M:5:TYR:HE1	15:M:46:LEU:HD13	1.65	0.61
2:9:40:C:N3	37:9:203:CL:CL	2.71	0.61
13:K:32:ILE:HD11	13:K:56:SER:HB2	1.83	0.61
1:0:2474:A:N3	40:0:9378:HOH:O	2.31	0.61
3:A:232:ARG:NH2	3:A:236:GLY:O	2.26	0.61
4:B:195:ARG:HG2	4:B:323:LEU:HD22	1.83	0.60
1:0:1666:C:H2'	1:0:1667:A:H8	1.66	0.60
1:0:1834:C:H2'	1:0:1840:A:N6	2.17	0.60
1:0:2353:A:H1'	19:Q:21:ARG:HH22	1.67	0.60
26:X:2:ALA:N	26:X:4:ASP:OD1	2.34	0.60
33:6:18:GLN:NE2	40:6:101:HOH:O	2.32	0.60
1:0:1964:U:H2'	1:0:1965:C:C6	2.36	0.60
9:G:184:ILE:HG22	9:G:186:PRO:HD3	1.83	0.60
30:2:38:LYS:NZ	40:2:101:HOH:O	2.33	0.60
5:C:118:THR:O	5:C:136:VAL:HG13	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:D:135:VAL:HG21	6:D:139:TYR:CD1	2.37	0.60
16:N:82:TYR:OH	16:N:176:ARG:NH1	2.35	0.59
2:9:92:G:H2'	2:9:93:A:C8	2.37	0.59
1:0:1197:G:H21	1:0:1202:A:H2	1.49	0.59
5:C:5:ILE:HD11	5:C:16:VAL:HG13	1.85	0.59
1:0:327:A:OP1	5:C:149:LYS:NZ	2.20	0.59
6:D:173:GLU:C	6:D:174:VAL:HG23	2.27	0.59
7:E:170:ARG:NH1	40:E:209:HOH:O	2.35	0.59
10:H:173:GLU:CD	10:H:173:GLU:H	2.10	0.59
1:0:2251:G:H2'	1:0:2252:A:C8	2.38	0.59
1:0:1676:G:OP1	30:2:8:LYS:NZ	2.36	0.59
1:0:2414:A:H2'	1:0:2415:A:C8	2.37	0.59
6:D:52:THR:HG22	6:D:70:GLY:H	1.66	0.59
1:0:1025:C:P	25:W:108:ARG:HH12	2.26	0.58
1:0:1197:G:O2'	1:0:1202:A:N6	2.35	0.58
15:M:164:THR:HG22	15:M:167:GLY:N	2.10	0.58
26:X:3:SER:HB2	26:X:79:GLU:HG3	1.85	0.58
1:0:1500:U:OP2	18:P:41:ARG:NH2	2.37	0.58
11:I:54:VAL:HG22	11:I:64:ILE:HG13	1.85	0.58
1:0:1151:G:OP1	9:G:63:ARG:NH1	2.36	0.57
1:0:820:G:C6	3:A:171:LYS:HB2	2.39	0.57
1:0:2866:U:H4'	1:0:2867:G:H5'	1.86	0.57
1:0:969:G:H1	1:0:999:C:N4	2.03	0.57
12:J:19:MET:HE2	12:J:79:PHE:HA	1.87	0.57
27:Y:126:PRO:HG2	27:Y:128:PHE:CZ	2.40	0.57
1:0:112:G:OP1	29:1:20:ARG:NH1	2.38	0.56
1:0:1008:C:H5''	10:H:19:ARG:HH12	1.69	0.56
1:0:2637:A:H2	1:0:2639:G:OP2	1.88	0.56
11:I:114:SER:OG	11:I:123:GLU:OE1	2.22	0.56
26:X:30:MET:HE1	26:X:54:ILE:HG22	1.87	0.56
6:D:38:GLU:OE2	6:D:51:ARG:NH1	2.38	0.56
1:0:969:G:H1	1:0:999:C:H42	1.52	0.56
4:B:162:MET:SD	4:B:310:ARG:HD3	2.45	0.56
13:K:34:VAL:HG22	13:K:47:ALA:HB2	1.88	0.56
1:0:1500:U:P	18:P:41:ARG:HH22	2.29	0.55
9:G:75:ASP:HA	9:G:78:LEU:N	2.20	0.55
11:I:75:ILE:HD11	11:I:131:LEU:HD13	1.88	0.55
9:G:171:VAL:HG23	9:G:175:LEU:HD23	1.88	0.55
3:A:30:ARG:HD3	3:A:38:ILE:HG21	1.88	0.55
9:G:71:LEU:HD13	9:G:81:LEU:HB3	1.88	0.55
9:G:214:UNK:O	9:G:218:UNK:N	2.39	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:0:2346:C:O2'	6:D:52:THR:HG21	2.06	0.55
1:0:92:G:H5'	24:V:42:ASN:HD22	1.72	0.55
1:0:1733:A:H4'	4:B:212:GLN:HA	1.88	0.55
26:X:30:MET:HE3	26:X:58:ALA:HB3	1.89	0.55
1:0:588:G:O6	25:W:154:ARG:NH1	2.39	0.55
3:A:132:ASP:HB3	3:A:135:VAL:H	1.72	0.55
6:D:99:ASP:N	6:D:103:ASN:O	2.29	0.55
7:E:84:MET:HE2	7:E:166:VAL:HG12	1.89	0.55
7:E:157:LYS:NZ	40:E:210:HOH:O	2.39	0.55
3:A:121:ALA:O	3:A:124:VAL:HG22	2.06	0.55
4:B:87:TYR:OH	4:B:163:GLU:OE2	2.24	0.55
9:G:139:VAL:HG21	9:G:151:ILE:HD11	1.89	0.55
1:0:553:G:OP2	27:Y:204:ARG:NH2	2.30	0.55
9:G:150:ARG:HH21	9:G:159:LEU:HD13	1.72	0.55
7:E:3:VAL:HG22	7:E:49:ILE:HB	1.89	0.54
1:0:1766:U:O2	1:0:1778:A:H5'	2.07	0.54
40:0:9667:HOH:O	10:H:61:ARG:HG2	2.06	0.54
1:0:1494:A:O2'	1:0:1505:U:O2	2.25	0.54
1:0:1926:G:H2'	1:0:1927:A:C8	2.42	0.54
2:9:28:U:H2'	2:9:29:C:C6	2.41	0.54
20:R:109:MET:HG2	20:R:148:GLU:C	2.31	0.54
26:X:23:HIS:CE1	26:X:24:LYS:HG3	2.43	0.54
7:E:84:MET:HB2	7:E:131:LEU:HB2	1.89	0.54
25:W:81:ASP:OD1	25:W:92:ASP:HB2	2.08	0.54
9:G:65:THR:O	9:G:69:ARG:HG2	2.07	0.54
15:M:169:ARG:NH2	40:M:351:HOH:O	2.40	0.54
1:0:371:U:H2'	1:0:372:A:C8	2.43	0.54
10:H:85:ASP:OD2	10:H:142:ASN:ND2	2.37	0.54
1:0:2072:G:C6	1:0:2533:C:H1'	2.43	0.53
9:G:53:LEU:O	9:G:54:HIS:ND1	2.40	0.53
1:0:416:G:H5''	1:0:417:G:OP2	2.07	0.53
7:E:49:ILE:HD11	7:E:69:ILE:HD12	1.90	0.53
12:J:45:VAL:HG11	12:J:121:LEU:HD22	1.89	0.53
9:G:85:ILE:HG23	9:G:89:VAL:HG11	1.90	0.53
16:N:69:TYR:OH	16:N:185:GLU:O	2.24	0.53
16:N:91:ARG:HG2	16:N:186:LEU:HB3	1.91	0.53
31:3:65:THR:HG21	31:3:83:TRP:CZ3	2.42	0.53
1:0:1120:U:H5''	1:0:1120:U:C6	2.44	0.53
10:H:61:ARG:HA	10:H:127:ALA:HA	1.91	0.53
23:U:33:SER:O	23:U:37:GLU:HG3	2.09	0.53
1:0:309:C:OP1	22:T:97:ARG:NH2	2.35	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:0:2666:U:H2'	1:0:2667:G:C8	2.44	0.53
1:0:1548:U:O2'	1:0:1798:C:O2	2.25	0.53
1:0:2637:A:H1'	1:0:2638:G:C5'	2.39	0.53
8:F:38:LYS:HE3	15:M:3:SER:HA	1.91	0.53
1:0:688:A:H61	1:0:698:A:H5''	1.73	0.53
3:A:211:LYS:HE3	40:A:484:HOH:O	2.09	0.52
16:N:71:TRP:CE2	16:N:73:ALA:HB3	2.44	0.52
17:O:54:GLU:HG3	37:O:203:CL:CL	2.47	0.52
30:2:26:MET:HG2	30:2:33:VAL:O	2.08	0.52
1:0:86:A:H4'	1:0:87:C:H5''	1.91	0.52
9:G:101:LEU:HD12	9:G:196:VAL:HG21	1.90	0.52
1:0:2911:C:H2'	1:0:2912:C:C6	2.45	0.52
6:D:57:THR:HB	6:D:64:ARG:HG3	1.91	0.52
13:K:7:ASP:OD2	13:K:81:ARG:NH2	2.41	0.52
17:O:32:ARG:HG2	17:O:61:PRO:HB2	1.91	0.52
24:V:12:THR:HG23	24:V:14:ALA:H	1.75	0.52
1:0:1181:A:N1	1:0:1192:A:O2'	2.41	0.52
1:0:1657:A:H2'	1:0:1658:A:C8	2.44	0.52
1:0:2878:U:H2'	1:0:2879:A:O4'	2.09	0.52
1:0:119:A:H2'	1:0:120:A:H5''	1.92	0.52
20:R:18:LEU:HG	20:R:91:LEU:HD13	1.90	0.52
22:T:9:LYS:HG2	22:T:13:ARG:CZ	2.39	0.52
9:G:45:GLN:O	9:G:49:MET:HG3	2.10	0.52
9:G:142:LEU:HD12	9:G:149:ALA:HB2	1.90	0.52
9:G:164:VAL:HG22	9:G:165:LEU:HG	1.90	0.52
1:0:2064:U:H5'	1:0:2652:U:H4'	1.92	0.52
1:0:2716:G:H5''	4:B:206:THR:HG21	1.91	0.52
1:0:2815:G:H21	12:J:2:SER:HA	1.75	0.52
18:P:80:ARG:HD2	18:P:87:ARG:NE	2.25	0.52
25:W:2:HIS:HE1	38:W:201:ACY:H1	1.74	0.52
31:3:5:ARG:NE	31:3:92:GLU:OE1	2.28	0.52
1:0:966:U:H2'	38:0:8213:ACY:H2	1.92	0.51
1:0:2637:A:H1'	1:0:2638:G:H5''	1.92	0.51
9:G:151:ILE:HA	9:G:156:ILE:HA	1.91	0.51
1:0:2916:G:H2'	1:0:2917:C:H5''	1.93	0.51
19:Q:42:LYS:NZ	19:Q:57:ASP:OD2	2.42	0.51
26:X:52:PRO:O	26:X:56:GLU:HG3	2.10	0.51
6:D:31:GLY:C	6:D:33:ASP:H	2.17	0.51
16:N:43:VAL:HG11	16:N:81:ALA:HA	1.92	0.51
1:0:952:G:OP1	19:Q:42:LYS:HE2	2.11	0.51
7:E:69:ILE:HA	7:E:72:MET:HE2	1.90	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:0:564:G:N3	40:0:9147:HOH:O	2.34	0.51
1:0:1020:A:H1'	40:Q:203:HOH:O	2.10	0.51
22:T:49:GLU:OE2	22:T:97:ARG:NH1	2.35	0.51
1:0:1169:U:H3	1:0:1173:A:H2'	1.75	0.51
10:H:41:LYS:O	10:H:87:LYS:NZ	2.33	0.51
11:I:125:VAL:HG21	11:I:142:PHE:CD2	2.45	0.51
25:W:64:THR:O	25:W:68:THR:HG22	2.10	0.51
9:G:30:TYR:HB3	9:G:94:THR:N	2.26	0.51
20:R:39:THR:HG23	20:R:107:GLU:O	2.11	0.51
1:0:1218:U:H2'	1:0:1219:U:C6	2.45	0.51
1:0:2256:G:H2'	1:0:2257:G:H5'	1.93	0.51
1:0:397:A:O2'	1:0:417:G:N3	2.34	0.51
6:D:59:GLY:N	6:D:60:GLU:O	2.44	0.51
9:G:56:THR:O	9:G:94:THR:OG1	2.25	0.51
1:0:877:G:H5'	1:0:878:G:OP1	2.11	0.50
1:0:1705:C:P	18:P:59:ARG:HH12	2.34	0.50
9:G:54:HIS:O	9:G:56:THR:N	2.44	0.50
9:G:23:ILE:HD13	9:G:67:LEU:HD23	1.94	0.50
14:L:76:LEU:O	14:L:80:ASP:N	2.36	0.50
16:N:12:ARG:HD3	16:N:18:THR:OG1	2.11	0.50
1:0:1383:U:OP1	40:0:9192:HOH:O	2.20	0.50
1:0:1558:C:O2	1:0:1563:G:N2	2.39	0.50
1:0:2339:A:H5''	6:D:107:GLY:HA3	1.92	0.50
1:0:2438:G:H2'	1:0:2439:C:O4'	2.12	0.50
1:0:1185:U:H2'	1:0:1186:C:C6	2.46	0.50
3:A:186:TRP:CG	3:A:187:PRO:HA	2.47	0.50
6:D:140:ARG:O	6:D:144:ARG:HG2	2.12	0.50
25:W:68:THR:HG23	25:W:69:ARG:HG2	1.92	0.50
4:B:166:VAL:O	4:B:174:ARG:HD3	2.12	0.50
1:0:2265:U:H2'	1:0:2266:A:C8	2.47	0.50
5:C:214:THR:HG23	40:C:532:HOH:O	2.11	0.50
28:Z:53:GLY:HA2	28:Z:67:GLY:O	2.11	0.50
1:0:1184:C:H2'	1:0:1185:U:H6	1.76	0.50
11:I:125:VAL:HA	11:I:128:CYS:HB2	1.94	0.50
1:0:2372:A:H2'	1:0:2373:U:C6	2.47	0.49
21:S:76:GLU:OE1	21:S:80:ARG:NH1	2.41	0.49
1:0:2300:A:H4'	1:0:2301:A:O5'	2.12	0.49
6:D:65:GLU:O	6:D:67:ASP:N	2.45	0.49
1:0:68:U:OP1	21:S:1:SER:OG	2.29	0.49
18:P:97:ARG:HD2	40:P:218:HOH:O	2.12	0.49
1:0:656:G:H5'	17:O:3:THR:OG1	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:0:1666:C:H2'	1:0:1667:A:C8	2.47	0.49
1:0:1992:U:OP2	13:K:66:ARG:HD3	2.13	0.49
1:0:1025:C:OP1	25:W:108:ARG:NH1	2.43	0.49
1:0:1164:U:N3	1:0:1192:A:H2	2.03	0.49
1:0:2359:G:N7	40:0:8507:HOH:O	2.35	0.49
6:D:20:LYS:HG2	6:D:22:VAL:HG23	1.95	0.49
10:H:135:GLN:NE2	10:H:169:GLU:OE2	2.45	0.49
1:0:2760:C:OP1	4:B:209:LYS:HE2	2.11	0.49
2:9:114:G:H2'	2:9:115:C:C6	2.48	0.49
4:B:51:VAL:HG22	4:B:330:VAL:HG22	1.94	0.49
8:F:60:VAL:HG22	8:F:63:ILE:HD12	1.94	0.49
1:0:1406:A:N6	1:0:1701:A:H5''	2.28	0.49
5:C:168:ARG:NH2	5:C:190:ALA:O	2.46	0.49
26:X:30:MET:HE2	26:X:33:ILE:HD12	1.95	0.49
1:0:485:A:N3	1:0:487:G:H5''	2.28	0.49
1:0:2591:C:H2'	1:0:2592:G:O4'	2.13	0.48
1:0:166:A:N7	14:L:25:GLY:HA2	2.28	0.48
1:0:2083:A:N6	12:J:90:LYS:HE2	2.27	0.48
1:0:2361:A:H2'	1:0:2362:A:C8	2.48	0.48
15:M:64:ARG:HD2	40:M:379:HOH:O	2.12	0.48
16:N:48:VAL:CG1	16:N:55:ASP:HB3	2.43	0.48
27:Y:117:LEU:HA	27:Y:174:VAL:HG21	1.95	0.48
31:3:14:CYS:HB3	31:3:16:GLU:HG2	1.95	0.48
4:B:24:PRO:HG2	4:B:310:ARG:HG3	1.95	0.48
1:0:1209:C:H2'	1:0:1210:G:H8	1.77	0.48
1:0:1528:A:H2'	1:0:1529:G:O4'	2.13	0.48
8:F:16:ALA:HA	8:F:111:ILE:HD13	1.95	0.48
1:0:1613:C:H2'	1:0:1614:G:O4'	2.13	0.48
9:G:37:ASN:C	9:G:39:ALA:H	2.22	0.48
9:G:103:GLN:C	9:G:105:LEU:H	2.20	0.48
6:D:9:ASP:N	6:D:9:ASP:OD1	2.47	0.48
23:U:6:CYS:HB2	23:U:32:CYS:HB3	1.96	0.48
27:Y:189:ASN:HA	27:Y:217:ILE:HD11	1.95	0.48
9:G:40:GLY:O	9:G:41:ILE:HG13	2.13	0.48
1:0:1704:G:O3'	18:P:59:ARG:NH1	2.47	0.48
1:0:2353:A:H4'	1:0:2354:A:O5'	2.13	0.48
6:D:9:ASP:O	6:D:12:GLU:HG2	2.14	0.48
12:J:88:PRO:O	12:J:94:GLY:HA3	2.14	0.48
1:0:371:U:H2'	1:0:372:A:H8	1.78	0.48
1:0:1118:A:H3'	1:0:1118:A:N3	2.29	0.48
3:A:125:ASN:HB3	3:A:158:VAL:HG12	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:G:47:GLN:O	9:G:51:ARG:N	2.47	0.48
12:J:26:VAL:HG13	12:J:36:VAL:HG11	1.94	0.48
1:O:1524:U:O2'	1:O:1525:G:OP2	2.24	0.47
9:G:164:VAL:HG13	9:G:165:LEU:H	1.79	0.47
1:O:834:G:H4'	1:O:835:U:OP2	2.14	0.47
17:O:32:ARG:NH2	17:O:35:LYS:HG3	2.23	0.47
1:O:816:G:C6	1:O:817:G:N1	2.82	0.47
1:O:2146:C:N3	1:O:2147:C:N4	2.62	0.47
2:9:107:C:H2'	2:9:108:C:C6	2.49	0.47
1:O:1304:U:H2'	1:O:1305:C:C6	2.50	0.47
10:H:72:ALA:HB2	10:H:156:ALA:HB2	1.96	0.47
22:T:51:LEU:HD11	22:T:97:ARG:HB2	1.96	0.47
6:D:111:HIS:CE1	6:D:126:TYR:H	2.33	0.47
14:L:149:ARG:HH11	14:L:149:ARG:HB3	1.80	0.47
15:M:169:ARG:NH1	40:M:369:HOH:O	2.48	0.47
27:Y:165:GLU:HG2	40:Y:477:HOH:O	2.14	0.47
27:Y:187:VAL:HG22	27:Y:192:ASP:HB3	1.97	0.47
1:O:1056:U:H2'	1:O:1057:A:O4'	2.15	0.47
1:O:1409:G:H5'	40:O:8530:HOH:O	2.13	0.47
1:O:1588:G:C6	1:O:1589:G:N1	2.83	0.47
2:9:56:A:OP2	40:9:393:HOH:O	2.21	0.47
5:C:153:VAL:O	5:C:157:LEU:HG	2.15	0.47
5:C:242:GLU:HB2	40:C:489:HOH:O	2.14	0.47
10:H:165:ARG:HH21	10:H:167:LYS:HD2	1.80	0.47
11:I:49:GLU:OE1	11:I:49:GLU:N	2.48	0.47
15:M:164:THR:HG23	40:M:426:HOH:O	2.15	0.47
1:O:1333:U:H2'	1:O:1334:C:C6	2.50	0.47
1:O:1684:A:H1'	30:2:43:ARG:HH22	1.80	0.47
6:D:27:ILE:HB	6:D:69:ILE:O	2.15	0.47
6:D:40:ILE:O	6:D:44:ILE:HG22	2.14	0.47
6:D:110:GLU:O	6:D:113:GLU:HG2	2.14	0.47
7:E:68:HIS:O	7:E:72:MET:HG3	2.14	0.47
10:H:4:LYS:HD3	10:H:8:MET:HE1	1.97	0.47
24:V:11:MET:HE1	24:V:19:GLU:HG2	1.97	0.47
9:G:58:GLU:O	9:G:92:ILE:HD12	2.15	0.47
1:O:2509:A:N7	40:O:9275:HOH:O	2.36	0.46
4:B:243:ASN:HA	4:B:244:PRO:C	2.40	0.46
6:D:58:VAL:HB	6:D:60:GLU:O	2.14	0.46
1:O:1097:A:H5''	25:W:125:HIS:CE1	2.51	0.46
6:D:99:ASP:HB3	6:D:103:ASN:H	1.80	0.46
1:O:2112:A:H2'	1:O:2113:G:C8	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:L:80:ASP:N	14:L:80:ASP:OD1	2.49	0.46
1:0:1842:A:C4	1:0:1979:G:C6	3.04	0.46
37:9:203:CL:CL	6:D:53:LYS:NZ	2.85	0.46
9:G:52:ASP:O	9:G:54:HIS:N	2.48	0.46
15:M:64:ARG:NH2	40:M:356:HOH:O	2.38	0.46
5:C:193:LEU:HB2	5:C:231:ARG:HD3	1.96	0.46
9:G:10:GLU:HG2	9:G:11:THR:H	1.80	0.46
1:0:2906:A:H5'	1:0:2907:C:O4'	2.15	0.46
40:0:8766:HOH:O	22:T:82:THR:HA	2.15	0.46
4:B:51:VAL:CG1	4:B:53:LEU:HD13	2.46	0.46
1:0:2011:A:H4'	1:0:2012:U:O5'	2.15	0.46
1:0:2237:G:O2'	1:0:2238:A:H8	1.99	0.46
1:0:2791:U:H1'	1:0:2792:A:H5''	1.98	0.46
11:I:17:GLY:HA3	11:I:18:PRO:HD3	1.73	0.46
31:3:87:ARG:NH1	31:3:89:GLU:OE2	2.49	0.46
1:0:1883:U:OP2	3:A:190:ARG:HD2	2.16	0.46
1:0:2503:A:OP1	10:H:154:ARG:NH2	2.35	0.46
4:B:5:ARG:HD2	13:K:41:LYS:HD3	1.98	0.46
6:D:135:VAL:CG2	40:D:209:HOH:O	2.63	0.46
14:L:61:ALA:HB2	14:L:105:TYR:CZ	2.51	0.46
22:T:26:THR:HA	22:T:39:ASN:HB3	1.98	0.46
1:0:1948:G:H2'	1:0:1949:G:O4'	2.15	0.46
1:0:2115:U:H2'	1:0:2116:U:C6	2.51	0.46
1:0:2348:C:H1'	6:D:131:THR:HG21	1.98	0.46
1:0:2666:U:O2'	1:0:2667:G:O4'	2.23	0.46
6:D:168:SER:C	6:D:170:TYR:H	2.24	0.46
1:0:1370:G:O5'	20:R:62:HIS:HB3	2.16	0.46
1:0:1574:C:H2'	1:0:1575:C:C6	2.51	0.46
1:0:2234:U:H2'	1:0:2235:G:H8	1.81	0.46
3:A:128:LEU:HD21	3:A:131:HIS:HE1	1.80	0.46
9:G:52:ASP:C	9:G:54:HIS:H	2.24	0.46
11:I:76:LYS:HD2	11:I:82:GLU:HA	1.98	0.46
26:X:34:ARG:NH1	26:X:48:VAL:O	2.43	0.46
1:0:1104:C:H4'	12:J:88:PRO:HD3	1.98	0.45
5:C:115:LEU:HD13	5:C:223:LEU:HD21	1.96	0.45
13:K:98:VAL:HG13	13:K:102:GLU:HA	1.97	0.45
1:0:226:A:H1'	1:0:393:G:C5	2.51	0.45
1:0:790:A:H2'	1:0:791:A:O4'	2.17	0.45
1:0:2081:A:H4'	12:J:69:TYR:CE2	2.52	0.45
23:U:31:PHE:CG	23:U:37:GLU:HG2	2.51	0.45
6:D:84:LEU:HD22	6:D:84:LEU:HA	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:D:135:VAL:HG23	40:D:208:HOH:O	2.15	0.45
7:E:96:ASN:HA	7:E:126:ILE:HD13	1.98	0.45
6:D:75:LEU:HD22	6:D:79:MET:HB3	1.98	0.45
1:0:932:U:H2'	1:0:933:C:C6	2.52	0.45
1:0:1007:A:H2'	10:H:22:TYR:CZ	2.51	0.45
7:E:69:ILE:HA	7:E:72:MET:CE	2.47	0.45
1:0:138:U:H4'	1:0:139:C:OP2	2.16	0.45
1:0:1175:G:H1'	1:0:1193:A:H2'	1.98	0.45
1:0:1180:U:H2'	1:0:1181:A:O4'	2.16	0.45
11:I:109:HIS:HB3	11:I:110:PRO:HD3	1.98	0.45
16:N:86:LEU:HD12	16:N:125:ALA:HB2	1.98	0.45
1:0:538:C:H5''	1:0:539:G:C8	2.52	0.45
2:9:24:U:H2'	2:9:25:G:O4'	2.16	0.45
1:0:1535:G:H2'	1:0:1536:C:C6	2.52	0.45
1:0:2664:A:H8	1:0:2664:A:OP1	2.00	0.45
14:L:144:ASP:O	14:L:148:GLU:HG3	2.17	0.45
15:M:5:TYR:CE1	15:M:46:LEU:HD13	2.49	0.45
17:O:57:THR:HB	17:O:111:VAL:HG12	1.97	0.45
28:Z:81:ARG:HG2	28:Z:84:ARG:HH21	1.82	0.45
1:0:749:C:OP2	40:0:8806:HOH:O	2.21	0.45
1:0:2866:U:C5	23:U:50:GLU:HB3	2.51	0.45
4:B:248:ARG:O	4:B:251:VAL:HG13	2.16	0.45
29:1:26:SER:HB2	29:1:34:CYS:SG	2.57	0.45
1:0:2637:A:H4'	1:0:2638:G:OP1	2.15	0.45
10:H:165:ARG:HD3	40:H:342:HOH:O	2.17	0.45
1:0:1010:C:H4'	16:N:4:PRO:HB2	1.99	0.44
1:0:1829:A:N6	28:Z:18:TYR:HA	2.32	0.44
9:G:35:VAL:HG22	9:G:89:VAL:HG22	1.99	0.44
9:G:120:ALA:HA	9:G:121:PRO:HD3	1.88	0.44
14:L:35:ARG:O	14:L:40:PHE:HA	2.18	0.44
1:0:88:G:H2'	1:0:89:G:C8	2.52	0.44
1:0:513:A:N3	40:0:8466:HOH:O	2.36	0.44
4:B:101:TRP:HB2	4:B:119:HIS:CD2	2.52	0.44
6:D:77:ASP:HB3	6:D:78:GLU:H	1.58	0.44
1:0:858:U:H2'	1:0:859:C:C6	2.52	0.44
10:H:143:VAL:HG13	10:H:175:LEU:HD23	1.99	0.44
1:0:2563:U:H2'	1:0:2565:C:O5'	2.17	0.44
1:0:2820:A:OP1	4:B:98:THR:HG23	2.17	0.44
3:A:55:VAL:HG22	3:A:68:ILE:O	2.17	0.44
6:D:168:SER:O	6:D:170:TYR:N	2.50	0.44
10:H:99:ARG:NH2	40:H:316:HOH:O	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:Q:18:PRO:O	19:Q:21:ARG:HG2	2.17	0.44
33:6:30:VAL:O	33:6:34:ARG:HG2	2.17	0.44
1:0:128:A:O2'	1:0:129:A:O4'	2.28	0.44
1:0:291:C:H2'	1:0:292:G:O4'	2.17	0.44
5:C:152:GLU:HG3	40:C:515:HOH:O	2.17	0.44
11:I:75:ILE:HG22	11:I:104:ILE:HD13	1.99	0.44
25:W:4:LEU:O	25:W:32:CYS:HA	2.17	0.44
1:0:486:A:OP2	22:T:81:LYS:NZ	2.46	0.44
1:0:2015:A:H2'	1:0:2016:U:O4'	2.18	0.44
2:9:73:A:N6	2:9:108:C:H42	2.14	0.44
5:C:65:ARG:HB3	5:C:67:GLN:HE21	1.83	0.44
6:D:116:SER:C	6:D:118:GLU:H	2.26	0.44
6:D:121:PRO:HB2	6:D:123:ILE:HG13	1.99	0.44
9:G:74:VAL:O	9:G:75:ASP:HB3	2.17	0.44
15:M:164:THR:HB	40:M:325:HOH:O	2.17	0.44
16:N:143:ARG:NH1	16:N:169:PRO:HB3	2.32	0.44
31:3:65:THR:HG22	31:3:82:GLY:HA2	1.99	0.44
1:0:1415:G:H5'	29:1:12:ASN:O	2.18	0.44
7:E:81:GLU:O	7:E:172:PRO:HG3	2.18	0.44
13:K:64:MET:O	13:K:67:GLN:HB2	2.17	0.44
20:R:125:ARG:NH2	40:R:369:HOH:O	2.42	0.44
22:T:18:GLU:O	22:T:21:LYS:HG2	2.17	0.44
22:T:103:LEU:HD23	22:T:103:LEU:HA	1.76	0.44
24:V:20:LEU:HG	24:V:24:LYS:HE3	1.99	0.44
9:G:150:ARG:NE	9:G:159:LEU:HD22	2.30	0.44
11:I:122:LYS:HA	11:I:125:VAL:HG22	2.00	0.44
1:0:89:G:H2'	1:0:90:A:O4'	2.18	0.44
1:0:278:A:H2'	1:0:279:C:O4'	2.18	0.44
1:0:1339:G:C6	1:0:1340:G:N1	2.85	0.44
4:B:212:GLN:HB2	4:B:257:THR:CG2	2.44	0.44
5:C:123:LEU:HD23	5:C:123:LEU:HA	1.84	0.44
9:G:106:GLU:HA	9:G:109:LYS:HD3	2.00	0.44
11:I:120:ALA:O	11:I:124:VAL:HG23	2.18	0.44
1:0:165:A:H5''	14:L:33:ALA:HB2	2.00	0.43
9:G:139:VAL:O	9:G:143:GLN:HG3	2.18	0.43
25:W:73:LEU:HD12	25:W:73:LEU:HA	1.81	0.43
28:Z:23:ARG:NH1	40:Z:203:HOH:O	2.45	0.43
1:0:484:A:N1	1:0:506:G:H4'	2.33	0.43
1:0:1181:A:C2	1:0:1192:A:C8	3.06	0.43
1:0:1594:C:O2'	1:0:1607:A:H4'	2.18	0.43
2:9:31:C:H2'	2:9:32:G:O4'	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:R:39:THR:HB	20:R:42:GLU:CD	2.43	0.43
30:2:39:ARG:HA	30:2:39:ARG:HD2	1.67	0.43
1:0:23:G:C6	1:0:24:G:N1	2.86	0.43
1:0:1095:U:O2	25:W:120:PRO:HG2	2.18	0.43
1:0:1221:G:N7	40:0:8631:HOH:O	2.37	0.43
1:0:1546:G:H1'	1:0:1702:U:O2	2.18	0.43
1:0:1847:A:OP1	3:A:175:LYS:HG3	2.19	0.43
6:D:116:SER:O	6:D:118:GLU:N	2.51	0.43
1:0:646:G:H2'	1:0:647:U:C6	2.54	0.43
1:0:776:A:H1'	1:0:779:U:O4	2.18	0.43
1:0:941:G:C5	1:0:942:U:C4	3.07	0.43
1:0:1006:A:N1	1:0:2311:A:H1'	2.32	0.43
1:0:1252:A:H4'	33:6:47:ARG:HD2	2.00	0.43
5:C:65:ARG:HB3	5:C:67:GLN:NE2	2.33	0.43
8:F:56:PRO:HA	37:F:201:CL:CL	2.55	0.43
15:M:35:GLY:O	40:M:341:HOH:O	2.21	0.43
16:N:66:LEU:HB3	16:N:71:TRP:CB	2.49	0.43
19:Q:10:THR:HB	19:Q:14:LEU:HG	2.00	0.43
1:0:1968:A:H2'	1:0:1969:A:C8	2.54	0.43
6:D:36:ASN:O	6:D:40:ILE:HG12	2.19	0.43
17:O:73:ASP:HA	17:O:92:VAL:O	2.19	0.43
1:0:1198:U:O4	40:0:9554:HOH:O	2.21	0.43
1:0:2094:G:H4'	4:B:245:SER:HB3	2.01	0.43
6:D:31:GLY:O	6:D:33:ASP:N	2.52	0.43
12:J:90:LYS:HB2	37:J:203:CL:CL	2.55	0.43
25:W:65:VAL:HA	25:W:68:THR:HG22	2.00	0.43
1:0:1130:U:O4	1:0:2523:U:H5''	2.19	0.43
6:D:54:ALA:HB2	6:D:69:ILE:HG13	2.01	0.43
25:W:150:LEU:HD23	25:W:150:LEU:HA	1.86	0.43
31:3:25:VAL:HG22	31:3:68:LYS:HG3	2.00	0.43
33:6:27:ASN:OD1	33:6:28:GLU:N	2.49	0.43
1:0:1527:A:H1'	1:0:1528:A:C8	2.54	0.43
1:0:2782:G:O6	1:0:2790:C:H5''	2.19	0.43
6:D:24:HIS:HB2	6:D:72:LYS:CB	2.49	0.43
7:E:94:GLN:HG3	40:E:221:HOH:O	2.17	0.43
11:I:73:GLU:HA	11:I:76:LYS:HG2	2.00	0.43
20:R:68:HIS:CD2	20:R:76:ASP:HB2	2.53	0.43
1:0:653:U:H2'	1:0:654:A:C8	2.53	0.43
1:0:1191:A:H3'	1:0:1192:A:H5''	2.01	0.43
1:0:2724:U:H2'	1:0:2725:G:O4'	2.19	0.43
1:0:2890:A:H1'	23:U:56:ARG:NH2	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:C:2:GLN:HG2	5:C:15:GLU:OE2	2.18	0.43
11:I:108:LYS:HB3	11:I:108:LYS:HE2	1.91	0.43
12:J:71:TYR:CG	12:J:72:PRO:HD2	2.54	0.43
1:0:64:G:H2'	1:0:65:C:O4'	2.18	0.42
1:0:1131:G:C6	1:0:1230:A:C4	3.07	0.42
1:0:1716:A:H4'	18:P:55:LYS:HD3	2.01	0.42
1:0:1909:A:N1	1:0:2128:G:H1'	2.34	0.42
1:0:2637:A:C2	1:0:2639:G:OP2	2.70	0.42
1:0:2911:C:H2'	1:0:2912:C:H6	1.81	0.42
3:A:72:GLU:OE2	28:Z:73:THR:OG1	2.35	0.42
4:B:75:GLU:C	4:B:77:PRO:HD3	2.44	0.42
11:I:68:VAL:HG11	11:I:108:LYS:HB3	2.01	0.42
11:I:89:GLN:HE21	11:I:89:GLN:HB3	1.63	0.42
22:T:9:LYS:HG2	22:T:13:ARG:NH1	2.34	0.42
1:0:177:A:H2'	1:0:178:U:O4'	2.19	0.42
1:0:401:C:H2'	1:0:402:U:C6	2.53	0.42
1:0:688:A:N6	1:0:698:A:H5''	2.34	0.42
1:0:2276:U:H2'	1:0:2277:U:C6	2.54	0.42
1:0:1123:A:C6	1:0:1238:C:H5'	2.54	0.42
1:0:1902:G:H2'	1:0:1903:U:O4'	2.19	0.42
1:0:2415:A:C2	16:N:25:ARG:HB2	2.54	0.42
3:A:180:LYS:HE3	3:A:180:LYS:HB3	1.87	0.42
9:G:156:ILE:H	9:G:156:ILE:HG13	1.66	0.42
14:L:142:LEU:HD12	14:L:142:LEU:HA	1.84	0.42
15:M:30:GLU:O	15:M:34:GLU:HG3	2.19	0.42
1:0:1393:A:H2'	1:0:1394:C:C6	2.55	0.42
1:0:1534:C:O2'	1:0:1656:A:OP1	2.32	0.42
4:B:268:ARG:NE	40:B:592:HOH:O	2.52	0.42
9:G:112:ALA:HB1	9:G:187:LYS:HE2	2.00	0.42
20:R:111:ILE:HG23	20:R:145:LEU:HD11	2.01	0.42
1:0:95:A:H5''	1:0:97:G:O4'	2.19	0.42
1:0:1014:A:H2'	1:0:1015:C:H5'	2.01	0.42
1:0:1098:A:H2'	1:0:1099:G:O4'	2.20	0.42
1:0:1184:C:H2'	1:0:1185:U:C6	2.54	0.42
5:C:115:LEU:HD12	5:C:115:LEU:HA	1.85	0.42
11:I:74:LEU:HD11	11:I:107:GLN:HG2	2.01	0.42
11:I:96:LEU:HD21	11:I:101:VAL:HG22	2.02	0.42
18:P:75:LYS:O	18:P:80:ARG:NH2	2.48	0.42
22:T:3:GLN:HA	22:T:4:PRO:HD3	1.81	0.42
24:V:39:ALA:HA	24:V:40:PRO:HD3	1.81	0.42
1:0:1980:U:O2	1:0:2008:U:H4'	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:2898:G:H4'	4:B:288:GLY:HA2	2.01	0.42
13:K:120:ARG:HG2	13:K:121:PHE:CZ	2.54	0.42
20:R:99:ALA:CB	20:R:109:MET:HE1	2.42	0.42
1:O:638:C:H2'	1:O:639:A:C8	2.54	0.42
1:O:1687:C:O2	29:1:9:GLY:HA2	2.20	0.42
1:O:2090:G:H2'	1:O:2091:G:C8	2.55	0.42
1:O:2570:G:H5''	40:O:9612:HOH:O	2.20	0.42
1:O:2896:A:H2'	38:O:8212:ACY:H2	2.02	0.42
6:D:23:VAL:HB	6:D:130:VAL:HG22	2.00	0.42
11:I:70:PRO:HG2	11:I:73:GLU:HB2	2.02	0.42
13:K:78:LYS:HA	13:K:79:PRO:HD3	1.90	0.42
17:O:53:GLN:HG2	17:O:56:GLU:OE1	2.20	0.42
26:X:79:GLU:O	26:X:82:GLU:N	2.52	0.42
28:Z:57:CYS:SG	28:Z:59:TYR:HB3	2.60	0.42
1:O:417:G:H8	1:O:417:G:OP1	2.03	0.42
1:O:1132:A:N6	1:O:1229:C:H2'	2.35	0.42
1:O:2356:A:H2'	1:O:2357:G:O4'	2.19	0.42
6:D:81:GLU:OE2	6:D:170:TYR:OH	2.24	0.42
27:Y:122:ARG:NH1	40:Y:480:HOH:O	2.41	0.42
31:3:11:CYS:HB2	31:3:20:HIS:CE1	2.55	0.42
1:O:2587:OMU:H2'	1:O:2589:U:H5''	2.01	0.42
13:K:74:VAL:HB	13:K:113:ILE:HD12	2.02	0.42
18:P:80:ARG:HD2	18:P:87:ARG:CZ	2.50	0.42
25:W:3:ALA:O	25:W:54:PHE:HA	2.20	0.42
1:O:188:C:H5''	15:M:163:LEU:HD21	2.01	0.42
1:O:806:A:H2'	1:O:807:A:O4'	2.20	0.42
1:O:1449:G:C6	1:O:1450:C:N4	2.88	0.42
1:O:1515:A:H2'	1:O:1516:U:C6	2.55	0.42
1:O:1964:U:H2'	1:O:1965:C:H6	1.82	0.42
2:9:97:U:O2'	25:W:134:GLU:OE2	2.29	0.42
8:F:28:ALA:H	8:F:100:ASP:HB3	1.85	0.42
14:L:143:THR:HG21	40:L:327:HOH:O	2.20	0.42
1:O:2141:G:N2	1:O:2235:G:C5	2.88	0.41
2:9:91:C:H2'	2:9:92:G:O4'	2.20	0.41
3:A:86:ALA:HB3	3:A:94:LEU:HG	2.01	0.41
4:B:267:LYS:NZ	4:B:300:SER:O	2.44	0.41
25:W:126:ASP:HB3	25:W:135:GLY:O	2.19	0.41
2:9:35:C:H5''	40:9:370:HOH:O	2.19	0.41
4:B:209:LYS:HB2	4:B:257:THR:OG1	2.20	0.41
6:D:60:GLU:HG2	6:D:61:PHE:CD1	2.55	0.41
9:G:113:PRO:HG3	9:G:185:GLU:OE2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:O:32:ARG:O	17:O:35:LYS:HB2	2.20	0.41
1:O:2668:G:H2'	1:O:2669:U:C6	2.55	0.41
6:D:44:ILE:HD11	6:D:114:PHE:CZ	2.55	0.41
9:G:205:PRO:O	9:G:209:GLU:HB2	2.20	0.41
13:K:27:ARG:NH2	40:K:334:HOH:O	2.38	0.41
17:O:22:GLY:O	40:O:325:HOH:O	2.21	0.41
25:W:113:SER:HA	25:W:114:PRO:HD3	1.86	0.41
1:O:591:A:OP2	40:O:9919:HOH:O	2.21	0.41
1:O:1257:C:H2'	1:O:1258:G:O4'	2.21	0.41
1:O:1589:G:N2	1:O:1605:G:H1'	2.35	0.41
1:O:2272:G:OP1	3:A:223:ARG:HD2	2.21	0.41
10:H:77:ILE:HG23	10:H:82:GLU:HA	2.01	0.41
10:H:99:ARG:HB3	10:H:121:GLY:HA3	2.02	0.41
10:H:175:LEU:HD23	10:H:175:LEU:HA	1.79	0.41
15:M:34:GLU:HB3	15:M:38:GLU:HG3	2.03	0.41
19:Q:25:PRO:HA	19:Q:26:PRO:HD3	1.80	0.41
21:S:18:MET:HE2	21:S:18:MET:HB3	1.90	0.41
26:X:6:GLU:HG3	26:X:76:ARG:HB3	2.02	0.41
1:O:734:U:O2'	1:O:737:A:N6	2.53	0.41
1:O:2365:G:O2'	1:O:2370:A:N1	2.49	0.41
6:D:127:GLY:C	6:D:128:LEU:HD13	2.46	0.41
9:G:51:ARG:HG2	11:I:115:TYR:HE1	1.84	0.41
10:H:49:GLN:O	10:H:169:GLU:N	2.46	0.41
10:H:50:ILE:HD12	10:H:149:VAL:CG2	2.50	0.41
25:W:80:ASP:HB2	40:W:325:HOH:O	2.20	0.41
29:1:25:LYS:HD2	30:2:48:ASP:HA	2.03	0.41
1:O:344:C:H2'	1:O:345:G:O4'	2.20	0.41
1:O:644:G:N3	1:O:644:G:H5'	2.35	0.41
1:O:1249:U:H2'	1:O:1250:C:C6	2.56	0.41
1:O:1684:A:O2'	1:O:1685:A:H5''	2.21	0.41
2:9:96:C:H2'	2:9:97:U:C6	2.56	0.41
5:C:214:THR:HB	40:C:439:HOH:O	2.20	0.41
9:G:131:THR:OG1	9:G:132:GLY:N	2.54	0.41
1:O:255:A:H2'	1:O:256:C:C6	2.55	0.41
1:O:797:A:H5'	28:Z:9:GLY:O	2.21	0.41
1:O:821:U:H2'	1:O:822:C:H6	1.86	0.41
1:O:858:U:H2'	1:O:859:C:H6	1.86	0.41
1:O:1006:A:H2'	1:O:1007:A:C8	2.56	0.41
1:O:1066:U:H2'	1:O:1067:A:C8	2.56	0.41
1:O:1418:U:OP1	30:2:42:TRP:HB3	2.21	0.41
1:O:2365:G:H4'	19:Q:45:PRO:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:0:2597:U:OP2	40:0:8619:HOH:O	2.21	0.41
1:0:2626:C:H2'	1:0:2627:G:C8	2.56	0.41
2:9:28:U:OP2	16:N:39:SER:HB2	2.20	0.41
2:9:29:C:O3'	6:D:138:GLY:HA2	2.21	0.41
2:9:92:G:C6	2:9:93:A:C6	3.09	0.41
6:D:14:ARG:HH21	6:D:136:ARG:CZ	2.33	0.41
6:D:135:VAL:HG23	40:D:209:HOH:O	2.20	0.41
8:F:18:GLU:O	8:F:22:VAL:HG12	2.20	0.41
11:I:64:ILE:HG12	11:I:65:GLU:N	2.36	0.41
11:I:102:LYS:O	11:I:106:GLU:HG2	2.20	0.41
15:M:49:ALA:C	15:M:54:TYR:HB3	2.46	0.41
20:R:68:HIS:CG	20:R:76:ASP:HB2	2.55	0.41
1:0:1167:G:H4'	11:I:131:LEU:HG	2.02	0.41
1:0:1185:U:H2'	1:0:1186:C:H6	1.85	0.41
2:9:23:U:O2'	2:9:24:U:H5'	2.20	0.41
2:9:27:C:H2'	2:9:28:U:O4'	2.20	0.41
4:B:179:LEU:O	4:B:183:GLU:HG2	2.21	0.41
1:0:1497:G:H4'	1:0:1627:G:O2'	2.21	0.40
6:D:60:GLU:C	6:D:62:ASP:H	2.28	0.40
9:G:112:ALA:HB3	9:G:187:LYS:O	2.20	0.40
26:X:30:MET:HE3	26:X:58:ALA:CB	2.50	0.40
27:Y:117:LEU:CA	27:Y:174:VAL:HG21	2.51	0.40
1:0:615:G:OP1	12:J:91:LYS:HE3	2.22	0.40
1:0:625:U:H5''	1:0:1044:C:N4	2.37	0.40
1:0:2105:C:H2'	1:0:2106:C:C6	2.56	0.40
1:0:2575:C:H2'	1:0:2576:A:O4'	2.21	0.40
1:0:2717:C:OP1	4:B:207:LYS:HG3	2.21	0.40
2:9:93:A:C5	2:9:94:G:H1'	2.57	0.40
3:A:175:LYS:HE2	40:A:475:HOH:O	2.20	0.40
19:Q:42:LYS:HA	19:Q:42:LYS:HD2	1.98	0.40
24:V:37:GLY:N	24:V:38:GLY:HA3	2.37	0.40
27:Y:126:PRO:HG2	27:Y:128:PHE:CE2	2.56	0.40
1:0:1422:U:H2'	1:0:1423:C:C6	2.56	0.40
1:0:2401:A:H2'	1:0:2402:A:C8	2.57	0.40
22:T:50:VAL:HG12	22:T:56:ALA:HA	2.03	0.40
1:0:113:A:C8	1:0:114:A:C8	3.09	0.40
1:0:1385:G:H5'	26:X:50:LEU:O	2.20	0.40
1:0:1596:U:H2'	1:0:1598:A:OP2	2.20	0.40
1:0:1985:U:C2	1:0:1996:U:O4'	2.75	0.40
1:0:2252:A:C5	1:0:2253:G:H1'	2.56	0.40
1:0:2694:A:H4'	7:E:91:PHE:CE2	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:W:146:ILE:HD12	25:W:146:ILE:HA	1.93	0.40
27:Y:213:LYS:HE3	27:Y:213:LYS:HB2	1.94	0.40
1:O:1993:C:H5''	13:K:66:ARG:NH1	2.37	0.40
1:O:2698:G:H2'	1:O:2699:A:C8	2.57	0.40
1:O:2819:C:H2'	1:O:2820:A:C8	2.56	0.40
2:9:73:A:H61	2:9:108:C:N4	2.16	0.40
6:D:44:ILE:HD11	6:D:114:PHE:CE1	2.57	0.40
9:G:75:ASP:O	9:G:76:ASP:C	2.63	0.40
10:H:30:LYS:HG3	10:H:62:HIS:CE1	2.57	0.40
10:H:102:LYS:HG3	10:H:103:GLN:H	1.87	0.40
12:J:39:VAL:HG22	12:J:107:ASN:HA	2.04	0.40
16:N:180:LEU:HD23	16:N:180:LEU:HA	1.92	0.40
20:R:106:GLY:HA2	20:R:109:MET:CE	2.52	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	
3	A	235/240 (98%)	223 (95%)	11 (5%)	1 (0%)	30	43
4	B	335/338 (99%)	327 (98%)	8 (2%)	0	100	100
5	C	244/246 (99%)	238 (98%)	6 (2%)	0	100	100
6	D	164/177 (93%)	143 (87%)	18 (11%)	3 (2%)	6	9
7	E	172/178 (97%)	168 (98%)	4 (2%)	0	100	100
8	F	117/120 (98%)	115 (98%)	2 (2%)	0	100	100
9	G	204/255 (80%)	171 (84%)	28 (14%)	5 (2%)	4	5
10	H	161/177 (91%)	156 (97%)	4 (2%)	1 (1%)	21	32
11	I	141/162 (87%)	122 (86%)	17 (12%)	2 (1%)	9	13
12	J	142/145 (98%)	138 (97%)	4 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
13	K	130/132 (98%)	124 (95%)	6 (5%)	0	100	100
14	L	145/165 (88%)	139 (96%)	5 (3%)	1 (1%)	18	28
15	M	193/196 (98%)	189 (98%)	4 (2%)	0	100	100
16	N	184/187 (98%)	179 (97%)	5 (3%)	0	100	100
17	O	113/116 (97%)	112 (99%)	1 (1%)	0	100	100
18	P	142/149 (95%)	142 (100%)	0	0	100	100
19	Q	93/96 (97%)	93 (100%)	0	0	100	100
20	R	148/155 (96%)	146 (99%)	2 (1%)	0	100	100
21	S	79/85 (93%)	78 (99%)	1 (1%)	0	100	100
22	T	117/120 (98%)	116 (99%)	1 (1%)	0	100	100
23	U	54/67 (81%)	53 (98%)	1 (2%)	0	100	100
24	V	63/71 (89%)	62 (98%)	1 (2%)	0	100	100
25	W	152/154 (99%)	152 (100%)	0	0	100	100
26	X	88/92 (96%)	83 (94%)	5 (6%)	0	100	100
27	Y	142/241 (59%)	142 (100%)	0	0	100	100
28	Z	81/92 (88%)	78 (96%)	3 (4%)	0	100	100
29	1	54/57 (95%)	53 (98%)	1 (2%)	0	100	100
30	2	47/50 (94%)	46 (98%)	1 (2%)	0	100	100
31	3	90/92 (98%)	88 (98%)	2 (2%)	0	100	100
33	6	53/56 (95%)	48 (91%)	5 (9%)	0	100	100
All	All	4083/4411 (93%)	3924 (96%)	146 (4%)	13 (0%)	36	50

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
9	G	38	ILE
6	D	173	GLU
9	G	74	VAL
9	G	53	LEU
9	G	75	ASP
10	H	176	ILE
11	I	25	THR
11	I	28	PRO
3	A	170	VAL
14	L	81	VAL

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Mol	Chain	Res	Type
6	D	63	ILE
6	D	66	GLY
9	G	114	ILE

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	
3	A	179/182 (98%)	169 (94%)	10 (6%)	19	33
4	B	282/283 (100%)	262 (93%)	20 (7%)	13	23
5	C	193/193 (100%)	181 (94%)	12 (6%)	16	29
6	D	139/148 (94%)	125 (90%)	14 (10%)	7	11
7	E	154/156 (99%)	147 (96%)	7 (4%)	24	42
8	F	93/94 (99%)	86 (92%)	7 (8%)	12	21
9	G	174/174 (100%)	154 (88%)	20 (12%)	5	8
10	H	138/145 (95%)	133 (96%)	5 (4%)	31	52
11	I	120/130 (92%)	111 (92%)	9 (8%)	12	21
12	J	120/121 (99%)	114 (95%)	6 (5%)	22	38
13	K	106/106 (100%)	101 (95%)	5 (5%)	23	41
14	L	114/127 (90%)	106 (93%)	8 (7%)	14	24
15	M	159/160 (99%)	153 (96%)	6 (4%)	29	49
16	N	149/150 (99%)	136 (91%)	13 (9%)	9	16
17	O	93/94 (99%)	85 (91%)	8 (9%)	10	16
18	P	114/117 (97%)	109 (96%)	5 (4%)	25	43
19	Q	79/80 (99%)	78 (99%)	1 (1%)	61	80
20	R	117/122 (96%)	111 (95%)	6 (5%)	21	37
21	S	71/74 (96%)	70 (99%)	1 (1%)	59	79
22	T	105/106 (99%)	100 (95%)	5 (5%)	23	40
23	U	47/53 (89%)	45 (96%)	2 (4%)	26	44

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
24	V	51/57 (90%)	49 (96%)	2 (4%)	28	48
25	W	130/130 (100%)	126 (97%)	4 (3%)	35	57
26	X	72/74 (97%)	68 (94%)	4 (6%)	19	33
27	Y	122/196 (62%)	116 (95%)	6 (5%)	22	39
28	Z	67/74 (90%)	65 (97%)	2 (3%)	36	58
29	1	46/47 (98%)	46 (100%)	0	100	100
30	2	45/46 (98%)	44 (98%)	1 (2%)	45	67
31	3	79/79 (100%)	75 (95%)	4 (5%)	21	37
33	6	48/49 (98%)	48 (100%)	0	100	100
All	All	3406/3567 (96%)	3213 (94%)	193 (6%)	18	33

All (193) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	30	ARG
3	A	31	LYS
3	A	38	ILE
3	A	55	VAL
3	A	69	LEU
3	A	105	VAL
3	A	133	ARG
3	A	175	LYS
3	A	192	VAL
3	A	217	ARG
4	B	11	LEU
4	B	51	VAL
4	B	53	LEU
4	B	71	VAL
4	B	84	LEU
4	B	97	LEU
4	B	98	THR
4	B	102	THR
4	B	175	LEU
4	B	195	ARG
4	B	245	SER
4	B	251	VAL
4	B	254	GLN
4	B	257	THR
4	B	264	GLU

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Mol	Chain	Res	Type
4	B	265	LEU
4	B	279	THR
4	B	280	VAL
4	B	289	GLU
4	B	310	ARG
5	C	21	VAL
5	C	59	GLU
5	C	115	LEU
5	C	136	VAL
5	C	140	VAL
5	C	162	VAL
5	C	214	THR
5	C	223	LEU
5	C	234	VAL
5	C	236	THR
5	C	240	LEU
5	C	243	VAL
6	D	18	ILE
6	D	23	VAL
6	D	24	HIS
6	D	34	LEU
6	D	44	ILE
6	D	52	THR
6	D	55	LYS
6	D	58	VAL
6	D	60	GLU
6	D	69	ILE
6	D	82	GLU
6	D	84	LEU
6	D	128	LEU
6	D	172	VAL
7	E	3	VAL
7	E	12	ASP
7	E	46	THR
7	E	86	VAL
7	E	102	VAL
7	E	132	THR
7	E	154	ILE
8	F	1	PRO
8	F	12	LEU
8	F	22	VAL
8	F	60	VAL

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Mol	Chain	Res	Type
8	F	65	GLU
8	F	99	THR
8	F	118	LEU
9	G	10	GLU
9	G	36	VAL
9	G	46	LEU
9	G	51	ARG
9	G	74	VAL
9	G	75	ASP
9	G	81	LEU
9	G	86	THR
9	G	89	VAL
9	G	101	LEU
9	G	106	GLU
9	G	110	THR
9	G	124	ILE
9	G	153	GLU
9	G	164	VAL
9	G	167	THR
9	G	169	GLU
9	G	173	GLN
9	G	188	GLU
9	G	210	LEU
10	H	4	LYS
10	H	21	GLU
10	H	61	ARG
10	H	65	LEU
10	H	157	TYR
11	I	7	VAL
11	I	31	VAL
11	I	41	GLN
11	I	74	LEU
11	I	96	LEU
11	I	98	VAL
11	I	112	LEU
11	I	135	ILE
11	I	136	GLU
12	J	39	VAL
12	J	45	VAL
12	J	52	GLN
12	J	54	VAL
12	J	80	LYS

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Mol	Chain	Res	Type
12	J	120	SER
13	K	55	VAL
13	K	63	GLU
13	K	98	VAL
13	K	115	ARG
13	K	130	MET
14	L	18	HIS
14	L	75	LEU
14	L	80	ASP
14	L	91	VAL
14	L	99	GLU
14	L	140	VAL
14	L	145	LEU
14	L	149	ARG
15	M	22	GLU
15	M	46	LEU
15	M	61	ILE
15	M	93	ARG
15	M	94	ARG
15	M	164	THR
16	N	17	ARG
16	N	26	LEU
16	N	43	VAL
16	N	49	THR
16	N	50	LEU
16	N	66	LEU
16	N	110	THR
16	N	127	LEU
16	N	135	VAL
16	N	154	LEU
16	N	168	LEU
16	N	175	LEU
16	N	186	LEU
17	O	21	SER
17	O	25	VAL
17	O	28	ASP
17	O	32	ARG
17	O	42	GLU
17	O	47	ARG
17	O	96	VAL
17	O	98	LEU
18	P	16	VAL

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Mol	Chain	Res	Type
18	P	21	VAL
18	P	52	LYS
18	P	80	ARG
18	P	91	LYS
19	Q	64	GLU
20	R	39	THR
20	R	55	GLN
20	R	82	GLU
20	R	109	MET
20	R	119	VAL
20	R	143	VAL
21	S	21	GLN
22	T	9	LYS
22	T	26	THR
22	T	48	VAL
22	T	71	VAL
22	T	96	VAL
23	U	28	THR
23	U	52	THR
24	V	12	THR
24	V	21	ASP
25	W	35	VAL
25	W	48	VAL
25	W	52	VAL
25	W	73	LEU
26	X	8	ARG
26	X	43	VAL
26	X	86	GLU
26	X	91	GLU
27	Y	172	THR
27	Y	186	ARG
27	Y	187	VAL
27	Y	203	VAL
27	Y	236	VAL
27	Y	238	VAL
28	Z	83	ILE
28	Z	87	LEU
30	2	8	LYS
31	3	3	MET
31	3	65	THR
31	3	73	GLU
31	3	92	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	5	GLN
3	A	134	ASN
3	A	218	ASN
5	C	39	GLN
6	D	158	ASN
7	E	66	GLN
7	E	150	GLN
7	E	173	ASN
10	H	75	HIS
10	H	131	GLN
11	I	15	ASN
11	I	103	GLN
12	J	142	ASN
13	K	119	GLN
16	N	22	GLN
20	R	101	HIS
21	S	51	GLN
22	T	11	GLN
27	Y	149	GLN
31	3	20	HIS
31	3	91	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	0	2804/2910 (96%)	268 (9%)	23 (0%)
2	9	121/122 (99%)	16 (13%)	1 (0%)
All	All	2925/3032 (96%)	284 (9%)	24 (0%)

All (284) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	0	9	C
1	0	17	G
1	0	31	C
1	0	67	A
1	0	70	A
1	0	71	G
1	0	87	C
1	0	92	G

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Mol	Chain	Res	Type
1	0	113	A
1	0	114	A
1	0	115	U
1	0	120	A
1	0	126	C
1	0	127	U
1	0	128	A
1	0	138	U
1	0	140	G
1	0	141	C
1	0	151	A
1	0	166	A
1	0	169	A
1	0	170	U
1	0	185	G
1	0	186	A
1	0	191	A
1	0	192	A
1	0	210	U
1	0	219	G
1	0	237	G
1	0	256	C
1	0	271	C
1	0	273	G
1	0	308	U
1	0	309	C
1	0	331	A
1	0	337	A
1	0	338	C
1	0	345	G
1	0	358	G
1	0	396	U
1	0	397	A
1	0	417	G
1	0	418	C
1	0	457	U
1	0	461	C
1	0	487	G
1	0	510	U
1	0	511	A
1	0	514	G
1	0	537	G

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Mol	Chain	Res	Type
1	0	538	C
1	0	539	G
1	0	553	G
1	0	581	G
1	0	588	G
1	0	604	G
1	0	620	A
1	0	632	A
1	0	644	G
1	0	660	A
1	0	698	A
1	0	699	C
1	0	700	A
1	0	701	U
1	0	702	G
1	0	715	U
1	0	716	G
1	0	717	C
1	0	759	C
1	0	777	U
1	0	809	G
1	0	821	U
1	0	835	U
1	0	840	U
1	0	857	A
1	0	858	U
1	0	868	G
1	0	869	G
1	0	875	A
1	0	877	G
1	0	878	G
1	0	884	C
1	0	885	G
1	0	898	G
1	0	905	C
1	0	920	C
1	0	921	G
1	0	923	A
1	0	925	C
1	0	961	A
1	0	998	U
1	0	1003	U

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Mol	Chain	Res	Type
1	0	1006	A
1	0	1008	C
1	0	1029	U
1	0	1045	G
1	0	1052	G
1	0	1059	G
1	0	1072	G
1	0	1081	A
1	0	1087	G
1	0	1088	A
1	0	1109	U
1	0	1110	G
1	0	1119	G
1	0	1120	U
1	0	1130	U
1	0	1137	G
1	0	1164	U
1	0	1166	A
1	0	1173	A
1	0	1174	A
1	0	1175	G
1	0	1176	C
1	0	1187	U
1	0	1192	A
1	0	1193	A
1	0	1194	A
1	0	1202	A
1	0	1216	G
1	0	1226	G
1	0	1237	U
1	0	1238	C
1	0	1239	G
1	0	1242	A
1	0	1252	A
1	0	1279	U
1	0	1280	A
1	0	1289	C
1	0	1299	G
1	0	1342	C
1	0	1360	C
1	0	1380	U
1	0	1386	G

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Mol	Chain	Res	Type
1	0	1406	A
1	0	1407	A
1	0	1409	G
1	0	1451	C
1	0	1477	C
1	0	1485	A
1	0	1492	A
1	0	1525	G
1	0	1526	A
1	0	1528	A
1	0	1592	G
1	0	1611	G
1	0	1626	A
1	0	1633	C
1	0	1634	G
1	0	1635	U
1	0	1640	C
1	0	1656	A
1	0	1684	A
1	0	1685	A
1	0	1692	C
1	0	1693	A
1	0	1701	A
1	0	1702	U
1	0	1703	G
1	0	1710	A
1	0	1722	U
1	0	1723	G
1	0	1725	C
1	0	1730	G
1	0	1731	C
1	0	1752	G
1	0	1778	A
1	0	1798	C
1	0	1820	G
1	0	1829	A
1	0	1856	C
1	0	1857	A
1	0	1878	G
1	0	1919	A
1	0	1958	U
1	0	1959	G

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Mol	Chain	Res	Type
1	0	1963	C
1	0	1971	G
1	0	1978	A
1	0	1979	G
1	0	1996	U
1	0	2004	U
1	0	2005	G
1	0	2008	U
1	0	2011	A
1	0	2012	U
1	0	2013	G
1	0	2033	G
1	0	2034	U
1	0	2064	U
1	0	2072	G
1	0	2073	G
1	0	2074	A
1	0	2096	A
1	0	2101	A
1	0	2102	G
1	0	2110	G
1	0	2144	C
1	0	2146	C
1	0	2148	G
1	0	2227	C
1	0	2228	G
1	0	2229	G
1	0	2238	A
1	0	2243	C
1	0	2253	G
1	0	2258	A
1	0	2271	G
1	0	2272	G
1	0	2291	A
1	0	2294	C
1	0	2313	C
1	0	2317	C
1	0	2321	A
1	0	2339	A
1	0	2342	G
1	0	2354	A
1	0	2369	A

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Mol	Chain	Res	Type
1	0	2378	U
1	0	2379	G
1	0	2404	G
1	0	2420	G
1	0	2422	U
1	0	2462	G
1	0	2467	A
1	0	2468	A
1	0	2469	A
1	0	2476	C
1	0	2480	G
1	0	2483	A
1	0	2505	G
1	0	2527	U
1	0	2536	C
1	0	2537	G
1	0	2541	U
1	0	2553	A
1	0	2564	G
1	0	2589	U
1	0	2601	A
1	0	2602	G
1	0	2608	C
1	0	2637	A
1	0	2638	G
1	0	2665	A
1	0	2676	C
1	0	2681	A
1	0	2726	U
1	0	2747	C
1	0	2748	G
1	0	2749	U
1	0	2750	G
1	0	2762	C
1	0	2768	A
1	0	2792	A
1	0	2800	A
1	0	2811	A
1	0	2812	A
1	0	2825	C
1	0	2827	A
1	0	2837	U

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Mol	Chain	Res	Type
1	0	2852	A
1	0	2876	G
1	0	2890	A
1	0	2896	A
1	0	2903	C
1	0	2909	G
1	0	2914	A
1	0	2915	A
2	9	3	A
2	9	7	G
2	9	11	A
2	9	23	U
2	9	24	U
2	9	32	G
2	9	44	A
2	9	52	A
2	9	55	U
2	9	56	A
2	9	57	A
2	9	66	G
2	9	77	A
2	9	94	G
2	9	114	G
2	9	122	C

All (24) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	0	8	A
1	0	169	A
1	0	417	G
1	0	509	A
1	0	699	C
1	0	715	U
1	0	716	G
1	0	834	G
1	0	857	A
1	0	877	G
1	0	1080	C
1	0	1246	A
1	0	1504	A
1	0	1524	U

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Mol	Chain	Res	Type
1	0	1684	A
1	0	1692	C
1	0	1856	C
1	0	1979	G
1	0	2467	A
1	0	2536	C
1	0	2553	A
1	0	2637	A
1	0	2791	U
2	9	2	U

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

5 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$
1	OMG	0	2588	1	23,26,27	1.16	3 (13%)	32,38,41	1.93	6 (18%)
1	OMU	0	2587	1	19,22,23	1.22	3 (15%)	25,31,34	1.83	6 (24%)
1	PSU	0	2621	1	18,21,22	1.40	2 (11%)	21,30,33	1.71	4 (19%)
1	1MA	0	628	36,1	21,25,26	1.28	2 (9%)	30,37,40	1.77	6 (20%)
1	UR3	0	2619	1	19,22,23	1.03	1 (5%)	26,32,35	1.60	3 (11%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	OMG	0	2588	1	-	0/9/27/28	0/3/3/3
1	OMU	0	2587	1	-	0/9/27/28	0/2/2/2
1	PSU	0	2621	1	-	0/7/25/26	0/2/2/2
1	1MA	0	628	36,1	-	2/7/25/26	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	UR3	0	2619	1	-	0/7/25/26	0/2/2/2

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	0	2621	PSU	C6-C5	4.02	1.39	1.35
1	0	628	1MA	C6-N6	3.21	1.35	1.28
1	0	2588	OMG	C5-C4	3.19	1.47	1.38
1	0	628	1MA	C5-C4	2.98	1.46	1.38
1	0	2587	OMU	C4-N3	-2.59	1.34	1.38
1	0	2619	UR3	C6-C5	2.32	1.40	1.35
1	0	2588	OMG	C6-N1	-2.14	1.34	1.38
1	0	2587	OMU	O2-C2	2.12	1.26	1.23
1	0	2587	OMU	C6-C5	2.10	1.40	1.35
1	0	2621	PSU	O2-C2	2.02	1.27	1.23
1	0	2588	OMG	C4-N9	-2.00	1.33	1.38

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	0	2619	UR3	C4-N3-C2	-6.47	119.37	124.58
1	0	2588	OMG	C5-C4-N3	-5.38	119.83	128.39
1	0	628	1MA	C5-C4-N3	-5.03	119.87	127.27
1	0	2621	PSU	N1-C2-N3	4.98	120.43	115.17
1	0	2587	OMU	C4-N3-C2	-4.87	120.57	126.61
1	0	628	1MA	C2-N3-C4	4.64	121.63	112.53
1	0	2588	OMG	C2-N3-C4	4.57	120.17	112.30
1	0	2587	OMU	N3-C2-N1	4.19	120.35	114.89
1	0	2588	OMG	N9-C4-N3	4.04	134.02	125.95
1	0	2587	OMU	C5-C4-N3	3.80	120.12	114.80
1	0	2588	OMG	C6-C5-N7	3.53	136.72	130.29
1	0	2587	OMU	O2-C2-N1	-3.31	118.49	122.80
1	0	2621	PSU	C4-N3-C2	-3.13	122.05	126.37
1	0	2621	PSU	O2-C2-N1	-3.08	119.62	122.79
1	0	628	1MA	N9-C4-N3	2.99	133.70	126.90
1	0	2619	UR3	C5-C4-N3	2.95	118.93	115.04
1	0	2587	OMU	O4-C4-C5	-2.93	120.11	125.16
1	0	628	1MA	C6-C5-N7	2.74	137.00	132.16
1	0	628	1MA	C4-C5-N7	-2.73	106.35	110.67
1	0	2588	OMG	C4-C5-N7	-2.60	106.54	110.67
1	0	628	1MA	N1-C2-N3	-2.42	123.12	126.00
1	0	2621	PSU	C6-C5-C4	-2.38	116.57	118.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	0	2619	UR3	C3U-N3-C4	2.32	121.08	117.87
1	0	2588	OMG	O6-C6-C5	-2.16	120.82	126.53
1	0	2587	OMU	C5-C6-N1	-2.05	118.50	121.84

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	0	628	1MA	C2'-C1'-N9-C8
1	0	628	1MA	C2'-C1'-N9-C4

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	0	2587	OMU	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 265 ligands modelled in this entry, 261 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
38	ACY	W	201	-	3,3,3	0.80	0	3,3,3	0.87	0
38	ACY	0	8213	-	3,3,3	0.73	0	3,3,3	1.10	0
38	ACY	0	8212	-	3,3,3	0.81	0	3,3,3	0.92	0
38	ACY	0	8214	-	3,3,3	0.91	0	3,3,3	0.55	0

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.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
38	W	201	ACY	1	0
38	0	8213	ACY	3	0
38	0	8212	ACY	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
9	G	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	G	211:ASP	C	213:UNK	N	3.25

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	0	2803/2910 (96%)	-0.99	19 (0%) 84 81	9, 28, 77, 221	0
2	9	122/122 (100%)	-0.69	3 (2%) 58 54	24, 44, 61, 127	0
3	A	237/240 (98%)	-0.37	3 (1%) 75 71	12, 30, 63, 98	0
4	B	337/338 (99%)	-0.25	1 (0%) 90 88	14, 35, 60, 72	0
5	C	246/246 (100%)	-0.47	0 100 100	10, 27, 48, 58	0
6	D	166/177 (93%)	1.38	40 (24%) 2 1	33, 81, 106, 113	0
7	E	174/178 (97%)	0.20	2 (1%) 78 74	30, 47, 64, 83	0
8	F	119/120 (99%)	0.16	0 100 100	27, 46, 72, 78	0
9	G	206/255 (80%)	2.06	97 (47%) 0 0	57, 105, 173, 186	0
10	H	165/177 (93%)	-0.06	5 (3%) 52 48	22, 37, 70, 94	0
11	I	145/162 (89%)	2.22	78 (53%) 0 0	81, 115, 155, 179	0
12	J	144/145 (99%)	-0.21	3 (2%) 63 59	23, 33, 55, 82	0
13	K	132/132 (100%)	-0.30	1 (0%) 82 80	19, 32, 49, 63	0
14	L	149/165 (90%)	0.12	10 (6%) 24 20	12, 41, 80, 88	0
15	M	195/196 (99%)	-0.66	0 100 100	14, 22, 34, 47	0
16	N	186/187 (99%)	0.04	3 (1%) 70 66	25, 41, 76, 101	0
17	O	115/116 (99%)	-0.15	1 (0%) 81 78	23, 34, 50, 55	0
18	P	144/149 (96%)	-0.25	1 (0%) 84 81	23, 34, 48, 60	0
19	Q	95/96 (98%)	-0.51	1 (1%) 78 74	19, 26, 41, 50	0
20	R	150/155 (96%)	-0.64	0 100 100	16, 26, 45, 56	0
21	S	81/85 (95%)	-0.19	1 (1%) 76 73	23, 36, 54, 63	0
22	T	119/120 (99%)	-0.24	0 100 100	21, 36, 55, 76	0
23	U	56/67 (83%)	-0.15	2 (3%) 46 42	28, 36, 60, 103	0
24	V	65/71 (91%)	0.23	4 (6%) 26 23	30, 46, 78, 114	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
25	W	154/154 (100%)	-0.31	0 100 100	22, 32, 47, 56	0
26	X	90/92 (97%)	0.42	10 (11%) 10 7	26, 39, 101, 111	0
27	Y	144/241 (59%)	-0.41	5 (3%) 47 43	15, 28, 53, 104	0
28	Z	83/92 (90%)	0.13	3 (3%) 46 42	25, 39, 85, 108	0
29	1	56/57 (98%)	-0.94	0 100 100	11, 16, 23, 24	0
30	2	49/50 (98%)	0.10	1 (2%) 65 60	20, 37, 64, 73	0
31	3	92/92 (100%)	-0.55	0 100 100	16, 32, 46, 52	0
32	4	0/70	-	-	-	-
32	5	0/70	-	-	-	-
33	6	55/56 (98%)	1.69	17 (30%) 1 1	56, 98, 117, 122	0
All	All	7074/7583 (93%)	-0.37	311 (4%) 39 34	9, 33, 102, 221	0

All (311) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
6	D	174	VAL	7.3
26	X	2	ALA	6.7
9	G	208	LEU	6.6
27	Y	238	VAL	6.0
11	I	81	PHE	5.9
11	I	5	ILE	5.8
16	N	154	LEU	5.6
9	G	53	LEU	5.5
9	G	210	LEU	5.5
9	G	133	VAL	5.3
10	H	176	ILE	5.3
10	H	177	ALA	5.2
9	G	135	PRO	5.2
9	G	110	THR	4.9
9	G	151	ILE	4.9
11	I	11	GLY	4.9
6	D	63	ILE	4.8
11	I	8	LEU	4.6
9	G	38	ILE	4.6
11	I	29	VAL	4.5
11	I	7	VAL	4.5
14	L	151	ALA	4.5
6	D	172	VAL	4.3
9	G	201	VAL	4.3

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Mol	Chain	Res	Type	RSRZ
14	L	84	VAL	4.3
33	6	56	ALA	4.3
9	G	74	VAL	4.3
11	I	50	VAL	4.3
9	G	137	PRO	4.2
6	D	125	ILE	4.2
9	G	126	ILE	4.2
9	G	131	THR	4.1
9	G	159	LEU	4.1
11	I	135	ILE	3.9
9	G	203	PHE	3.9
11	I	137	GLY	3.9
9	G	101	LEU	3.9
11	I	105	ALA	3.9
11	I	9	VAL	3.8
11	I	98	VAL	3.8
6	D	10	PHE	3.8
26	X	91	GLU	3.8
11	I	110	PRO	3.8
26	X	3	SER	3.7
9	G	211	ASP	3.7
11	I	43	ALA	3.7
9	G	145	VAL	3.7
12	J	2	SER	3.7
9	G	179	LEU	3.7
6	D	119	TYR	3.7
1	0	2637	A	3.6
11	I	18	PRO	3.6
11	I	56	TYR	3.6
9	G	149	ALA	3.6
11	I	74	LEU	3.6
11	I	38	ILE	3.5
33	6	55	VAL	3.5
11	I	156	ALA	3.5
4	B	1	PRO	3.5
11	I	31	VAL	3.5
33	6	23	VAL	3.5
1	0	127	U	3.5
11	I	51	PRO	3.5
33	6	26	PRO	3.5
6	D	11	HIS	3.5
9	G	77	GLY	3.4

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Mol	Chain	Res	Type	RSRZ
11	I	25	THR	3.4
9	G	138	PHE	3.4
9	G	139	VAL	3.4
11	I	34	VAL	3.4
3	A	237	GLY	3.4
11	I	75	ILE	3.4
6	D	170	TYR	3.4
6	D	61	PHE	3.4
28	Z	7	LYS	3.3
24	V	39	ALA	3.3
1	O	715	U	3.3
17	O	1	SER	3.3
9	G	121	PRO	3.3
6	D	121	PRO	3.2
9	G	142	LEU	3.2
9	G	202	LEU	3.2
9	G	147	ALA	3.2
28	Z	87	LEU	3.2
11	I	155	PHE	3.2
9	G	75	ASP	3.2
9	G	175	LEU	3.2
11	I	154	VAL	3.2
9	G	114	ILE	3.2
9	G	124	ILE	3.2
9	G	113	PRO	3.1
11	I	139	PRO	3.1
12	J	3	VAL	3.1
11	I	32	GLN	3.1
11	I	89	GLN	3.1
11	I	10	PRO	3.1
16	N	186	LEU	3.1
9	G	168	GLY	3.1
9	G	39	ALA	3.1
11	I	14	ALA	3.1
9	G	160	SER	3.1
9	G	189	VAL	3.1
14	L	82	ALA	3.0
24	V	40	PRO	3.0
9	G	191	LEU	3.0
21	S	81	ILE	3.0
27	Y	95	THR	3.0
11	I	28	PRO	3.0

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Mol	Chain	Res	Type	RSRZ
9	G	155	SER	3.0
11	I	83	THR	3.0
11	I	132	GLY	3.0
11	I	66	VAL	3.0
9	G	165	LEU	3.0
2	9	1	U	3.0
6	D	123	ILE	2.9
12	J	4	ALA	2.9
1	0	2917	C	2.9
9	G	99	PHE	2.9
9	G	136	GLY	2.9
9	G	205	PRO	2.9
11	I	94	ALA	2.9
6	D	69	ILE	2.9
27	Y	237	GLU	2.9
9	G	171	VAL	2.8
1	0	8	A	2.8
9	G	98	PRO	2.8
11	I	72	ALA	2.8
9	G	134	ASP	2.8
16	N	184	ILE	2.8
9	G	162	SER	2.8
11	I	42	THR	2.8
11	I	62	PHE	2.8
6	D	126	TYR	2.8
11	I	68	VAL	2.8
9	G	194	ARG	2.8
9	G	6	GLU	2.7
1	0	126	C	2.7
9	G	156	ILE	2.7
11	I	93	VAL	2.7
11	I	101	VAL	2.7
9	G	76	ASP	2.7
1	0	128	A	2.7
9	G	144	SER	2.7
10	H	114	ASP	2.7
33	6	11	PRO	2.7
9	G	164	VAL	2.7
11	I	20	LEU	2.7
11	I	17	GLY	2.6
9	G	102	PHE	2.6
9	G	127	PRO	2.6

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Mol	Chain	Res	Type	RSRZ
9	G	184	ILE	2.6
33	6	13	ARG	2.6
9	G	198	ALA	2.6
14	L	55	GLN	2.6
6	D	66	GLY	2.6
14	L	81	VAL	2.6
26	X	80	GLU	2.6
14	L	78	ALA	2.6
11	I	16	PRO	2.6
11	I	19	PRO	2.6
23	U	1	PRO	2.6
33	6	16	PRO	2.6
9	G	71	LEU	2.6
9	G	182	LEU	2.6
11	I	117	LEU	2.6
9	G	119	VAL	2.6
9	G	178	VAL	2.6
27	Y	236	VAL	2.6
6	D	116	SER	2.6
24	V	1	THR	2.6
9	G	111	PRO	2.6
33	6	15	GLY	2.6
9	G	153	GLU	2.6
9	G	36	VAL	2.6
9	G	183	GLY	2.5
6	D	118	GLU	2.5
11	I	92	PHE	2.5
26	X	79	GLU	2.5
9	G	86	THR	2.5
11	I	129	THR	2.5
11	I	60	GLY	2.5
24	V	43	PRO	2.5
11	I	113	LEU	2.5
11	I	109	HIS	2.5
9	G	195	ALA	2.5
6	D	169	THR	2.5
9	G	167	THR	2.5
11	I	61	SER	2.5
11	I	71	THR	2.5
11	I	49	GLU	2.5
33	6	14	ASP	2.5
1	0	971	G	2.5

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Mol	Chain	Res	Type	RSRZ
6	D	68	PRO	2.5
9	G	115	GLY	2.5
9	G	132	GLY	2.5
9	G	190	GLY	2.5
13	K	119	GLN	2.5
23	U	2	ARG	2.5
11	I	33	ALA	2.5
6	D	122	SER	2.5
33	6	48	THR	2.5
6	D	124	GLY	2.4
33	6	50	ILE	2.4
11	I	131	LEU	2.4
10	H	103	GLN	2.4
6	D	114	PHE	2.4
18	P	144	ASN	2.4
26	X	82	GLU	2.4
26	X	81	GLY	2.4
14	L	79	ASP	2.4
7	E	174	ARG	2.4
9	G	196	VAL	2.4
11	I	125	VAL	2.4
26	X	65	ASN	2.4
1	0	996	C	2.4
6	D	57	THR	2.4
33	6	3	THR	2.4
3	A	31	LYS	2.4
6	D	173	GLU	2.4
11	I	130	SER	2.4
6	D	28	GLY	2.4
14	L	87	GLY	2.4
11	I	59	ASP	2.4
9	G	78	LEU	2.4
9	G	91	LEU	2.4
9	G	197	PHE	2.4
11	I	79	ALA	2.4
6	D	9	ASP	2.4
9	G	81	LEU	2.4
9	G	37	ASN	2.4
26	X	83	ALA	2.4
6	D	64	ARG	2.3
11	I	64	ILE	2.3
9	G	54	HIS	2.3

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Mol	Chain	Res	Type	RSRZ
27	Y	108	ASP	2.3
9	G	84	TYR	2.3
9	G	112	ALA	2.3
6	D	85	GLN	2.3
9	G	105	LEU	2.3
11	I	52	VAL	2.3
26	X	78	GLU	2.3
1	0	2231	G	2.3
6	D	32	ARG	2.3
6	D	117	GLN	2.3
9	G	154	GLY	2.3
1	0	1130	U	2.3
2	9	23	U	2.3
9	G	107	ALA	2.3
11	I	15	ASN	2.3
6	D	31	GLY	2.3
1	0	1952	U	2.3
6	D	21	VAL	2.3
9	G	35	VAL	2.3
9	G	150	ARG	2.3
6	D	101	THR	2.2
11	I	24	LEU	2.2
11	I	35	VAL	2.2
9	G	94	THR	2.2
11	I	104	ILE	2.2
1	0	998	U	2.2
1	0	2234	U	2.2
1	0	514	G	2.2
11	I	133	VAL	2.2
11	I	96	LEU	2.2
33	6	52	ILE	2.2
9	G	95	ASP	2.2
9	G	148	ASP	2.2
6	D	91	ALA	2.2
6	D	112	THR	2.2
11	I	80	GLY	2.2
28	Z	83	ILE	2.2
10	H	13	ASP	2.2
14	L	99	GLU	2.2
11	I	69	PRO	2.2
9	G	24	VAL	2.1
11	I	54	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
11	I	138	ASN	2.1
9	G	85	ILE	2.1
1	0	997	C	2.1
1	0	2144	C	2.1
6	D	56	ARG	2.1
9	G	170	GLU	2.1
6	D	120	ASP	2.1
1	0	2915	A	2.1
9	G	93	GLY	2.1
33	6	45	LEU	2.1
9	G	206	GLU	2.1
33	6	4	TYR	2.1
11	I	22	PRO	2.1
9	G	173	GLN	2.1
11	I	6	GLU	2.1
11	I	136	GLU	2.1
3	A	64	ASP	2.1
6	D	165	PHE	2.1
14	L	100	ALA	2.1
6	D	34	LEU	2.1
19	Q	95	GLU	2.1
1	0	2233	C	2.1
30	2	37	HIS	2.1
6	D	58	VAL	2.0
9	G	181	GLU	2.0
11	I	12	GLY	2.0
33	6	21	LYS	2.0
6	D	171	ASP	2.0
7	E	87	PHE	2.0
9	G	33	VAL	2.0
33	6	6	VAL	2.0
2	9	2	U	2.0
9	G	52	ASP	2.0
9	G	199	ASP	2.0
11	I	88	PRO	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. 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($\text{\AA}^2$)	Q<0.9
1	OMU	0	2587	21/22	0.98	0.05	15,18,20,22	0
1	UR3	0	2619	21/22	0.98	0.05	16,18,21,26	0
1	OMG	0	2588	24/25	0.99	0.04	15,17,20,22	0
1	1MA	0	628	23/24	0.99	0.04	13,15,17,18	0
1	PSU	0	2621	20/21	0.99	0.05	14,17,20,21	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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($\text{\AA}^2$)	Q<0.9
36	NA	0	8178	1/1	0.60	0.24	60,60,60,60	0
36	NA	R	202	1/1	0.73	0.15	48,48,48,48	0
34	MG	A	301	1/1	0.82	0.09	42,42,42,42	0
36	NA	0	8134	1/1	0.82	0.13	46,46,46,46	0
34	MG	0	8050	1/1	0.82	0.15	73,73,73,73	0
34	MG	0	8203	1/1	0.82	0.12	43,43,43,43	0
36	NA	0	8168	1/1	0.83	0.24	43,43,43,43	0
34	MG	0	8047	1/1	0.85	0.12	49,49,49,49	0
34	MG	0	8201	1/1	0.85	0.10	42,42,42,42	0
38	ACY	0	8214	4/4	0.85	0.22	28,35,37,39	0
36	NA	9	202	1/1	0.86	0.08	42,42,42,42	0
35	K	O	204	1/1	0.86	0.07	77,77,77,77	0
34	MG	0	8205	1/1	0.86	0.15	43,43,43,43	0
36	NA	0	8144	1/1	0.87	0.13	37,37,37,37	0
37	CL	O	203	1/1	0.88	0.12	56,56,56,56	0
36	NA	0	8179	1/1	0.88	0.10	43,43,43,43	0
36	NA	0	8174	1/1	0.89	0.08	28,28,28,28	0
36	NA	0	8176	1/1	0.89	0.11	37,37,37,37	0
38	ACY	W	201	4/4	0.89	0.12	46,47,48,49	0
34	MG	0	8094	1/1	0.90	0.07	32,32,32,32	0
36	NA	0	8166	1/1	0.90	0.08	38,38,38,38	0
38	ACY	0	8213	4/4	0.90	0.13	34,39,40,49	0
36	NA	0	8118	1/1	0.90	0.09	41,41,41,41	0
34	MG	0	8198	1/1	0.90	0.06	39,39,39,39	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
37	CL	0	8209	1/1	0.91	0.09	54,54,54,54	0
37	CL	G	301	1/1	0.91	0.08	58,58,58,58	0
36	NA	0	8171	1/1	0.91	0.08	39,39,39,39	0
36	NA	0	8172	1/1	0.91	0.12	33,33,33,33	0
34	MG	0	8208	1/1	0.91	0.06	45,45,45,45	0
34	MG	0	8098	1/1	0.91	0.05	33,33,33,33	0
36	NA	0	8180	1/1	0.92	0.14	37,37,37,37	0
36	NA	0	8110	1/1	0.92	0.09	26,26,26,26	0
36	NA	R	201	1/1	0.92	0.06	29,29,29,29	0
34	MG	0	8195	1/1	0.92	0.07	45,45,45,45	0
36	NA	0	8129	1/1	0.92	0.14	37,37,37,37	0
37	CL	0	8210	1/1	0.92	0.09	49,49,49,49	0
37	CL	9	203	1/1	0.92	0.11	59,59,59,59	0
36	NA	0	8132	1/1	0.92	0.08	23,23,23,23	0
34	MG	N	202	1/1	0.92	0.17	58,58,58,58	0
34	MG	0	8095	1/1	0.92	0.08	34,34,34,34	0
36	NA	0	8149	1/1	0.92	0.09	28,28,28,28	0
36	NA	0	8156	1/1	0.92	0.35	33,33,33,33	0
36	NA	0	8145	1/1	0.93	0.10	31,31,31,31	0
34	MG	0	8086	1/1	0.93	0.08	38,38,38,38	0
34	MG	0	8199	1/1	0.93	0.14	55,55,55,55	0
36	NA	0	8161	1/1	0.93	0.08	29,29,29,29	0
34	MG	9	201	1/1	0.93	0.10	54,54,54,54	0
38	ACY	0	8212	4/4	0.93	0.16	35,39,41,43	0
34	MG	0	8200	1/1	0.93	0.06	37,37,37,37	0
34	MG	0	8091	1/1	0.93	0.06	20,20,20,20	0
34	MG	0	8197	1/1	0.93	0.10	31,31,31,31	0
34	MG	0	8097	1/1	0.94	0.06	34,34,34,34	0
36	NA	0	8175	1/1	0.94	0.11	30,30,30,30	0
34	MG	0	8076	1/1	0.94	0.10	38,38,38,38	0
34	MG	0	8104	1/1	0.94	0.07	26,26,26,26	0
34	MG	0	8107	1/1	0.94	0.04	32,32,32,32	0
34	MG	Y	304	1/1	0.94	0.05	33,33,33,33	0
36	NA	0	8154	1/1	0.94	0.10	24,24,24,24	0
36	NA	H	202	1/1	0.94	0.09	39,39,39,39	0
34	MG	0	8202	1/1	0.94	0.07	33,33,33,33	0
36	NA	0	8158	1/1	0.94	0.07	28,28,28,28	0
36	NA	0	8160	1/1	0.94	0.11	26,26,26,26	0
34	MG	0	8087	1/1	0.94	0.05	22,22,22,22	0
36	NA	0	8164	1/1	0.94	0.07	39,39,39,39	0
36	NA	0	8115	1/1	0.94	0.08	32,32,32,32	0
36	NA	0	8167	1/1	0.94	0.12	29,29,29,29	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
34	MG	0	8088	1/1	0.94	0.06	40,40,40,40	0
36	NA	0	8170	1/1	0.94	0.14	38,38,38,38	0
34	MG	0	8207	1/1	0.94	0.06	33,33,33,33	0
36	NA	0	8130	1/1	0.94	0.12	31,31,31,31	0
36	NA	0	8162	1/1	0.95	0.05	25,25,25,25	0
34	MG	0	8093	1/1	0.95	0.09	42,42,42,42	0
34	MG	0	8061	1/1	0.95	0.05	30,30,30,30	0
34	MG	0	8046	1/1	0.95	0.09	37,37,37,37	0
35	K	0	8211	1/1	0.95	0.13	51,51,51,51	0
36	NA	0	8169	1/1	0.95	0.05	30,30,30,30	0
34	MG	0	8204	1/1	0.95	0.05	33,33,33,33	0
36	NA	0	8146	1/1	0.95	0.08	22,22,22,22	0
34	MG	0	8096	1/1	0.95	0.14	43,43,43,43	0
34	MG	0	8079	1/1	0.95	0.04	27,27,27,27	0
34	MG	0	8084	1/1	0.95	0.07	37,37,37,37	0
36	NA	0	8119	1/1	0.95	0.08	32,32,32,32	0
36	NA	0	8124	1/1	0.95	0.08	32,32,32,32	0
34	MG	0	8100	1/1	0.95	0.08	42,42,42,42	0
36	NA	0	8147	1/1	0.96	0.17	23,23,23,23	0
34	MG	L	204	1/1	0.96	0.12	26,26,26,26	0
36	NA	L	201	1/1	0.96	0.14	24,24,24,24	0
36	NA	M	201	1/1	0.96	0.08	15,15,15,15	0
34	MG	0	8040	1/1	0.96	0.04	26,26,26,26	0
36	NA	0	8131	1/1	0.96	0.06	28,28,28,28	0
36	NA	S	101	1/1	0.96	0.13	25,25,25,25	0
36	NA	0	8157	1/1	0.96	0.12	24,24,24,24	0
34	MG	Y	303	1/1	0.96	0.10	38,38,38,38	0
34	MG	0	8196	1/1	0.96	0.04	39,39,39,39	0
37	CL	B	402	1/1	0.96	0.26	62,62,62,62	0
36	NA	0	8135	1/1	0.96	0.05	29,29,29,29	0
37	CL	J	203	1/1	0.96	0.06	42,42,42,42	0
37	CL	J	204	1/1	0.96	0.10	36,36,36,36	0
35	K	0	8109	1/1	0.96	0.15	44,44,44,44	0
36	NA	0	8177	1/1	0.96	0.05	33,33,33,33	0
34	MG	0	8063	1/1	0.96	0.05	19,19,19,19	0
36	NA	0	8165	1/1	0.96	0.06	29,29,29,29	0
36	NA	0	8125	1/1	0.96	0.06	19,19,19,19	0
34	MG	0	8191	1/1	0.97	0.08	31,31,31,31	0
34	MG	0	8042	1/1	0.97	0.03	28,28,28,28	0
34	MG	0	8085	1/1	0.97	0.06	28,28,28,28	0
36	NA	0	8111	1/1	0.97	0.07	25,25,25,25	0
34	MG	0	8062	1/1	0.97	0.07	35,35,35,35	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
36	NA	0	8117	1/1	0.97	0.04	24,24,24,24	0
34	MG	0	8041	1/1	0.97	0.09	25,25,25,25	0
34	MG	0	8064	1/1	0.97	0.11	20,20,20,20	0
36	NA	0	8120	1/1	0.97	0.03	21,21,21,21	0
36	NA	0	8121	1/1	0.97	0.05	20,20,20,20	0
34	MG	0	8089	1/1	0.97	0.06	23,23,23,23	0
34	MG	0	8066	1/1	0.97	0.02	27,27,27,27	0
36	NA	0	8126	1/1	0.97	0.07	23,23,23,23	0
36	NA	0	8128	1/1	0.97	0.11	19,19,19,19	0
34	MG	0	8092	1/1	0.97	0.04	25,25,25,25	0
34	MG	0	8067	1/1	0.97	0.04	33,33,33,33	0
34	MG	0	8068	1/1	0.97	0.04	34,34,34,34	0
36	NA	A	304	1/1	0.97	0.08	28,28,28,28	0
36	NA	H	201	1/1	0.97	0.08	19,19,19,19	0
34	MG	0	8070	1/1	0.97	0.05	22,22,22,22	0
34	MG	0	8206	1/1	0.97	0.05	33,33,33,33	0
36	NA	L	202	1/1	0.97	0.15	24,24,24,24	0
34	MG	0	8072	1/1	0.97	0.03	15,15,15,15	0
36	NA	0	8141	1/1	0.97	0.04	30,30,30,30	0
34	MG	0	8075	1/1	0.97	0.05	28,28,28,28	0
34	MG	0	8051	1/1	0.97	0.08	32,32,32,32	0
37	CL	0	8189	1/1	0.97	0.04	35,35,35,35	0
34	MG	0	8053	1/1	0.97	0.12	22,22,22,22	0
34	MG	A	303	1/1	0.97	0.08	23,23,23,23	0
34	MG	0	8103	1/1	0.97	0.03	15,15,15,15	0
36	NA	0	8150	1/1	0.97	0.11	36,36,36,36	0
37	CL	F	201	1/1	0.97	0.09	33,33,33,33	0
34	MG	0	8081	1/1	0.97	0.11	42,42,42,42	0
34	MG	T	201	1/1	0.97	0.11	47,47,47,47	0
34	MG	Y	301	1/1	0.97	0.04	19,19,19,19	0
34	MG	0	8105	1/1	0.97	0.03	23,23,23,23	0
37	CL	Y	302	1/1	0.97	0.04	26,26,26,26	0
36	NA	0	8159	1/1	0.97	0.06	26,26,26,26	0
34	MG	0	8106	1/1	0.97	0.06	29,29,29,29	0
34	MG	Y	305	1/1	0.97	0.04	33,33,33,33	0
34	MG	0	8082	1/1	0.97	0.03	17,17,17,17	0
34	MG	0	8108	1/1	0.98	0.03	20,20,20,20	0
34	MG	0	8014	1/1	0.98	0.02	19,19,19,19	0
34	MG	0	8194	1/1	0.98	0.06	35,35,35,35	0
36	NA	0	8173	1/1	0.98	0.18	25,25,25,25	0
34	MG	0	8043	1/1	0.98	0.02	29,29,29,29	0
34	MG	0	8065	1/1	0.98	0.02	40,40,40,40	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
36	NA	0	8137	1/1	0.98	0.07	21,21,21,21	0
36	NA	0	8138	1/1	0.98	0.10	20,20,20,20	0
36	NA	0	8139	1/1	0.98	0.12	29,29,29,29	0
36	NA	0	8140	1/1	0.98	0.07	21,21,21,21	0
34	MG	0	8044	1/1	0.98	0.03	21,21,21,21	0
36	NA	0	8192	1/1	0.98	0.05	22,22,22,22	0
36	NA	0	8143	1/1	0.98	0.03	13,13,13,13	0
34	MG	0	8090	1/1	0.98	0.06	19,19,19,19	0
34	MG	0	8045	1/1	0.98	0.03	35,35,35,35	0
34	MG	0	8024	1/1	0.98	0.07	17,17,17,17	0
34	MG	0	8032	1/1	0.98	0.02	17,17,17,17	0
36	NA	0	8148	1/1	0.98	0.03	21,21,21,21	0
34	MG	0	8071	1/1	0.98	0.03	37,37,37,37	0
36	NA	Q	101	1/1	0.98	0.03	25,25,25,25	0
36	NA	0	8114	1/1	0.98	0.14	19,19,19,19	0
36	NA	0	8153	1/1	0.98	0.15	29,29,29,29	0
34	MG	0	8049	1/1	0.98	0.04	36,36,36,36	0
37	CL	0	8181	1/1	0.98	0.04	29,29,29,29	0
37	CL	0	8185	1/1	0.98	0.06	32,32,32,32	0
37	CL	0	8187	1/1	0.98	0.13	38,38,38,38	0
37	CL	0	8188	1/1	0.98	0.10	30,30,30,30	0
36	NA	0	8155	1/1	0.98	0.04	34,34,34,34	0
34	MG	0	8074	1/1	0.98	0.04	30,30,30,30	0
34	MG	0	8035	1/1	0.98	0.03	27,27,27,27	0
34	MG	0	8036	1/1	0.98	0.04	21,21,21,21	0
37	CL	A	305	1/1	0.98	0.08	35,35,35,35	0
34	MG	0	8052	1/1	0.98	0.04	36,36,36,36	0
34	MG	0	8037	1/1	0.98	0.03	29,29,29,29	0
36	NA	0	8122	1/1	0.98	0.05	16,16,16,16	0
37	CL	J	202	1/1	0.98	0.04	34,34,34,34	0
36	NA	0	8123	1/1	0.98	0.06	23,23,23,23	0
36	NA	0	8163	1/1	0.98	0.15	18,18,18,18	0
37	CL	L	203	1/1	0.98	0.08	29,29,29,29	0
37	CL	N	201	1/1	0.98	0.05	36,36,36,36	0
37	CL	O	202	1/1	0.98	0.05	36,36,36,36	0
34	MG	0	8056	1/1	0.98	0.03	25,25,25,25	0
37	CL	R	203	1/1	0.98	0.09	30,30,30,30	0
34	MG	0	8083	1/1	0.98	0.05	34,34,34,34	0
34	MG	0	8006	1/1	0.98	0.03	22,22,22,22	0
36	NA	0	8127	1/1	0.98	0.20	22,22,22,22	0
34	MG	K	201	1/1	0.98	0.03	25,25,25,25	0
34	MG	0	8012	1/1	0.98	0.05	23,23,23,23	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
39	CD	O	201	1/1	0.98	0.07	58,58,58,58	0
36	NA	0	8142	1/1	0.99	0.03	28,28,28,28	0
34	MG	3	101	1/1	0.99	0.03	27,27,27,27	0
34	MG	0	8078	1/1	0.99	0.02	27,27,27,27	0
36	NA	C	301	1/1	0.99	0.02	20,20,20,20	0
34	MG	0	8031	1/1	0.99	0.03	15,15,15,15	0
34	MG	0	8193	1/1	0.99	0.13	14,14,14,14	0
36	NA	J	201	1/1	0.99	0.02	27,27,27,27	0
34	MG	0	8003	1/1	0.99	0.02	12,12,12,12	0
34	MG	0	8054	1/1	0.99	0.04	10,10,10,10	0
36	NA	0	8112	1/1	0.99	0.07	19,19,19,19	0
36	NA	0	8113	1/1	0.99	0.06	19,19,19,19	0
36	NA	0	8152	1/1	0.99	0.03	17,17,17,17	0
34	MG	0	8013	1/1	0.99	0.03	16,16,16,16	0
34	MG	0	8057	1/1	0.99	0.02	23,23,23,23	0
36	NA	0	8116	1/1	0.99	0.03	29,29,29,29	0
37	CL	0	8182	1/1	0.99	0.09	26,26,26,26	0
37	CL	0	8183	1/1	0.99	0.05	25,25,25,25	0
37	CL	0	8184	1/1	0.99	0.06	26,26,26,26	0
34	MG	0	8058	1/1	0.99	0.02	18,18,18,18	0
37	CL	0	8186	1/1	0.99	0.07	25,25,25,25	0
34	MG	0	8059	1/1	0.99	0.06	21,21,21,21	0
34	MG	0	8060	1/1	0.99	0.03	22,22,22,22	0
34	MG	0	8001	1/1	0.99	0.03	16,16,16,16	0
37	CL	0	8190	1/1	0.99	0.26	31,31,31,31	0
34	MG	0	8015	1/1	0.99	0.04	15,15,15,15	0
34	MG	0	8039	1/1	0.99	0.02	27,27,27,27	0
34	MG	0	8017	1/1	0.99	0.05	11,11,11,11	0
34	MG	0	8018	1/1	0.99	0.03	22,22,22,22	0
37	CL	B	401	1/1	0.99	0.06	24,24,24,24	0
34	MG	0	8019	1/1	0.99	0.05	23,23,23,23	0
34	MG	0	8022	1/1	0.99	0.02	23,23,23,23	0
34	MG	0	8023	1/1	0.99	0.05	20,20,20,20	0
34	MG	0	8069	1/1	0.99	0.05	29,29,29,29	0
34	MG	0	8007	1/1	0.99	0.06	18,18,18,18	0
34	MG	A	302	1/1	0.99	0.02	20,20,20,20	0
34	MG	0	8025	1/1	0.99	0.02	25,25,25,25	0
37	CL	M	202	1/1	0.99	0.04	21,21,21,21	0
34	MG	0	8026	1/1	0.99	0.04	19,19,19,19	0
36	NA	0	8133	1/1	0.99	0.11	28,28,28,28	0
34	MG	0	8101	1/1	0.99	0.04	19,19,19,19	0
34	MG	0	8102	1/1	0.99	0.01	24,24,24,24	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
37	CL	R	204	1/1	0.99	0.05	34,34,34,34	0
36	NA	0	8136	1/1	0.99	0.05	20,20,20,20	0
37	CL	3	103	1/1	0.99	0.05	33,33,33,33	0
34	MG	0	8073	1/1	0.99	0.04	10,10,10,10	0
34	MG	0	8027	1/1	0.99	0.02	30,30,30,30	0
34	MG	0	8028	1/1	0.99	0.04	21,21,21,21	0
34	MG	0	8029	1/1	0.99	0.02	22,22,22,22	0
34	MG	0	8077	1/1	0.99	0.03	24,24,24,24	0
39	CD	U	8401	1/1	0.99	0.02	39,39,39,39	0
39	CD	Z	101	1/1	0.99	0.04	38,38,38,38	0
39	CD	3	102	1/1	0.99	0.03	39,39,39,39	0
34	MG	0	8048	1/1	1.00	0.03	33,33,33,33	0
34	MG	0	8099	1/1	1.00	0.04	26,26,26,26	0
34	MG	0	8034	1/1	1.00	0.04	25,25,25,25	0
34	MG	0	8005	1/1	1.00	0.05	13,13,13,13	0
34	MG	0	8002	1/1	1.00	0.01	19,19,19,19	0
34	MG	0	8004	1/1	1.00	0.03	13,13,13,13	0
34	MG	0	8038	1/1	1.00	0.02	13,13,13,13	0
34	MG	0	8008	1/1	1.00	0.03	13,13,13,13	0
34	MG	0	8055	1/1	1.00	0.02	31,31,31,31	0
34	MG	0	8016	1/1	1.00	0.04	20,20,20,20	0
34	MG	0	8009	1/1	1.00	0.02	15,15,15,15	0
34	MG	0	8010	1/1	1.00	0.02	14,14,14,14	0
34	MG	0	8011	1/1	1.00	0.05	16,16,16,16	0
36	NA	0	8151	1/1	1.00	0.01	11,11,11,11	0
34	MG	0	8030	1/1	1.00	0.02	17,17,17,17	0
34	MG	0	8020	1/1	1.00	0.03	17,17,17,17	0
34	MG	0	8021	1/1	1.00	0.02	17,17,17,17	0
34	MG	0	8033	1/1	1.00	0.03	16,16,16,16	0
39	CD	1	101	1/1	1.00	0.04	37,37,37,37	0
34	MG	0	8080	1/1	1.00	0.02	31,31,31,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.