



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 10:33 pm BST

PDB ID : 6MKN
Title : Structure of the *Thermus thermophilus* 30S ribosomal subunit complexed with an inosine (I34) modified anticodon stem loop (ASL) of *Escherichia coli* transfer RNA Arginine 2 (TRNAARG2) bound to an mRNA with an CGU-codon in the A-site and paromomycin
Authors : Cantara, W.A.; DeMirci, H.; Agris, P.F.
Deposited on : 2018-09-25
Resolution : 3.46 Å (reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

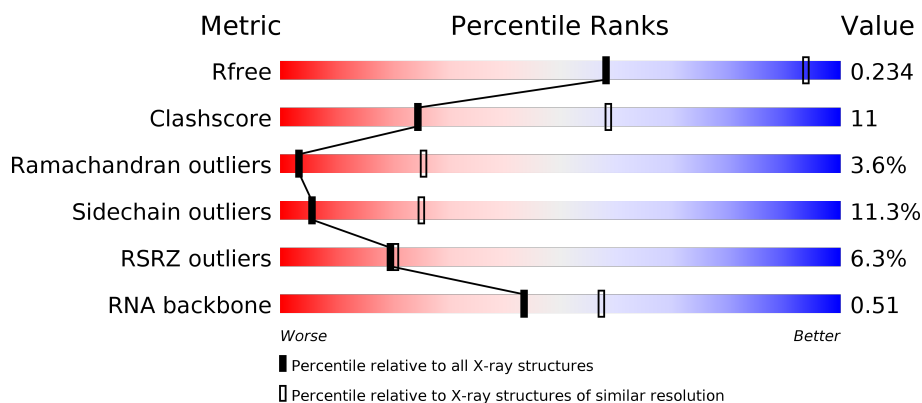
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.46 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1291 (3.52-3.40)
Clashscore	141614	1372 (3.52-3.40)
Ramachandran outliers	138981	1337 (3.52-3.40)
Sidechain outliers	138945	1338 (3.52-3.40)
RSRZ outliers	127900	1205 (3.52-3.40)
RNA backbone	3102	1036 (3.96-2.96)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	1507	<div> <div>2%</div> <div> <div></div> <div>49%</div> <div>37%</div> <div>11%</div> <div>.</div> </div> </div>
2	B	256	<div> <div>7%</div> <div> <div></div> <div>55%</div> <div>30%</div> <div>6%</div> <div>9%</div> </div> </div>
3	C	239	<div> <div>7%</div> <div> <div></div> <div>52%</div> <div>27%</div> <div>6%</div> <div>.</div> <div>14%</div> </div> </div>
4	D	209	<div> <div>6%</div> <div> <div></div> <div>66%</div> <div>29%</div> <div>5%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
5	E	162	
6	F	101	
7	G	156	
8	H	138	
9	I	128	
10	J	105	
11	K	129	
12	L	132	
13	M	126	
14	N	61	
15	O	89	
16	P	88	
17	Q	105	
18	R	88	
19	S	93	
20	T	106	
21	V	27	
22	X	17	
23	W	6	

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
25	MG	A	1602	-	-	-	X
25	MG	A	1604	-	-	-	X
25	MG	A	1606	-	-	-	X
25	MG	A	1609	-	-	-	X
25	MG	A	1614	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
25	MG	A	1619	-	-	-	X
25	MG	A	1624	-	-	-	X
25	MG	A	1629	-	-	-	X
25	MG	A	1632	-	-	-	X
25	MG	A	1643	-	-	-	X
25	MG	A	1644	-	-	-	X
25	MG	A	1662	-	-	-	X
25	MG	A	1663	-	-	-	X
25	MG	A	1664	-	-	-	X
25	MG	A	1667	-	-	-	X
25	MG	A	1671	-	-	-	X
25	MG	A	1675	-	-	-	X
25	MG	E	201	-	-	-	X

2 Entry composition

There are 25 unique types of molecules in this entry. The entry contains 52043 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 16S rRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1507	Total	C	N	O	P	0	0	0
			32380	14414	5990	10470	1506			

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	G	deletion	GB 55771382
A	?	-	G	deletion	GB 55771382
A	?	-	C	deletion	GB 55771382
A	?	-	G	deletion	GB 55771382
A	?	-	A	deletion	GB 55771382
A	?	-	C	deletion	GB 55771382
A	1540	U	-	insertion	GB 55771382
A	1541	U	-	insertion	GB 55771382
A	1542	U	-	insertion	GB 55771382
A	1543	C	-	insertion	GB 55771382
A	1544	U	-	insertion	GB 55771382

- Molecule 2 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	234	Total	C	N	O	S	0	0	0
			1900	1213	341	341	5			

- Molecule 3 is a protein called 30S ribosomal protein S3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	206	Total	C	N	O	S	0	0	0
			1612	1016	314	281	1			

- Molecule 4 is a protein called 30S ribosomal protein S4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	208	Total	C	N	O	S	0	0	0
			1703	1066	339	291	7			

- Molecule 5 is a protein called 30S ribosomal protein S5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	150	Total	C	N	O	S	0	0	0
			1146	724	217	201	4			

- Molecule 6 is a protein called 30S ribosomal protein S6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	101	Total	C	N	O	S	0	0	0
			843	531	155	154	3			

- Molecule 7 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	155	Total	C	N	O	S	0	0	0
			1257	781	252	218	6			

- Molecule 8 is a protein called 30S ribosomal protein S8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	138	Total	C	N	O	S	0	0	0
			1116	705	215	193	3			

- Molecule 9 is a protein called 30S ribosomal protein S9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	I	127	Total	C	N	O		0	0	0
			1011	639	198	174				

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I	58	ARG	HIS	conflict	UNP P80374

- Molecule 10 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	J	98	Total	C	N	O	S	0	0	0
			792	498	156	137	1			

- Molecule 11 is a protein called 30S ribosomal protein S11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	K	119	Total	C	N	O	S	0	0	0
			885	549	168	165	3			

- Molecule 12 is a protein called 30S ribosomal protein S12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	L	124	Total	C	N	O	S	0	0	0
			970	611	195	163	1			

- Molecule 13 is a protein called 30S ribosomal protein S13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	M	125	Total	C	N	O	S	0	0	0
			997	617	207	171	2			

- Molecule 14 is a protein called 30S ribosomal protein S14 type Z.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	N	60	Total	C	N	O	S	0	0	0
			492	312	104	72	4			

- Molecule 15 is a protein called 30S ribosomal protein S15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
15	O	88	Total	C	N	O	S	0	0	0
			734	459	147	126	2			

- Molecule 16 is a protein called 30S ribosomal protein S16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
16	P	83	Total	C	N	O	S	0	0	0
			700	443	139	117	1			

- Molecule 17 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
17	Q	104	Total	C	N	O	S	0	0	0
			857	547	161	147	2			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Q	96	GLN	GLU	conflict	UNP P0DOY7

- Molecule 18 is a protein called 30S ribosomal protein S18.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
18	R	73	Total	C	N	O		0	0	0
			597	380	118	99				

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
R	60	GLY	ALA	conflict	UNP Q5SLQ0

- Molecule 19 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
19	S	80	Total	C	N	O	S	0	0	0
			647	414	119	112	2			

- Molecule 20 is a protein called 30S ribosomal protein S20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
20	T	99	Total	C	N	O	S	0	0	0
			762	469	162	129	2			

- Molecule 21 is a protein called 30S ribosomal protein Thx.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
21	V	24	Total	C	N	O	0	0	0
			208	128	50	30			

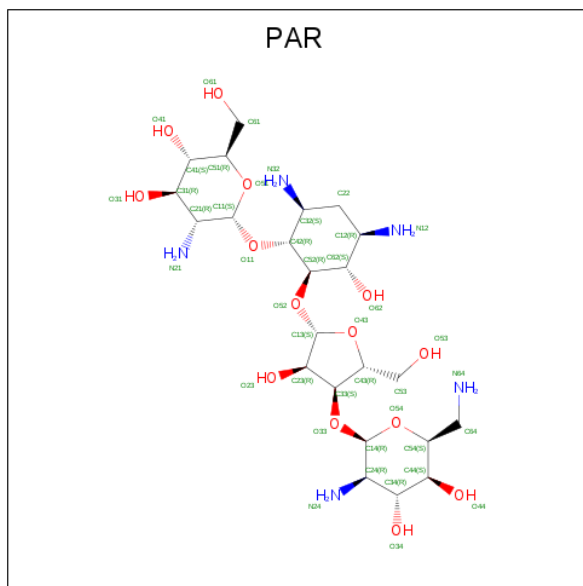
- Molecule 22 is a RNA chain called tRNA ASL Escherichia coli Arg2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
22	X	11	Total	C	N	O	P	0	0	0
			232	105	43	74	10			

- Molecule 23 is a RNA chain called mRNA A-site fragment.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
23	W	4	Total	C	N	O	P	0	0
			82	38	15	26	3		

- Molecule 24 is PAROMOMYCIN (three-letter code: PAR) (formula: $C_{23}H_{45}N_5O_{14}$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
24	A	1	Total	C	N	O	0	0
			42	23	5	14		

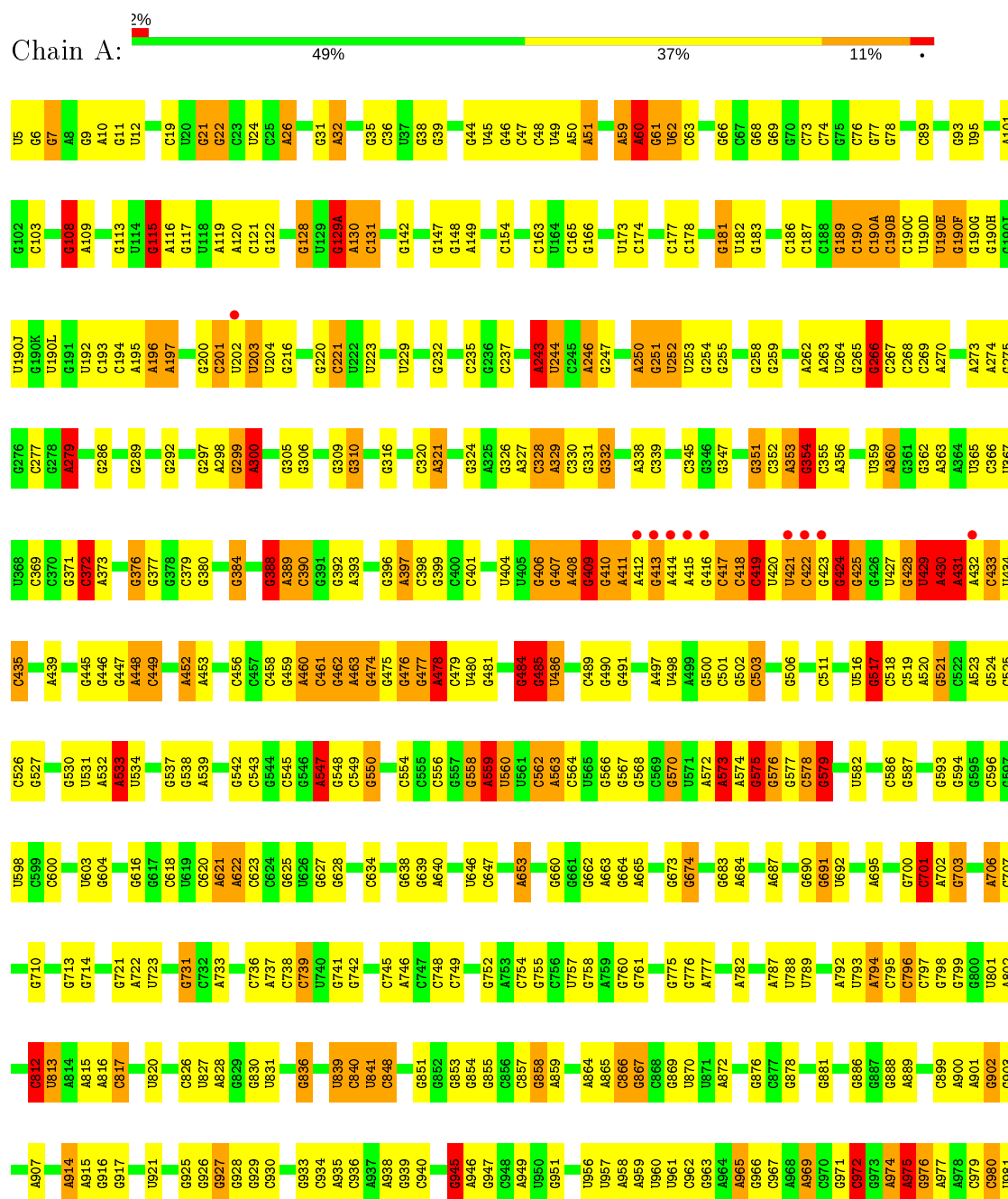
- Molecule 25 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

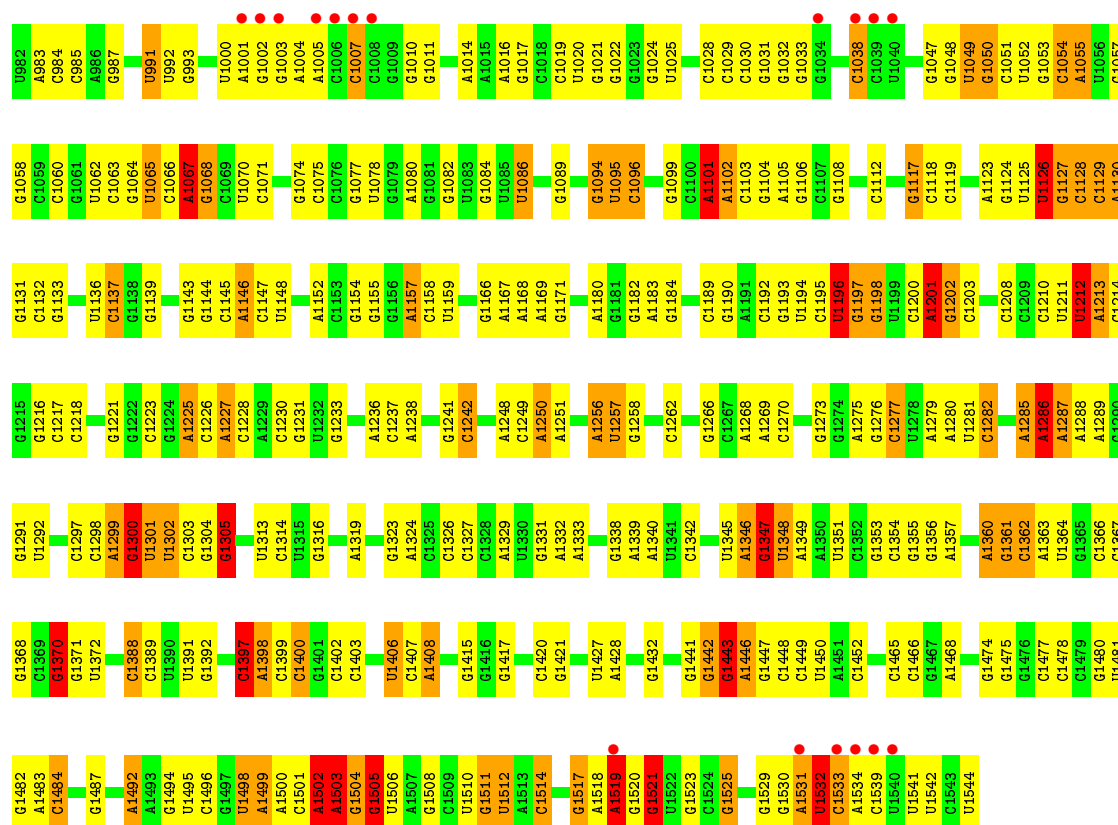
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
25	J	1	Total	Mg	0	0
			1	1		
25	A	76	Total	Mg	0	0
			76	76		
25	E	1	Total	Mg	0	0
			1	1		

3 Residue-property plots

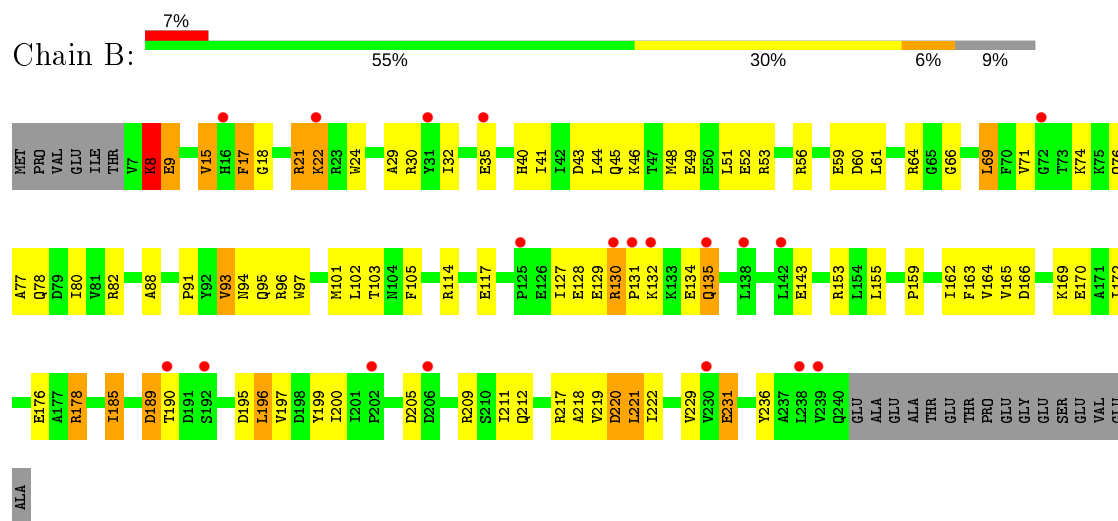
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: 16S rRNA

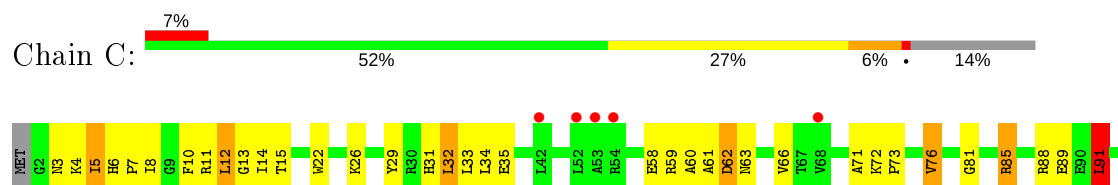


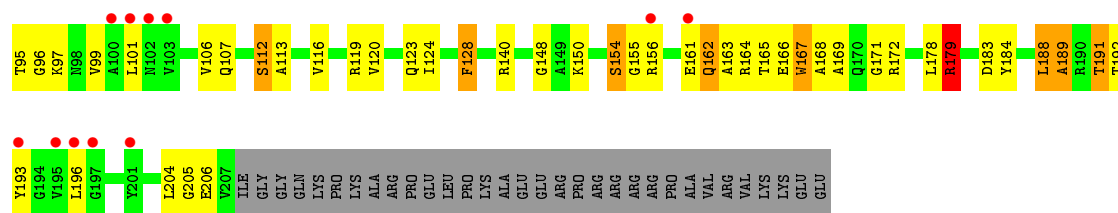


• Molecule 2: 30S ribosomal protein S2

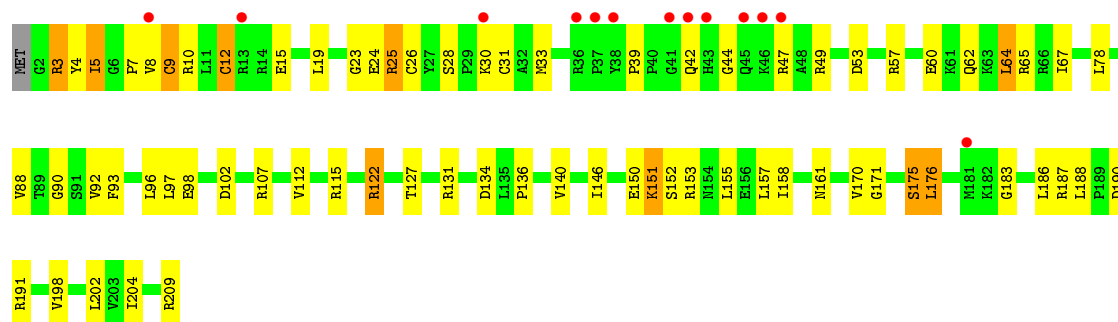


• Molecule 3: 30S ribosomal protein S3

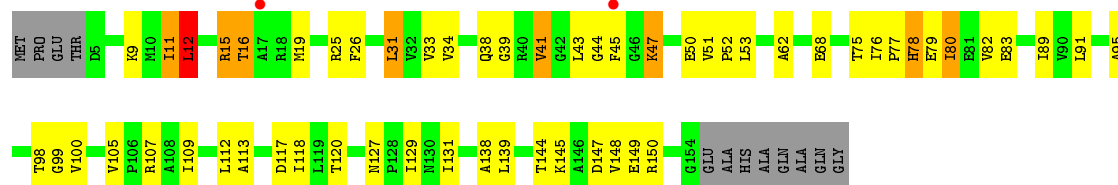




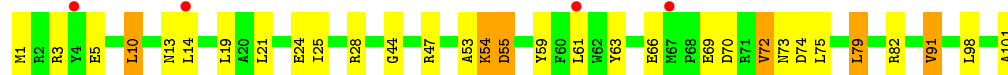
• Molecule 4: 30S ribosomal protein S4



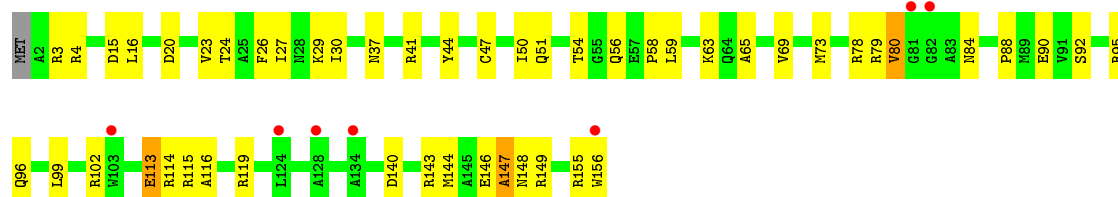
• Molecule 5: 30S ribosomal protein S5



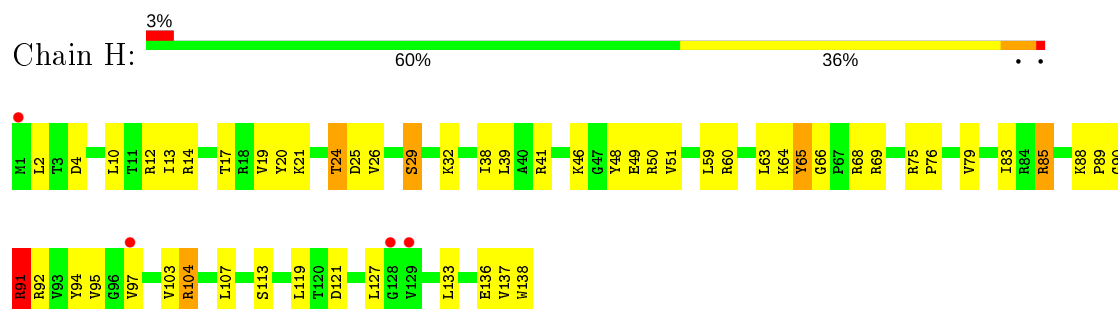
• Molecule 6: 30S ribosomal protein S6



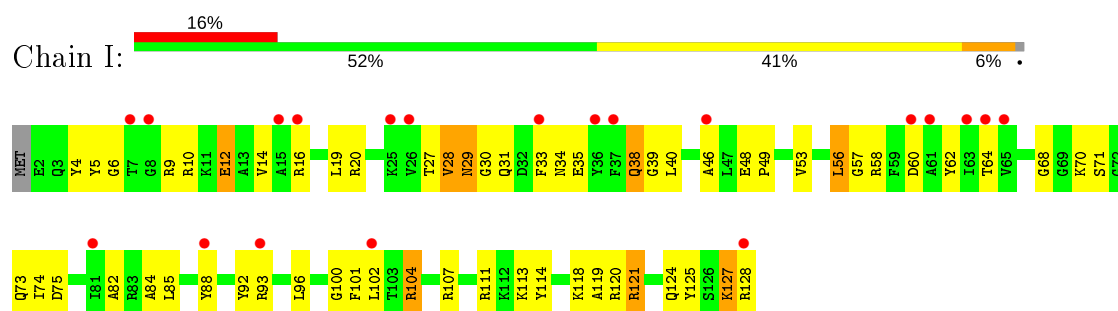
• Molecule 7: 30S ribosomal protein S7



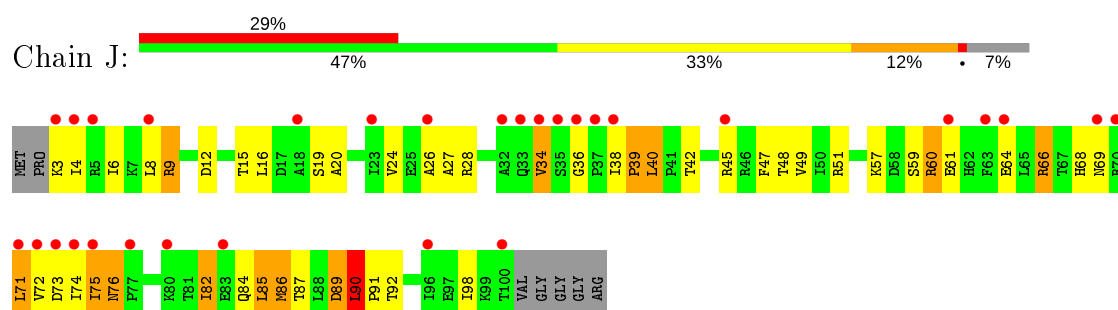
- Molecule 8: 30S ribosomal protein S8



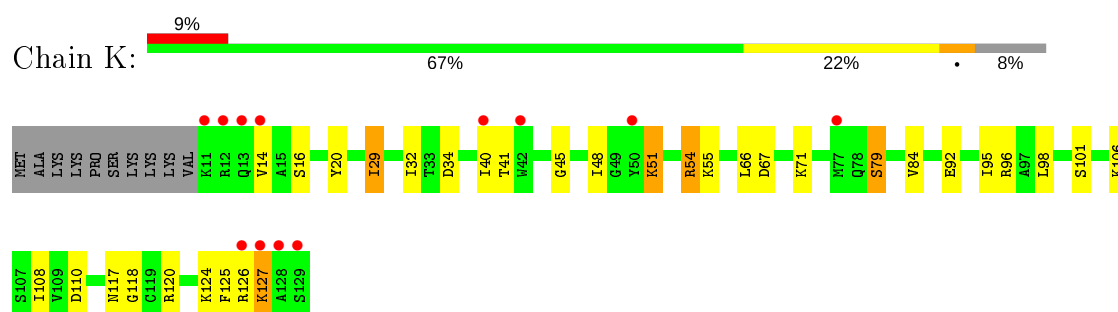
- Molecule 9: 30S ribosomal protein S9



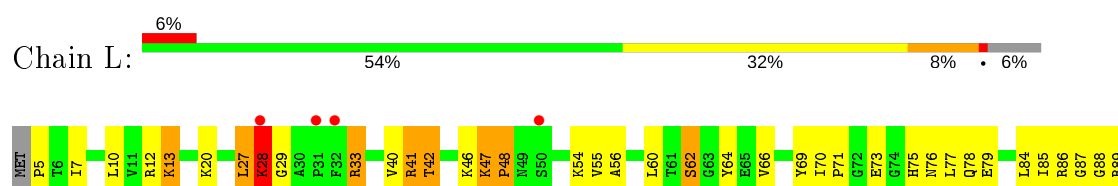
- Molecule 10: 30S ribosomal protein S10



- Molecule 11: 30S ribosomal protein S11



- Molecule 12: 30S ribosomal protein S12

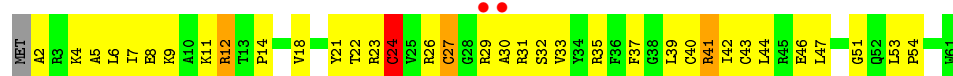




- Molecule 13: 30S ribosomal protein S13



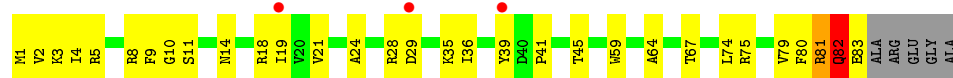
- Molecule 14: 30S ribosomal protein S14 type Z



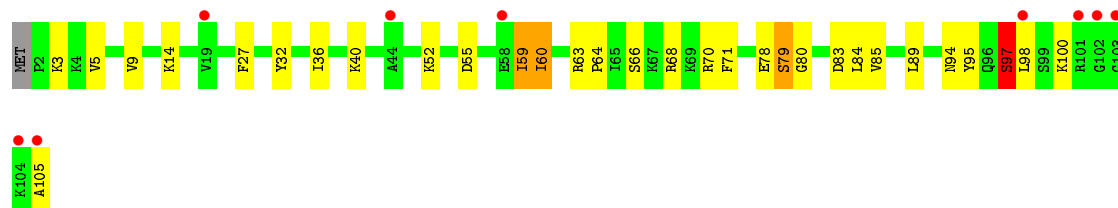
- Molecule 15: 30S ribosomal protein S15



- Molecule 16: 30S ribosomal protein S16



- Molecule 17: 30S ribosomal protein S17



- Molecule 18: 30S ribosomal protein S18



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	403.73 Å 403.73 Å 177.42 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.80 – 3.46 49.80 – 3.46	Depositor EDS
% Data completeness (in resolution range)	87.2 (49.80-3.46) 83.4 (49.80-3.46)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.42 (at 3.48 Å)	Xtriage
Refinement program	PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.203 , 0.234 0.203 , 0.234	Depositor DCC
R_{free} test set	1778 reflections (1.13%)	wwPDB-VP
Wilson B-factor (Å ²)	136.3	Xtriage
Anisotropy	0.186	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 99.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	52043	wwPDB-VP
Average B, all atoms (Å ²)	141.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.24% 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: PAR, MG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.80	9/36244 (0.0%)	1.39	409/56567 (0.7%)
2	B	0.45	0/1935	0.58	0/2609
3	C	0.49	1/1636 (0.1%)	0.62	0/2205
4	D	0.51	0/1733	0.62	1/2318 (0.0%)
5	E	0.57	0/1162	0.68	1/1564 (0.1%)
6	F	0.43	0/856	0.56	0/1154
7	G	0.40	0/1276	0.54	0/1709
8	H	0.64	0/1136	0.72	0/1527
9	I	0.47	0/1029	0.62	0/1378
10	J	0.41	0/805	0.62	0/1082
11	K	0.47	0/900	0.62	0/1213
12	L	0.62	0/986	0.72	0/1320
13	M	0.45	0/1008	0.58	0/1347
14	N	0.53	0/501	0.73	1/664 (0.2%)
15	O	0.50	0/745	0.62	0/992
16	P	0.57	0/716	0.70	0/963
17	Q	0.55	0/870	0.67	0/1159
18	R	0.46	0/603	0.59	0/799
19	S	0.41	0/661	0.59	0/890
20	T	0.45	0/764	0.64	0/1006
21	V	0.47	0/212	0.65	0/277
22	X	0.80	0/258	1.37	2/398 (0.5%)
23	W	0.85	0/91	1.45	0/140
All	All	0.71	10/56127 (0.0%)	1.20	414/83281 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
20	T	0	1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	119	A	N9-C4	-7.68	1.33	1.37
1	A	1362	C	P-O5'	6.89	1.66	1.59
1	A	1502	A	N7-C5	-6.63	1.35	1.39
1	A	1361	G	O3'-P	6.31	1.68	1.61
1	A	1502	A	C5-C6	-6.06	1.35	1.41
3	C	167	TRP	CB-CG	-5.78	1.39	1.50
1	A	1361	G	P-O5'	5.51	1.65	1.59
1	A	190	C	N1-C2	5.43	1.45	1.40
1	A	429	U	C2-N3	5.20	1.41	1.37
1	A	867	G	N9-C4	5.17	1.42	1.38

All (414) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1502	A	N1-C6-N6	11.64	125.58	118.60
1	A	1502	A	C6-C5-N7	-11.26	124.42	132.30
1	A	742	G	N1-C6-O6	10.97	126.48	119.90
1	A	1370	G	C5-C6-N1	-10.55	106.22	111.50
1	A	1347	G	C8-N9-C4	10.03	110.41	106.40
1	A	108	G	N1-C6-O6	9.55	125.63	119.90
1	A	60	A	P-O3'-C3'	9.50	131.10	119.70
1	A	848	C	C6-N1-C2	-9.45	116.52	120.30
1	A	190	C	C6-N1-C2	-9.31	116.58	120.30
1	A	573	A	N9-C4-C5	9.28	109.51	105.80
1	A	174	C	C6-N1-C2	-9.17	116.63	120.30
1	A	117	G	N1-C6-O6	9.13	125.38	119.90
1	A	1208	C	C6-N1-C2	-9.10	116.66	120.30
1	A	867	G	N3-C4-N9	9.07	131.44	126.00
1	A	328	C	O5'-P-OP1	-8.92	97.67	105.70
1	A	812	C	C6-N1-C2	-8.84	116.77	120.30
1	A	435	C	C6-N1-C2	-8.76	116.80	120.30
1	A	429	U	C6-N1-C2	-8.70	115.78	121.00
1	A	1511	G	C5-C6-N1	-8.66	107.17	111.50
1	A	575	G	P-O3'-C3'	8.63	130.06	119.70
1	A	1502	A	C4-C5-C6	8.60	121.30	117.00
1	A	299	G	C5-C6-N1	-8.57	107.22	111.50
1	A	867	G	N3-C4-C5	-8.53	124.34	128.60
1	A	867	G	C4-N9-C1'	8.49	137.53	126.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1397	C	C6-N1-C2	-8.46	116.92	120.30
1	A	1101	A	P-O3'-C3'	8.39	129.77	119.70
1	A	328	C	C6-N1-C2	-8.37	116.95	120.30
1	A	742	G	C5-C6-O6	-8.24	123.66	128.60
1	A	181	G	C8-N9-C4	-8.20	103.12	106.40
1	A	1501	C	N3-C4-C5	8.19	125.17	121.90
1	A	1512	U	C6-N1-C2	-8.13	116.12	121.00
1	A	1502	A	C5-N7-C8	-8.11	99.84	103.90
1	A	1505	G	C8-N9-C4	-8.04	103.19	106.40
1	A	1051	C	C6-N1-C2	-8.03	117.09	120.30
1	A	573	A	C8-N9-C4	-8.00	102.60	105.80
1	A	59	A	N1-C6-N6	-7.99	113.81	118.60
1	A	1498	U	P-O3'-C3'	7.98	129.28	119.70
1	A	5	U	N1-C2-O2	7.98	128.38	122.80
1	A	1103	C	C6-N1-C2	7.89	123.45	120.30
1	A	831	U	C5-C4-O4	7.85	130.61	125.90
1	A	254	G	O5'-P-OP1	-7.80	98.68	105.70
1	A	1196	U	C6-N1-C2	-7.78	116.33	121.00
1	A	812	C	C2-N1-C1'	7.74	127.32	118.80
1	A	691	G	N1-C6-O6	7.73	124.54	119.90
1	A	1389	C	C6-N1-C2	7.71	123.38	120.30
1	A	1523	G	C5-C6-N1	-7.63	107.68	111.50
1	A	867	G	C8-N9-C1'	-7.58	117.14	127.00
1	A	1502	A	N7-C8-N9	7.57	117.58	113.80
1	A	429	U	C5-C6-N1	7.55	126.48	122.70
1	A	108	G	C6-C5-N7	-7.55	125.87	130.40
1	A	867	G	C6-C5-N7	-7.52	125.89	130.40
1	A	1237	C	C6-N1-C2	-7.50	117.30	120.30
1	A	812	C	P-O3'-C3'	7.45	128.64	119.70
1	A	1502	A	C2-N3-C4	-7.39	106.90	110.60
1	A	1502	A	C5-C6-N1	-7.39	114.00	117.70
1	A	431	A	N1-C6-N6	-7.37	114.18	118.60
1	A	1400	C	C6-N1-C2	-7.35	117.36	120.30
1	A	916	G	OP1-P-OP2	7.32	130.58	119.60
1	A	1126	U	C5-C6-N1	7.26	126.33	122.70
1	A	980	C	N1-C2-O2	7.21	123.23	118.90
1	A	360	A	O5'-P-OP1	-7.18	99.24	105.70
1	A	406	G	N3-C4-N9	7.17	130.30	126.00
1	A	1502	A	C4-C5-N7	7.16	114.28	110.70
1	A	634	C	C6-N1-C2	-7.14	117.44	120.30
1	A	638	G	N1-C6-O6	7.11	124.17	119.90
1	A	328	C	C5-C6-N1	7.07	124.54	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1158	C	C2-N1-C1'	6.99	126.49	118.80
1	A	917	G	C8-N9-C4	6.99	109.20	106.40
1	A	12	U	N3-C4-C5	-6.96	110.42	114.60
1	A	1346	A	P-O3'-C3'	6.96	128.05	119.70
1	A	1484	C	C6-N1-C2	-6.95	117.52	120.30
1	A	1361	G	C8-N9-C4	-6.95	103.62	106.40
1	A	867	G	N1-C6-O6	6.94	124.06	119.90
1	A	108	G	C5-N7-C8	-6.92	100.84	104.30
1	A	328	C	P-O3'-C3'	6.92	128.00	119.70
1	A	867	G	C5-C6-O6	-6.92	124.45	128.60
1	A	401	C	C6-N1-C2	-6.91	117.54	120.30
1	A	1189	C	N3-C4-C5	6.91	124.66	121.90
1	A	10	A	N1-C6-N6	6.89	122.74	118.60
1	A	21	G	C5-C6-N1	-6.88	108.06	111.50
1	A	1189	C	C6-N1-C2	6.86	123.05	120.30
1	A	888	G	N1-C6-O6	6.84	124.01	119.90
1	A	1084	G	C5-C6-O6	6.84	132.71	128.60
1	A	243	A	P-O3'-C3'	6.83	127.90	119.70
1	A	310	G	N1-C6-O6	6.82	123.99	119.90
1	A	575	G	O4'-C1'-N9	-6.81	102.75	108.20
1	A	376	G	C5-C6-N1	-6.80	108.10	111.50
1	A	147	G	N1-C6-O6	6.80	123.98	119.90
1	A	1370	G	N3-C2-N2	-6.79	115.14	119.90
1	A	203	U	O5'-P-OP2	-6.79	99.59	105.70
1	A	1502	A	O5'-P-OP2	-6.78	99.60	105.70
1	A	292	G	C6-C5-N7	-6.75	126.35	130.40
1	A	1084	G	N1-C6-O6	-6.74	115.86	119.90
1	A	1054	C	C2-N1-C1'	6.73	126.20	118.80
1	A	128	G	C5-C6-O6	-6.73	124.56	128.60
1	A	305	G	N1-C6-O6	6.70	123.92	119.90
1	A	653	A	C8-N9-C4	-6.70	103.12	105.80
1	A	975	A	N7-C8-N9	6.68	117.14	113.80
1	A	329	A	N1-C6-N6	6.67	122.60	118.60
1	A	1300	G	P-O3'-C3'	6.66	127.69	119.70
1	A	190	C	N3-C2-O2	-6.65	117.24	121.90
1	A	578	C	C6-N1-C2	-6.64	117.64	120.30
1	A	1200	C	N1-C2-O2	6.64	122.89	118.90
1	A	700	G	C8-N9-C4	6.63	109.05	106.40
1	A	430	A	C8-N9-C4	6.63	108.45	105.80
1	A	1361	G	N3-C4-C5	-6.63	125.28	128.60
1	A	1501	C	C6-N1-C2	6.61	122.94	120.30
1	A	1415	G	N1-C6-O6	6.59	123.86	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1067	A	O4'-C1'-N9	-6.59	102.93	108.20
1	A	1285	A	P-O3'-C3'	6.56	127.57	119.70
1	A	246	A	N1-C6-N6	-6.54	114.68	118.60
1	A	366	C	C2-N1-C1'	6.51	125.96	118.80
1	A	972	C	C6-N1-C2	-6.51	117.70	120.30
1	A	752	G	C5-C6-N1	-6.51	108.25	111.50
1	A	1198	G	C6-C5-N7	-6.50	126.50	130.40
1	A	1074	G	N1-C6-O6	6.49	123.80	119.90
1	A	1370	G	N1-C6-O6	6.49	123.80	119.90
1	A	799	G	N1-C6-O6	6.49	123.79	119.90
1	A	108	G	C4-C5-N7	6.48	113.39	110.80
1	A	417	C	C4-C5-C6	-6.47	114.16	117.40
1	A	1051	C	N3-C4-C5	-6.45	119.32	121.90
1	A	1084	G	N9-C4-C5	6.44	107.98	105.40
1	A	286	G	N1-C6-O6	6.44	123.77	119.90
1	A	108	G	C5-C6-O6	-6.44	124.74	128.60
1	A	888	G	C5-C6-O6	-6.43	124.74	128.60
1	A	1200	C	C2-N1-C1'	6.43	125.88	118.80
1	A	710	G	N1-C6-O6	6.43	123.76	119.90
1	A	181	G	N3-C4-C5	-6.42	125.39	128.60
1	A	62	U	N3-C2-O2	-6.42	117.71	122.20
1	A	1212	U	P-O3'-C3'	6.41	127.39	119.70
1	A	108	G	N7-C8-N9	6.41	116.30	113.10
1	A	1512	U	N3-C4-C5	-6.40	110.76	114.60
1	A	366	C	N1-C2-O2	6.39	122.73	118.90
1	A	328	C	N1-C2-O2	6.37	122.72	118.90
1	A	1511	G	N1-C6-O6	6.37	123.72	119.90
1	A	1521	G	C6-C5-N7	-6.36	126.58	130.40
1	A	550	G	N9-C4-C5	-6.36	102.86	105.40
1	A	362	G	N1-C6-O6	6.33	123.70	119.90
1	A	831	U	N3-C4-O4	-6.33	114.97	119.40
1	A	119	A	C8-N9-C4	6.32	108.33	105.80
22	X	30	G	N3-C4-N9	6.30	129.78	126.00
1	A	424	G	C2-N3-C4	6.30	115.05	111.90
1	A	559	A	P-O3'-C3'	6.30	127.26	119.70
1	A	1347	G	N7-C8-N9	-6.29	109.95	113.10
1	A	886	G	C5-C6-N1	-6.28	108.36	111.50
1	A	328	C	C2-N1-C1'	6.28	125.71	118.80
1	A	1305	G	C5-C6-N1	-6.24	108.38	111.50
1	A	1299	A	C6-C5-N7	-6.24	127.93	132.30
1	A	189	G	N1-C6-O6	6.24	123.64	119.90
1	A	406	G	N3-C4-C5	-6.22	125.49	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	449	C	C6-N1-C2	-6.22	117.81	120.30
1	A	975	A	C6-C5-N7	-6.22	127.94	132.30
1	A	300	A	C8-N9-C4	-6.19	103.33	105.80
1	A	594	G	N1-C6-O6	6.18	123.61	119.90
1	A	627	G	N3-C4-N9	6.18	129.71	126.00
1	A	1351	U	C5-C4-O4	6.17	129.60	125.90
1	A	131	C	N3-C2-O2	-6.15	117.59	121.90
1	A	576	G	O5'-P-OP1	-6.15	100.17	105.70
1	A	1542	U	N3-C2-O2	-6.13	117.91	122.20
1	A	366	C	N3-C2-O2	-6.12	117.62	121.90
1	A	917	G	OP1-P-OP2	6.12	128.78	119.60
1	A	186	C	C6-N1-C2	6.12	122.75	120.30
1	A	621	A	C8-N9-C4	-6.10	103.36	105.80
1	A	1370	G	C4-C5-C6	6.08	122.45	118.80
1	A	915	A	O5'-P-OP1	-6.06	100.24	105.70
1	A	411	A	C8-N9-C4	-6.05	103.38	105.80
1	A	1200	C	C2-N3-C4	6.04	122.92	119.90
1	A	1299	A	C8-N9-C4	-6.01	103.39	105.80
1	A	733	A	C8-N9-C4	6.00	108.20	105.80
1	A	1347	G	OP2-P-O3'	6.00	118.40	105.20
1	A	533	A	P-O3'-C3'	6.00	126.90	119.70
1	A	579	G	N1-C6-O6	5.99	123.49	119.90
1	A	1408	A	C8-N9-C4	5.98	108.19	105.80
1	A	1402	C	C6-N1-C2	-5.98	117.91	120.30
1	A	181	G	C4-N9-C1'	5.97	134.26	126.50
1	A	189	G	C5-C6-N1	-5.96	108.52	111.50
1	A	1519	A	N1-C6-N6	-5.95	115.03	118.60
1	A	1502	A	C4-N9-C1'	5.95	137.01	126.30
1	A	710	G	C5-C6-O6	-5.94	125.03	128.60
1	A	733	A	N1-C6-N6	5.94	122.16	118.60
1	A	1067	A	P-O3'-C3'	5.93	126.82	119.70
1	A	917	G	N3-C4-C5	5.93	131.56	128.60
1	A	372	C	N1-C2-O2	5.92	122.45	118.90
1	A	701	C	P-O3'-C3'	5.90	126.78	119.70
1	A	117	G	C6-C5-N7	-5.90	126.86	130.40
1	A	1299	A	C4-C5-C6	5.87	119.93	117.00
1	A	761	G	C4-N9-C1'	5.86	134.12	126.50
1	A	1450	U	C6-N1-C2	-5.85	117.49	121.00
1	A	889	A	O5'-P-OP2	-5.85	100.44	105.70
1	A	189	G	C8-N9-C4	-5.84	104.06	106.40
1	A	812	C	N1-C2-O2	5.84	122.40	118.90
1	A	190	C	C2-N1-C1'	5.83	125.22	118.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	550	G	C8-N9-C4	5.83	108.73	106.40
1	A	965	A	P-O3'-C3'	5.83	126.69	119.70
1	A	365	U	C2-N1-C1'	5.82	124.68	117.70
1	A	266	G	P-O3'-C3'	5.82	126.68	119.70
1	A	761	G	C8-N9-C1'	-5.80	119.46	127.00
1	A	558	G	C4-C5-C6	5.80	122.28	118.80
1	A	1521	G	N1-C6-O6	5.79	123.38	119.90
1	A	886	G	N1-C6-O6	5.78	123.37	119.90
1	A	1231	G	C2-N3-C4	-5.78	109.01	111.90
1	A	1074	G	C5-C6-N1	-5.78	108.61	111.50
1	A	380	G	N3-C4-N9	-5.77	122.54	126.00
1	A	429	U	N3-C4-C5	-5.77	111.14	114.60
1	A	1313	U	C6-N1-C2	-5.77	117.54	121.00
1	A	190	C	N1-C2-O2	5.76	122.36	118.90
1	A	417	C	C5-C6-N1	5.76	123.88	121.00
1	A	1286	A	C8-N9-C4	-5.76	103.50	105.80
1	A	872	A	O4'-C1'-N9	5.75	112.80	108.20
1	A	506	G	C8-N9-C4	-5.75	104.10	106.40
1	A	798	G	N1-C6-O6	5.74	123.34	119.90
1	A	1208	C	N3-C4-C5	-5.74	119.61	121.90
1	A	297	G	N1-C6-O6	5.72	123.33	119.90
1	A	331	G	N3-C2-N2	-5.71	115.90	119.90
1	A	380	G	C5-C6-N1	-5.70	108.65	111.50
1	A	5	U	C2-N1-C1'	5.69	124.53	117.70
1	A	1103	C	N3-C4-C5	5.67	124.17	121.90
1	A	190	C	O4'-C1'-N1	5.67	112.74	108.20
1	A	388	G	C8-N9-C4	-5.67	104.13	106.40
1	A	1242	C	C6-N1-C2	5.67	122.57	120.30
1	A	1299	A	N7-C8-N9	5.67	116.63	113.80
1	A	221	C	C6-N1-C2	-5.67	118.03	120.30
1	A	1198	G	N1-C6-O6	5.66	123.30	119.90
1	A	1250	A	O4'-C1'-N9	5.66	112.73	108.20
1	A	1544	U	N3-C2-O2	-5.66	118.23	122.20
1	A	866	C	N1-C2-O2	-5.66	115.50	118.90
1	A	32	A	C8-N9-C4	-5.64	103.54	105.80
1	A	739	C	C6-N1-C2	-5.64	118.04	120.30
1	A	1084	G	C4-C5-N7	-5.64	108.55	110.80
1	A	115	G	P-O3'-C3'	5.63	126.46	119.70
1	A	848	C	N3-C4-C5	-5.62	119.65	121.90
1	A	396	G	N9-C4-C5	5.62	107.65	105.40
1	A	1397	C	OP2-P-O3'	5.61	117.55	105.20
1	A	1408	A	N9-C4-C5	-5.61	103.56	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	622	A	O5'-P-OP1	-5.61	100.65	105.70
1	A	122	G	C5-C6-N1	-5.60	108.70	111.50
1	A	575	G	N1-C6-O6	-5.59	116.54	119.90
1	A	1084	G	N3-C4-N9	-5.59	122.64	126.00
1	A	246	A	C6-C5-N7	5.59	136.21	132.30
1	A	19	C	C6-N1-C2	-5.59	118.06	120.30
1	A	869	G	N1-C6-O6	5.58	123.25	119.90
1	A	1388	C	C2-N1-C1'	-5.58	112.66	118.80
1	A	299	G	C5-C6-O6	5.57	131.94	128.60
1	A	568	G	C8-N9-C4	-5.57	104.17	106.40
1	A	564	C	C6-N1-C2	-5.56	118.08	120.30
1	A	578	C	N3-C4-C5	-5.55	119.68	121.90
1	A	812	C	C5-C6-N1	5.55	123.78	121.00
1	A	975	A	C5-N7-C8	-5.55	101.12	103.90
1	A	1397	C	C6-N1-C1'	5.54	127.45	120.80
1	A	1103	C	C5-C6-N1	-5.54	118.23	121.00
1	A	1194	U	C6-N1-C2	-5.54	117.67	121.00
1	A	484	G	P-O3'-C3'	5.54	126.35	119.70
1	A	1388	C	C5-C6-N1	-5.54	118.23	121.00
1	A	419	C	C2-N1-C1'	5.53	124.89	118.80
1	A	917	G	O5'-P-OP2	-5.53	100.72	105.70
1	A	722	A	N1-C6-N6	5.53	121.92	118.60
1	A	11	G	N1-C6-O6	5.52	123.21	119.90
1	A	396	G	N3-C4-N9	-5.52	122.69	126.00
1	A	229	U	N3-C4-O4	5.51	123.26	119.40
1	A	1300	G	OP2-P-O3'	5.51	117.33	105.20
1	A	1094	G	C8-N9-C4	-5.51	104.20	106.40
1	A	1532	U	N3-C2-O2	-5.50	118.35	122.20
1	A	417	C	N3-C4-C5	5.50	124.10	121.90
1	A	503	C	C6-N1-C2	-5.49	118.10	120.30
1	A	1347	G	P-O3'-C3'	5.49	126.29	119.70
1	A	51	A	N1-C6-N6	-5.49	115.31	118.60
1	A	201	C	C6-N1-C2	-5.49	118.11	120.30
1	A	799	G	C6-C5-N7	-5.47	127.12	130.40
1	A	491	G	N1-C6-O6	5.47	123.18	119.90
1	A	789	U	N3-C2-O2	-5.47	118.37	122.20
1	A	354	G	O5'-P-OP1	-5.47	100.78	105.70
1	A	1362	C	O5'-P-OP1	5.46	117.25	110.70
1	A	1508	G	N9-C4-C5	5.45	107.58	105.40
1	A	1531	A	N7-C8-N9	5.44	116.52	113.80
1	A	575	G	C5-N7-C8	5.44	107.02	104.30
1	A	1492	A	O5'-P-OP2	-5.44	100.81	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1504	G	O5'-P-OP2	-5.43	100.81	105.70
1	A	255	G	C8-N9-C4	5.43	108.57	106.40
1	A	1158	C	N1-C2-O2	5.43	122.16	118.90
1	A	59	A	N9-C4-C5	5.42	107.97	105.80
1	A	558	G	C5-C6-N1	-5.42	108.79	111.50
1	A	556	C	N1-C2-O2	-5.42	115.65	118.90
1	A	812	C	N3-C2-O2	-5.41	118.11	121.90
1	A	1503	A	C4-C5-C6	5.41	119.70	117.00
1	A	300	A	N7-C8-N9	5.40	116.50	113.80
1	A	516	U	C6-N1-C2	-5.40	117.76	121.00
1	A	1314	C	C6-N1-C2	-5.40	118.14	120.30
1	A	1300	G	C4-N9-C1'	-5.40	119.48	126.50
1	A	742	G	N3-C2-N2	-5.40	116.12	119.90
1	A	362	G	C5-C6-N1	-5.39	108.81	111.50
1	A	1158	C	C6-N1-C1'	-5.39	114.33	120.80
1	A	1512	U	N1-C2-N3	5.39	118.13	114.90
1	A	1443	G	O5'-P-OP1	-5.38	100.86	105.70
1	A	638	G	N9-C4-C5	-5.37	103.25	105.40
1	A	1099	G	N3-C4-N9	-5.37	122.78	126.00
1	A	235	C	C6-N1-C2	5.37	122.45	120.30
1	A	1503	A	P-O3'-C3'	5.37	126.14	119.70
1	A	867	G	C4-C5-C6	5.36	122.02	118.80
1	A	1525	G	N1-C6-O6	5.36	123.12	119.90
1	A	916	G	C5-C6-O6	-5.36	125.38	128.60
1	A	816	A	C2-N3-C4	-5.36	107.92	110.60
1	A	1450	U	N3-C2-O2	-5.36	118.45	122.20
1	A	1347	G	N9-C4-C5	-5.35	103.26	105.40
1	A	129(A)	G	P-O3'-C3'	5.34	126.11	119.70
1	A	975	A	C4-C5-C6	5.34	119.67	117.00
1	A	593	G	N1-C6-O6	5.34	123.10	119.90
1	A	190(B)	C	C6-N1-C2	-5.33	118.17	120.30
1	A	563	A	O4'-C1'-N9	5.32	112.46	108.20
1	A	1503	A	N1-C6-N6	5.32	121.79	118.60
1	A	279	A	N1-C6-N6	5.32	121.79	118.60
1	A	691	G	C5-C6-N1	-5.32	108.84	111.50
1	A	796	C	C6-N1-C2	-5.32	118.17	120.30
1	A	706	A	C8-N9-C4	5.32	107.93	105.80
1	A	731	G	N1-C6-O6	5.31	123.09	119.90
1	A	196	A	N1-C6-N6	5.30	121.78	118.60
1	A	478	A	C8-N9-C4	-5.30	103.68	105.80
1	A	1300	G	N3-C4-C5	5.29	131.25	128.60
1	A	1201	A	P-O3'-C3'	5.29	126.05	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1389	C	N3-C4-C5	5.28	124.01	121.90
1	A	733	A	N9-C4-C5	-5.27	103.69	105.80
1	A	1503	A	C5-C6-N1	-5.25	115.07	117.70
1	A	653	A	N9-C4-C5	5.25	107.90	105.80
1	A	1038	C	C6-N1-C2	-5.24	118.20	120.30
1	A	858	G	C8-N9-C1'	-5.24	120.19	127.00
1	A	24	U	C6-N1-C2	5.23	124.14	121.00
1	A	618	C	C2-N1-C1'	-5.23	113.04	118.80
1	A	752	G	N1-C6-O6	5.23	123.04	119.90
1	A	406	G	C4-N9-C1'	5.23	133.29	126.50
1	A	523	A	N1-C6-N6	5.23	121.74	118.60
1	A	577	G	N1-C6-O6	5.22	123.03	119.90
1	A	1084	G	C6-C5-N7	5.22	133.53	130.40
1	A	108	G	C8-N9-C4	-5.22	104.31	106.40
1	A	277	C	N1-C2-O2	-5.21	115.78	118.90
1	A	554	C	N1-C2-O2	-5.20	115.78	118.90
1	A	89	C	C6-N1-C2	-5.20	118.22	120.30
1	A	547	A	C8-N9-C4	-5.20	103.72	105.80
1	A	292	G	N1-C6-O6	5.19	123.02	119.90
1	A	562	C	N3-C2-O2	-5.19	118.27	121.90
1	A	376	G	N1-C6-O6	5.18	123.01	119.90
1	A	570	G	C6-C5-N7	-5.18	127.29	130.40
1	A	408	A	C5-C6-N1	5.18	120.29	117.70
1	A	817	C	N3-C4-C5	-5.18	119.83	121.90
1	A	930	C	C6-N1-C2	-5.18	118.23	120.30
1	A	575	G	C4-C5-N7	-5.18	108.73	110.80
1	A	745	C	C6-N1-C2	5.18	122.37	120.30
1	A	1362	C	O5'-P-OP2	-5.18	101.04	105.70
1	A	691	G	C6-C5-N7	-5.17	127.30	130.40
1	A	788	U	C6-N1-C2	-5.17	117.90	121.00
1	A	388	G	N3-C4-C5	-5.17	126.02	128.60
1	A	799	G	C5-C6-O6	-5.17	125.50	128.60
1	A	1194	U	C5-C6-N1	5.17	125.28	122.70
1	A	945	G	C5-C6-O6	-5.17	125.50	128.60
1	A	925	G	C8-N9-C4	5.16	108.47	106.40
14	N	6	LEU	CA-CB-CG	5.16	127.17	115.30
1	A	485	G	P-O3'-C3'	5.16	125.89	119.70
1	A	858	G	C4-N9-C1'	5.16	133.21	126.50
1	A	625	G	N9-C4-C5	-5.16	103.34	105.40
1	A	628	G	N3-C4-N9	5.15	129.09	126.00
1	A	788	U	C5-C6-N1	5.15	125.27	122.70
1	A	1299	A	C4-N9-C1'	5.15	135.56	126.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	154	C	C6-N1-C2	-5.14	118.24	120.30
1	A	292	G	N9-C4-C5	-5.13	103.35	105.40
1	A	869	G	C5-N7-C8	-5.13	101.73	104.30
1	A	870	U	C2-N1-C1'	-5.13	111.54	117.70
1	A	864	A	C5-C6-N6	-5.13	119.60	123.70
1	A	1094	G	N3-C4-C5	-5.13	126.04	128.60
1	A	266	G	O4'-C1'-N9	-5.13	104.10	108.20
1	A	430	A	N9-C4-C5	-5.12	103.75	105.80
1	A	830	G	C5-C6-N1	-5.12	108.94	111.50
1	A	638	G	C5-C6-O6	-5.12	125.53	128.60
1	A	975	A	N1-C6-N6	5.12	121.67	118.60
1	A	409	G	N1-C2-N2	5.12	120.81	116.20
1	A	573	A	N3-C4-N9	-5.12	123.31	127.40
1	A	1075	C	N3-C4-C5	5.12	123.95	121.90
1	A	22	G	N1-C6-O6	5.11	122.97	119.90
1	A	328	C	N3-C2-O2	-5.11	118.32	121.90
1	A	1065	U	P-O3'-C3'	5.10	125.82	119.70
1	A	174	C	C5-C6-N1	5.10	123.55	121.00
1	A	189	G	N7-C8-N9	5.10	115.65	113.10
5	E	12	LEU	CA-CB-CG	5.10	127.02	115.30
1	A	813	U	N1-C2-O2	-5.09	119.23	122.80
1	A	914	A	C8-N9-C4	-5.09	103.76	105.80
1	A	1299	A	N1-C6-N6	5.09	121.65	118.60
1	A	190(F)	G	N9-C4-C5	5.08	107.43	105.40
1	A	1277	C	C6-N1-C2	-5.07	118.27	120.30
1	A	1082	G	C2-N3-C4	-5.07	109.36	111.90
1	A	1200	C	C6-N1-C1'	-5.07	114.72	120.80
1	A	1054	C	C6-N1-C2	-5.07	118.27	120.30
1	A	232	G	C6-C5-N7	-5.06	127.36	130.40
1	A	1128	C	C6-N1-C2	-5.06	118.28	120.30
1	A	31	G	C8-N9-C4	5.06	108.42	106.40
1	A	306	G	C5-C6-O6	-5.05	125.57	128.60
1	A	674	G	O5'-P-OP1	-5.05	101.15	105.70
1	A	190(B)	C	C5-C6-N1	5.04	123.52	121.00
1	A	1403	C	N1-C2-O2	-5.04	115.88	118.90
1	A	331	G	C2-N3-C4	-5.04	109.38	111.90
1	A	384	G	N9-C4-C5	5.04	107.41	105.40
1	A	1514	C	N3-C4-C5	5.04	123.91	121.90
1	A	229	U	C5-C4-O4	-5.03	122.88	125.90
1	A	517	G	N3-C4-C5	-5.03	126.09	128.60
1	A	1406	U	O5'-P-OP2	-5.03	101.18	105.70
1	A	975	A	C3'-C2'-C1'	5.02	105.52	101.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
22	X	35	C	C6-N1-C2	-5.02	118.29	120.30
1	A	419	C	C6-N1-C1'	-5.02	114.78	120.80
1	A	117	G	C5-C6-O6	-5.02	125.59	128.60
1	A	526	C	N3-C4-C5	5.02	123.91	121.90
1	A	980	C	N3-C2-O2	-5.02	118.39	121.90
4	D	12	CYS	CA-CB-SG	5.02	123.03	114.00
1	A	907	A	N1-C6-N6	-5.01	115.59	118.60
1	A	389	A	O5'-P-OP2	-5.01	101.19	105.70
1	A	448	A	C8-N9-C4	5.01	107.80	105.80
1	A	761	G	C6-C5-N7	-5.01	127.39	130.40
1	A	365	U	C6-N1-C1'	-5.01	114.19	121.20
1	A	1086	U	C5-C6-N1	5.00	125.20	122.70

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
20	T	12	ALA	Peptide

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	32380	0	16346	452	0
2	B	1900	0	1951	53	0
3	C	1612	0	1677	52	0
4	D	1703	0	1765	36	0
5	E	1146	0	1207	41	0
6	F	843	0	857	15	0
7	G	1257	0	1296	34	0
8	H	1116	0	1177	37	0
9	I	1011	0	1043	44	0
10	J	792	0	835	36	0
11	K	885	0	904	24	0
12	L	970	0	1057	38	0
13	M	997	0	1072	32	0
14	N	492	0	529	40	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
15	O	734	0	771	13	0
16	P	700	0	720	26	0
17	Q	857	0	930	22	0
18	R	597	0	668	22	0
19	S	647	0	673	15	0
20	T	762	0	856	27	0
21	V	208	0	221	11	0
22	X	232	0	121	1	0
23	W	82	0	45	2	0
24	A	42	0	45	2	0
25	A	76	0	0	0	0
25	E	1	0	0	0	0
25	J	1	0	0	0	0
All	All	52043	0	36766	937	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 11.

All (937) 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)
14:N:24:CYS:HB2	14:N:40:CYS:HB3	1.13	1.09
14:N:24:CYS:CB	14:N:40:CYS:HB3	1.90	1.00
14:N:24:CYS:HB2	14:N:40:CYS:CB	1.91	1.00
12:L:27:LEU:O	12:L:29:GLY:N	2.07	0.88
12:L:47:LYS:HB3	12:L:48:PRO:HD3	1.55	0.87
1:A:448:A:OP2	1:A:485:G:N2	2.08	0.85
10:J:90:LEU:H	10:J:91:PRO:HD2	1.43	0.84
1:A:1502:A:H2	1:A:1505:G:H1	1.24	0.83
1:A:542:G:OP1	4:D:10:ARG:NH2	2.11	0.83
11:K:110:ASP:HB2	18:R:88:LYS:HE2	1.61	0.82
1:A:927:G:O2'	1:A:1503:A:N7	2.11	0.82
8:H:69:ARG:NH1	8:H:75:ARG:O	2.12	0.82
1:A:664:G:H22	1:A:741:G:H1	1.28	0.80
1:A:1266:G:N2	1:A:1269:A:OP2	2.14	0.79
13:M:10:PRO:HB2	13:M:18:ALA:HB1	1.64	0.79
20:T:75:ASN:N	20:T:75:ASN:OD1	2.15	0.79
15:O:2:PRO:O	15:O:38:ARG:NH1	2.12	0.77
1:A:1356:G:H2'	1:A:1357:A:C8	2.19	0.77
1:A:1292:U:OP2	7:G:41:ARG:NH2	2.18	0.77
13:M:91:ARG:HB2	13:M:98:VAL:HG13	1.66	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:50:GLU:HG3	5:E:52:PRO:HD2	1.66	0.76
20:T:57:ARG:HH11	20:T:102:GLY:HA3	1.50	0.76
1:A:1003:G:N2	1:A:1038:C:O2	2.17	0.76
9:I:46:ALA:HB2	9:I:74:ILE:HG23	1.69	0.75
2:B:74:LYS:NZ	2:B:205:ASP:OD2	2.20	0.74
1:A:275:G:H5'	17:Q:14:LYS:HE2	1.69	0.74
1:A:477:G:H8	1:A:478:A:H4'	1.52	0.74
18:R:38:GLU:HA	18:R:41:LYS:HE2	1.68	0.74
8:H:10:LEU:HD22	8:H:83:ILE:HD11	1.70	0.73
18:R:59:SER:HB3	18:R:62:GLU:HG3	1.71	0.73
7:G:20:ASP:OD2	7:G:63:LYS:NZ	2.18	0.73
2:B:103:THR:N	2:B:176:GLU:OE1	2.17	0.73
19:S:13:ASP:HA	19:S:16:LEU:HB3	1.71	0.73
1:A:462:G:N2	16:P:82:GLN:HB3	2.04	0.73
3:C:58:GLU:HB3	10:J:92:THR:HG21	1.70	0.72
13:M:96:LEU:O	13:M:110:ARG:NH1	2.23	0.72
4:D:25:ARG:NH1	4:D:30:LYS:O	2.23	0.72
1:A:448:A:H62	1:A:486:U:H3	1.38	0.72
1:A:1392:G:H21	1:A:1502:A:H8	1.38	0.71
4:D:19:LEU:HD21	4:D:67:ILE:HG12	1.71	0.71
1:A:537:G:OP1	12:L:113:ARG:NH2	2.23	0.71
13:M:49:THR:HG22	13:M:51:ALA:H	1.54	0.71
1:A:1195:C:H3'	1:A:1196:U:H5''	1.72	0.71
8:H:4:ASP:OD2	8:H:85:ARG:NH1	2.23	0.71
1:A:1420:C:H2'	1:A:1421:G:H8	1.56	0.71
17:Q:66:SER:O	17:Q:70:ARG:NH1	2.24	0.71
11:K:40:ILE:HG22	11:K:41:THR:HG23	1.71	0.70
10:J:49:VAL:HG23	14:N:41:ARG:HB2	1.73	0.70
7:G:51:GLN:NE2	7:G:56:GLN:O	2.24	0.70
1:A:1291:G:OP1	7:G:37:ASN:ND2	2.24	0.70
15:O:26:GLU:OE1	15:O:77:ARG:NH1	2.25	0.70
3:C:155:GLY:HA3	3:C:163:ALA:HB1	1.72	0.70
5:E:33:VAL:HG12	5:E:112:LEU:HD12	1.74	0.70
1:A:975:A:H5'	1:A:975:A:H8	1.56	0.69
4:D:176:LEU:HA	4:D:183:GLY:HA2	1.74	0.69
4:D:64:LEU:HB2	4:D:198:VAL:HG11	1.74	0.69
1:A:474:G:O6	16:P:81:ARG:NH2	2.25	0.69
1:A:129(A):G:N3	1:A:190(E):U:H5''	2.08	0.69
1:A:259:G:OP2	20:T:83:ARG:NH1	2.26	0.69
1:A:462:G:C2	16:P:82:GLN:HB3	2.28	0.69
1:A:1443:G:H5''	1:A:1446:A:H5''	1.75	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:928:G:H5''	1:A:1533:C:H5''	1.74	0.68
1:A:243:A:H4'	1:A:244:U:O5'	1.93	0.68
3:C:11:ARG:O	3:C:13:GLY:N	2.27	0.68
1:A:1129:C:OP1	9:I:62:TYR:OH	2.10	0.68
1:A:1123:A:H4'	10:J:36:GLY:HA3	1.75	0.68
1:A:279:A:OP2	17:Q:95:TYR:OH	2.11	0.68
1:A:1227:A:C2	13:M:117:VAL:HG21	2.28	0.68
1:A:1221:G:OP2	19:S:37:ARG:NH2	2.26	0.68
1:A:476:G:C2	1:A:478:A:H5'	2.28	0.68
1:A:409:G:N1	1:A:433:C:H5'	2.09	0.67
1:A:477:G:C8	1:A:478:A:H4'	2.29	0.67
1:A:929:G:H1	1:A:1388:C:H42	1.43	0.67
1:A:452:A:HO2'	1:A:453:A:H8	1.42	0.66
1:A:1166:G:N2	1:A:1169:A:OP2	2.29	0.66
1:A:1372:U:H5''	9:I:71:SER:HB3	1.77	0.66
1:A:958:A:N3	1:A:985:C:O2'	2.27	0.66
3:C:14:ILE:HG22	3:C:15:THR:HG23	1.78	0.66
1:A:1203:C:OP1	14:N:2:ALA:N	2.29	0.66
12:L:124:LYS:HD2	12:L:125:PRO:HD2	1.76	0.66
13:M:22:ILE:HG22	13:M:24:GLY:H	1.61	0.66
9:I:100:GLY:O	9:I:102:LEU:N	2.29	0.66
10:J:89:ASP:HB2	10:J:91:PRO:HD2	1.78	0.65
10:J:34:VAL:HG22	10:J:74:ILE:HG23	1.78	0.65
1:A:190:C:N4	1:A:190(J):U:O2'	2.27	0.65
12:L:89:ARG:HA	12:L:97:ARG:HA	1.79	0.65
14:N:27:CYS:HB3	14:N:43:CYS:CB	2.26	0.65
1:A:974:A:OP2	14:N:41:ARG:NH1	2.29	0.65
1:A:1112:C:N3	3:C:178:LEU:N	2.36	0.65
1:A:316:G:OP2	1:A:351:G:O2'	2.15	0.65
1:A:353:A:H5'	1:A:353:A:H8	1.62	0.65
2:B:77:ALA:HB2	2:B:211:ILE:HD13	1.79	0.65
7:G:54:THR:HG22	7:G:56:GLN:H	1.62	0.65
3:C:148:GLY:HA2	3:C:171:GLY:HA3	1.79	0.64
4:D:98:GLU:OE2	4:D:107:ARG:NE	2.30	0.64
18:R:52:PRO:HB2	18:R:54:ARG:HD2	1.79	0.64
1:A:463:A:H1'	1:A:474:G:N2	2.12	0.64
1:A:477:G:H2'	1:A:479:C:OP2	1.98	0.64
1:A:1057:G:H5''	3:C:154:SER:HB2	1.80	0.64
1:A:1250:A:H2'	1:A:1251:A:C8	2.33	0.64
1:A:927:G:OP2	1:A:927:G:H4'	1.96	0.64
14:N:26:ARG:NH2	14:N:46:GLU:OE1	2.31	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:L:27:LEU:C	12:L:29:GLY:H	2.01	0.64
16:P:81:ARG:HA	16:P:81:ARG:HH11	1.63	0.64
8:H:46:LYS:HG3	8:H:64:LYS:HG3	1.78	0.63
16:P:81:ARG:NH1	16:P:81:ARG:HA	2.14	0.63
1:A:1420:C:H2'	1:A:1421:G:C8	2.32	0.63
10:J:82:ILE:HA	10:J:85:LEU:HD12	1.80	0.63
5:E:79:GLU:O	8:H:104:ARG:NH1	2.31	0.63
1:A:1298:C:C5	7:G:114:ARG:HD3	2.34	0.63
9:I:31:GLN:HB3	9:I:35:GLU:HB3	1.79	0.63
3:C:88:ARG:HG3	3:C:101:LEU:HD12	1.80	0.63
12:L:126:LYS:HD2	12:L:126:LYS:H	1.64	0.63
1:A:795:C:H5''	1:A:796:C:OP2	2.00	0.62
5:E:78:HIS:CD2	8:H:107:LEU:HD12	2.34	0.62
3:C:119:ARG:HE	3:C:140:ARG:HH12	1.47	0.62
3:C:35:GLU:HB2	3:C:59:ARG:HH22	1.63	0.62
2:B:97:TRP:HZ2	2:B:102:LEU:HD13	1.63	0.62
1:A:1367:C:H4'	10:J:48:THR:HG21	1.81	0.62
15:O:3:ILE:HD13	15:O:34:LEU:HD13	1.81	0.62
16:P:18:ARG:HD3	16:P:35:LYS:HD2	1.81	0.62
9:I:16:ARG:HD3	9:I:64:THR:HB	1.81	0.62
21:V:2:GLY:O	21:V:4:GLY:N	2.33	0.62
1:A:299:G:H2'	1:A:300:A:C8	2.35	0.61
1:A:410:G:OP1	1:A:431:A:N6	2.32	0.61
1:A:673:G:H2'	1:A:674:G:C8	2.34	0.61
1:A:1101:A:H4'	1:A:1102:A:O5'	2.01	0.61
7:G:78:ARG:HB2	7:G:156:TRP:HZ3	1.65	0.60
16:P:11:SER:HB3	16:P:14:ASN:HB3	1.82	0.60
1:A:237:C:OP2	17:Q:40:LYS:NZ	2.34	0.60
5:E:78:HIS:HD2	8:H:107:LEU:HD12	1.65	0.60
1:A:1118:C:OP1	9:I:104:ARG:NH1	2.34	0.60
10:J:42:THR:HG23	10:J:68:HIS:HA	1.83	0.60
8:H:91:ARG:HG2	12:L:7:ILE:HG13	1.84	0.60
1:A:1028:C:H42	1:A:1033:G:H1	1.48	0.60
2:B:132:LYS:HD3	2:B:135:GLN:HB2	1.83	0.60
1:A:1465:C:H2'	1:A:1466:C:O4'	2.02	0.59
2:B:18:GLY:HA2	2:B:40:HIS:O	2.02	0.59
6:F:10:LEU:HD12	6:F:59:TYR:HB3	1.84	0.59
19:S:28:LYS:HG2	19:S:31:ILE:HD11	1.84	0.59
21:V:17:THR:O	21:V:22:ARG:NH1	2.35	0.59
1:A:1057:G:H2'	1:A:1058:G:O4'	2.00	0.59
7:G:146:GLU:HA	7:G:149:ARG:HB2	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:L:70:ILE:HD13	12:L:77:LEU:HD12	1.84	0.59
16:P:10:GLY:HA3	16:P:14:ASN:O	2.02	0.59
10:J:38:ILE:HD12	10:J:71:LEU:HB3	1.84	0.59
1:A:1442:G:H2'	1:A:1446:A:N7	2.17	0.59
3:C:6:HIS:HD2	3:C:8:ILE:H	1.51	0.59
1:A:404:U:OP2	4:D:3:ARG:HD3	2.02	0.59
7:G:26:PHE:CE2	7:G:30:ILE:HD11	2.38	0.59
1:A:1360:A:OP2	14:N:35:ARG:NH2	2.35	0.59
4:D:12:CYS:HB3	4:D:33:MET:HG2	1.84	0.59
1:A:404:U:H5'	4:D:122:ARG:HD2	1.85	0.59
1:A:959:A:O2'	1:A:984:C:O2'	2.17	0.58
2:B:197:VAL:HB	2:B:200:ILE:HG23	1.85	0.58
16:P:75:ARG:HD3	16:P:82:GLN:HE22	1.67	0.58
1:A:738:C:H5''	6:F:69:GLU:HB3	1.85	0.58
9:I:10:ARG:HD3	9:I:75:ASP:HB3	1.83	0.58
20:T:57:ARG:NH1	20:T:102:GLY:HA3	2.16	0.58
2:B:80:ILE:HD12	2:B:80:ILE:H	1.68	0.58
1:A:1118:C:P	9:I:104:ARG:HH11	2.26	0.58
1:A:1269:A:H5'	21:V:19:GLY:HA2	1.84	0.58
14:N:37:PHE:HB2	14:N:39:LEU:HB2	1.84	0.58
10:J:9:ARG:NH2	10:J:69:ASN:OD1	2.34	0.58
1:A:269:C:H2'	1:A:270:A:C8	2.39	0.58
5:E:144:THR:HG23	5:E:147:ASP:H	1.68	0.58
1:A:1228:C:O2'	13:M:117:VAL:HG22	2.03	0.58
1:A:1510:U:H2'	1:A:1511:G:C8	2.38	0.58
1:A:6:G:H4'	1:A:298:A:H4'	1.84	0.58
13:M:23:TYR:HB3	13:M:67:GLU:HA	1.86	0.58
1:A:409:G:H22	1:A:433:C:P	2.26	0.57
1:A:1144:G:H21	1:A:1146:A:H62	1.50	0.57
1:A:1020:U:H2'	1:A:1021:G:C8	2.39	0.57
1:A:1089:G:H1	1:A:1096:C:H42	1.52	0.57
1:A:60:A:H4'	1:A:61:G:O5'	2.05	0.57
10:J:90:LEU:H	10:J:91:PRO:CD	2.13	0.57
13:M:49:THR:HG22	13:M:51:ALA:N	2.19	0.57
1:A:1129:C:H1'	1:A:1132:C:C5	2.39	0.57
1:A:409:G:H1	1:A:433:C:H5'	1.68	0.57
1:A:1052:U:O2'	1:A:1055:A:OP2	2.19	0.57
7:G:50:ILE:HG21	7:G:58:PRO:HA	1.86	0.57
1:A:1366:C:H2'	1:A:1367:C:H6	1.70	0.57
1:A:165:C:H2'	1:A:166:G:H8	1.70	0.57
5:E:80:ILE:HD12	5:E:91:LEU:HB2	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:V:8:THR:HB	21:V:11:GLY:H	1.69	0.57
1:A:1332:A:H2'	1:A:1333:A:C8	2.40	0.57
2:B:71:VAL:HG22	2:B:164:VAL:HA	1.86	0.57
1:A:900:A:H2'	1:A:901:A:C8	2.40	0.57
5:E:33:VAL:HG21	5:E:109:ILE:HG12	1.87	0.57
1:A:713:G:H2'	1:A:714:G:C8	2.40	0.56
2:B:97:TRP:CZ2	2:B:102:LEU:HD13	2.39	0.56
6:F:98:LEU:HD12	18:R:28:GLU:HB3	1.87	0.56
18:R:53:ARG:HD2	18:R:58:LEU:O	2.05	0.56
2:B:178:ARG:HH21	8:H:68:ARG:HH22	1.53	0.56
3:C:12:LEU:HD11	14:N:51:GLY:HA2	1.86	0.56
1:A:1477:C:H2'	1:A:1478:C:C6	2.40	0.56
3:C:11:ARG:C	3:C:13:GLY:H	2.09	0.56
11:K:45:GLY:HA3	11:K:55:LYS:HG2	1.88	0.56
12:L:55:VAL:HG12	12:L:56:ALA:H	1.69	0.56
1:A:1503:A:H5'	1:A:1531:A:H1'	1.87	0.56
5:E:76:ILE:HD13	5:E:118:ILE:HD12	1.86	0.56
1:A:1016:A:H2'	1:A:1017:G:O4'	2.05	0.56
1:A:1275:A:H2'	1:A:1276:G:C8	2.41	0.56
1:A:560:U:H5'	1:A:566:G:C2	2.40	0.56
3:C:156:ARG:H	3:C:163:ALA:HA	1.70	0.56
3:C:61:ALA:O	3:C:63:ASN:N	2.37	0.56
5:E:12:LEU:HD13	5:E:31:LEU:HB2	1.87	0.56
15:O:70:LEU:HB3	15:O:78:TYR:HB2	1.87	0.56
1:A:1086:U:O5'	1:A:1086:U:H6	1.89	0.56
8:H:104:ARG:HG3	8:H:138:TRP:CG	2.41	0.56
10:J:51:ARG:HD2	10:J:59:SER:HB3	1.86	0.56
1:A:1202:G:H21	14:N:27:CYS:HB2	1.71	0.56
1:A:1327:C:OP1	21:V:20:LYS:HB3	2.06	0.56
1:A:460:A:N6	1:A:463:A:N1	2.53	0.56
1:A:1117:G:O3'	9:I:104:ARG:HD2	2.06	0.56
9:I:20:ARG:HB2	9:I:60:ASP:HB3	1.87	0.56
1:A:1347:G:O2'	1:A:1348:U:P	2.64	0.56
1:A:424:G:H3'	1:A:424:G:N3	2.20	0.56
1:A:1077:G:N2	1:A:1080:A:OP2	2.36	0.55
1:A:252:U:H2'	1:A:253:U:C6	2.40	0.55
1:A:1397:C:H4'	1:A:1398:A:OP2	2.05	0.55
1:A:530:G:O6	23:W:3:U:O2'	2.24	0.55
3:C:150:LYS:HG3	3:C:169:ALA:HB2	1.89	0.55
1:A:1249:C:O2'	9:I:73:GLN:OE1	2.25	0.55
1:A:1286:A:H61	1:A:1355:G:P	2.30	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:175:SER:HB3	4:D:186:LEU:HD21	1.89	0.55
19:S:36:ARG:NH2	19:S:75:ALA:HB3	2.22	0.55
1:A:1427:U:H2'	1:A:1428:A:C8	2.42	0.55
9:I:9:ARG:HG3	9:I:14:VAL:HG13	1.89	0.55
10:J:20:ALA:O	10:J:24:VAL:HG23	2.07	0.55
1:A:1196:U:O2	3:C:162:GLN:NE2	2.38	0.55
1:A:192:U:C1'	20:T:103:GLY:HA2	2.37	0.55
14:N:39:LEU:HD11	14:N:47:LEU:HD12	1.89	0.55
1:A:448:A:H2'	1:A:449:C:H6	1.71	0.55
1:A:456:C:H42	1:A:476:G:H1'	1.72	0.55
5:E:100:VAL:O	5:E:107:ARG:NH2	2.40	0.54
12:L:41:ARG:HG2	12:L:42:THR:H	1.72	0.54
12:L:86:ARG:NH1	12:L:87:GLY:O	2.40	0.54
2:B:60:ASP:HB3	2:B:64:ARG:HH21	1.72	0.54
5:E:75:THR:OG1	5:E:76:ILE:N	2.40	0.54
9:I:102:LEU:HD23	9:I:102:LEU:H	1.72	0.54
17:Q:83:ASP:N	17:Q:83:ASP:OD1	2.39	0.54
1:A:192:U:H1'	20:T:103:GLY:HA2	1.89	0.54
1:A:190(A):C:H42	1:A:190(H):G:H1	1.54	0.54
1:A:109:A:H2'	1:A:326:G:N2	2.23	0.54
1:A:760:G:H22	17:Q:105:ALA:HA	1.71	0.54
1:A:1104:G:H2'	1:A:1105:A:H8	1.73	0.54
14:N:23:ARG:NH1	14:N:30:ALA:HB2	2.23	0.54
1:A:264:U:H4'	17:Q:63:ARG:HD3	1.90	0.54
1:A:432:A:H4'	1:A:433:C:C2	2.43	0.54
1:A:757:U:H2'	1:A:758:G:O4'	2.08	0.54
1:A:1112:C:O2	3:C:179:ARG:HD2	2.07	0.54
4:D:28:SER:O	4:D:30:LYS:N	2.39	0.54
1:A:1047:G:N2	1:A:1210:C:O2	2.36	0.53
1:A:1216:G:H2'	1:A:1217:C:C6	2.43	0.53
1:A:190(L):U:O2	20:T:105:SER:HB2	2.08	0.53
1:A:453:A:H62	1:A:479:C:H5	1.54	0.53
1:A:646:U:H2'	1:A:647:C:C6	2.43	0.53
4:D:140:VAL:HG11	4:D:146:ILE:HD11	1.89	0.53
1:A:723:U:O5'	1:A:723:U:H6	1.91	0.53
1:A:854:G:C2	1:A:855:G:C8	2.96	0.53
2:B:93:VAL:HG21	2:B:97:TRP:CD1	2.42	0.53
8:H:19:VAL:HG23	8:H:21:LYS:HG2	1.90	0.53
1:A:1511:G:H2'	1:A:1512:U:O4'	2.09	0.53
1:A:409:G:H5'	1:A:430:A:N7	2.24	0.53
1:A:662:G:O2'	1:A:836:G:H5'	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:127:ASN:OD1	5:E:129:ILE:N	2.39	0.53
13:M:101:GLN:OE1	13:M:101:GLN:N	2.41	0.53
1:A:1316:G:O2'	14:N:18:VAL:HG11	2.08	0.53
1:A:1329:A:P	13:M:28:ALA:HB3	2.48	0.53
13:M:88:ARG:HG3	13:M:98:VAL:HG12	1.91	0.53
1:A:1020:U:H2'	1:A:1021:G:H8	1.73	0.53
9:I:6:GLY:N	9:I:84:ALA:HB2	2.23	0.53
1:A:421:U:O2'	1:A:422:C:H3'	2.09	0.53
3:C:156:ARG:NH2	3:C:161:GLU:HA	2.24	0.53
2:B:95:GLN:O	2:B:96:ARG:HD2	2.09	0.53
8:H:38:ILE:HG13	8:H:41:ARG:NH1	2.24	0.53
16:P:4:ILE:HG13	16:P:64:ALA:HB1	1.90	0.53
16:P:81:ARG:C	16:P:83:GLU:H	2.13	0.53
1:A:1021:G:H2'	1:A:1022:G:H8	1.73	0.52
1:A:560:U:H5'	1:A:566:G:N2	2.23	0.52
5:E:145:LYS:O	5:E:149:GLU:HG2	2.08	0.52
11:K:14:VAL:O	11:K:16:SER:N	2.40	0.52
1:A:1342:C:O2'	9:I:124:GLN:HG3	2.09	0.52
1:A:407:G:O2'	1:A:431:A:P	2.67	0.52
8:H:90:GLY:O	8:H:91:ARG:HB2	2.09	0.52
9:I:93:ARG:HB3	9:I:93:ARG:NH1	2.23	0.52
5:E:80:ILE:H	5:E:80:ILE:HD12	1.75	0.52
1:A:1427:U:H2'	1:A:1428:A:H8	1.75	0.52
2:B:178:ARG:HH22	8:H:68:ARG:HH12	1.58	0.52
7:G:78:ARG:O	7:G:84:ASN:HA	2.09	0.52
13:M:82:MET:HE3	13:M:93:ARG:HE	1.75	0.52
1:A:1007:C:H42	1:A:1022:G:H1	1.56	0.52
3:C:148:GLY:HA3	3:C:172:ARG:H	1.74	0.52
3:C:3:ASN:N	3:C:3:ASN:OD1	2.41	0.52
8:H:29:SER:HB3	8:H:32:LYS:HG3	1.91	0.52
1:A:377:G:OP1	16:P:3:LYS:HD3	2.10	0.52
5:E:80:ILE:HG23	8:H:104:ARG:NH2	2.25	0.52
1:A:165:C:H2'	1:A:166:G:C8	2.45	0.52
1:A:836:G:C6	1:A:851:G:C6	2.98	0.52
2:B:43:ASP:HB3	2:B:46:LYS:HB2	1.92	0.52
3:C:123:GLN:O	3:C:128:PHE:HB2	2.09	0.52
4:D:24:GLU:HA	4:D:24:GLU:OE1	2.10	0.52
16:P:39:TYR:CE2	16:P:41:PRO:HB3	2.45	0.52
10:J:51:ARG:NH2	10:J:61:GLU:HB2	2.24	0.51
19:S:36:ARG:HH22	19:S:75:ALA:HB3	1.73	0.51
1:A:1227:A:O3'	13:M:115:LYS:NZ	2.43	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:975:A:H4'	1:A:976:G:O5'	2.09	0.51
2:B:18:GLY:HA2	2:B:41:ILE:HA	1.91	0.51
5:E:33:VAL:HG22	5:E:43:LEU:HD23	1.92	0.51
1:A:754:C:OP1	15:O:72:ARG:NH2	2.42	0.51
1:A:1128:C:HO2'	1:A:1130:A:H8	1.56	0.51
1:A:1196:U:OP1	1:A:1197:G:H5'	2.10	0.51
12:L:75:HIS:HD2	12:L:77:LEU:H	1.58	0.51
1:A:409:G:N3	1:A:409:G:H2'	2.26	0.51
2:B:229:VAL:HG12	2:B:231:GLU:HG3	1.92	0.51
2:B:30:ARG:O	2:B:46:LYS:NZ	2.29	0.51
1:A:1477:C:H2'	1:A:1478:C:H6	1.76	0.51
1:A:7:G:H5'	1:A:298:A:O4'	2.11	0.51
3:C:7:PRO:HG2	3:C:184:TYR:HB2	1.93	0.51
6:F:69:GLU:CD	6:F:69:GLU:H	2.13	0.51
14:N:27:CYS:CB	14:N:43:CYS:HB3	2.41	0.51
1:A:1225:A:N3	1:A:1225:A:H2'	2.26	0.51
3:C:154:SER:OG	3:C:155:GLY:N	2.44	0.51
4:D:152:SER:O	4:D:158:ILE:HD12	2.10	0.51
6:F:1:MET:HB3	6:F:66:GLU:HG2	1.93	0.51
7:G:147:ALA:H	7:G:149:ARG:H	1.57	0.51
1:A:130:A:H5'	17:Q:63:ARG:NE	2.25	0.51
1:A:542:G:H2'	1:A:543:C:H6	1.76	0.50
1:A:501:C:H1'	1:A:549:C:H1'	1.92	0.50
2:B:219:VAL:C	2:B:221:LEU:H	2.14	0.50
10:J:60:ARG:O	10:J:61:GLU:HB3	2.11	0.50
1:A:369:C:OP2	1:A:388:G:N2	2.39	0.50
1:A:448:A:H2'	1:A:449:C:C6	2.47	0.50
1:A:447:G:H2'	1:A:485:G:N2	2.26	0.50
1:A:521:G:OP1	12:L:54:LYS:HE3	2.12	0.50
1:A:620:C:H2'	1:A:621:A:C8	2.46	0.50
2:B:48:MET:HA	2:B:51:LEU:HD12	1.94	0.50
4:D:187:ARG:NH1	4:D:188:LEU:HB2	2.27	0.50
4:D:88:VAL:O	4:D:92:VAL:HG23	2.11	0.50
1:A:972:C:O5'	10:J:57:LYS:HD3	2.11	0.50
19:S:9:VAL:HG11	19:S:39:THR:HG21	1.93	0.50
1:A:980:C:H3'	1:A:981:U:H6	1.76	0.50
5:E:31:LEU:HD23	5:E:44:GLY:O	2.11	0.50
1:A:103:C:OP1	20:T:17:ARG:NH1	2.40	0.50
1:A:1062:U:H2'	1:A:1063:C:C6	2.47	0.50
1:A:409:G:N2	1:A:433:C:OP1	2.42	0.50
2:B:91:PRO:HG2	2:B:155:LEU:HG	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:60:GLU:HG2	4:D:202:LEU:HD12	1.93	0.50
8:H:24:THR:HG22	8:H:63:LEU:HD21	1.93	0.50
10:J:51:ARG:HG3	10:J:60:ARG:O	2.10	0.50
10:J:60:ARG:N	10:J:60:ARG:HD2	2.21	0.50
20:T:57:ARG:HD3	20:T:102:GLY:HA3	1.93	0.50
1:A:1147:C:HO2'	9:I:5:TYR:HH	1.60	0.50
1:A:517:G:N1	1:A:533:A:OP2	2.35	0.50
1:A:1216:G:H2'	1:A:1217:C:H6	1.76	0.50
1:A:1286:A:N6	1:A:1354:C:O3'	2.44	0.50
24:A:1601:PAR:O43	24:A:1601:PAR:N21	2.45	0.50
4:D:62:GLN:OE1	4:D:65:ARG:NH1	2.44	0.50
1:A:1300:G:O2'	1:A:1301:U:P	2.69	0.50
1:A:501:C:H2'	1:A:502:G:C8	2.46	0.50
7:G:15:ASP:OD1	7:G:16:LEU:N	2.44	0.50
1:A:21:G:H2'	1:A:22:G:C8	2.47	0.50
14:N:42:ILE:C	14:N:44:LEU:H	2.15	0.50
22:X:36:G:H1	23:W:1:C:H42	1.60	0.50
9:I:39:GLY:O	9:I:40:LEU:HD23	2.12	0.50
15:O:18:PHE:HD1	15:O:19:PRO:O	1.94	0.50
19:S:53:ASN:HB3	19:S:75:ALA:HB1	1.93	0.50
1:A:1129:C:H1'	1:A:1132:C:H5	1.75	0.49
1:A:62:U:H2'	1:A:63:C:C6	2.48	0.49
1:A:690:G:O6	11:K:51:LYS:NZ	2.35	0.49
2:B:17:PHE:HB3	2:B:44:LEU:HD21	1.94	0.49
11:K:84:VAL:HG21	11:K:95:ILE:HD11	1.94	0.49
20:T:50:GLU:HA	20:T:100:ILE:HG22	1.94	0.49
1:A:1000:U:H2'	1:A:1001:A:O4'	2.12	0.49
3:C:189:ALA:HB3	3:C:196:LEU:HB2	1.93	0.49
10:J:24:VAL:O	10:J:28:ARG:HB2	2.11	0.49
13:M:15:VAL:O	13:M:18:ALA:N	2.45	0.49
1:A:1441:G:H4'	1:A:1442:G:C5	2.48	0.49
1:A:1057:G:H5''	3:C:154:SER:CB	2.43	0.49
1:A:1136:U:H5''	1:A:1137:C:H5	1.78	0.49
2:B:189:ASP:OD1	2:B:189:ASP:N	2.45	0.49
14:N:12:ARG:C	14:N:14:PRO:HD3	2.33	0.49
1:A:1067:A:O5'	1:A:1067:A:H8	1.96	0.49
1:A:706:A:H1'	11:K:29:ILE:HD11	1.95	0.49
12:L:10:LEU:HB3	17:Q:32:TYR:CE2	2.46	0.49
1:A:1332:A:H2'	1:A:1333:A:H8	1.78	0.49
1:A:1366:C:H2'	1:A:1367:C:C6	2.47	0.49
1:A:273:A:N6	1:A:274:A:C6	2.80	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:J:64:GLU:OE2	10:J:66:ARG:NH1	2.44	0.49
1:A:1226:C:H4'	1:A:1227:A:OP1	2.13	0.49
1:A:1288:A:N6	1:A:1289:A:N6	2.61	0.49
1:A:839:U:OP2	1:A:840:C:N4	2.24	0.49
2:B:71:VAL:O	2:B:165:VAL:HG23	2.12	0.49
1:A:409:G:N2	1:A:431:A:O2'	2.46	0.49
1:A:975:A:H5'	1:A:975:A:C8	2.42	0.49
16:P:28:ARG:HG2	16:P:29:ASP:OD1	2.12	0.49
21:V:5:ASP:O	21:V:11:GLY:HA3	2.13	0.49
1:A:683:G:H2'	1:A:684:A:C8	2.48	0.49
12:L:46:LYS:HE2	12:L:47:LYS:HB2	1.94	0.49
1:A:1514:C:H42	1:A:1521:G:H1	1.61	0.48
1:A:320:C:H2'	1:A:321:A:C8	2.48	0.48
1:A:128:G:O2'	17:Q:3:LYS:NZ	2.46	0.48
1:A:972:C:H4'	10:J:57:LYS:HD3	1.95	0.48
2:B:127:ILE:HG13	2:B:128:GLU:OE2	2.13	0.48
1:A:1406:U:O2	1:A:1517:G:N2	2.41	0.48
3:C:71:ALA:HA	3:C:106:VAL:HB	1.94	0.48
3:C:91:LEU:HD21	3:C:99:VAL:HG22	1.95	0.48
1:A:390:C:H4'	16:P:28:ARG:HH21	1.77	0.48
1:A:1408:A:N1	24:A:1601:PAR:O61	2.37	0.48
1:A:1499:A:C2	1:A:1500:A:C8	3.01	0.48
1:A:250:A:H4'	1:A:251:G:O5'	2.13	0.48
1:A:575:G:OP1	1:A:575:G:H4'	2.13	0.48
1:A:639:G:C2	1:A:640:A:C8	3.01	0.48
1:A:865:A:H2'	1:A:866:C:C6	2.49	0.48
3:C:22:TRP:CH2	3:C:32:LEU:HB3	2.48	0.48
10:J:12:ASP:HB3	10:J:15:THR:HG22	1.94	0.48
1:A:951:G:OP2	13:M:102:ARG:NH2	2.46	0.48
1:A:979:C:OP1	1:A:1223:C:N4	2.46	0.48
10:J:51:ARG:CZ	10:J:61:GLU:HB2	2.44	0.48
1:A:1049:U:H4'	1:A:1050:G:O5'	2.12	0.48
1:A:190:C:H4'	1:A:190(A):C:OP1	2.12	0.48
1:A:603:U:H2'	1:A:604:G:C8	2.49	0.48
3:C:63:ASN:HA	3:C:99:VAL:HG12	1.95	0.48
4:D:39:PRO:O	4:D:44:GLY:HA3	2.13	0.48
9:I:28:VAL:O	9:I:30:GLY:N	2.47	0.48
15:O:82:ILE:HG23	15:O:87:ILE:O	2.13	0.48
1:A:192:U:H4'	20:T:57:ARG:HD2	1.96	0.48
1:A:1347:G:O2'	1:A:1348:U:OP2	2.31	0.48
1:A:355:C:O4'	1:A:388:G:O2'	2.25	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:428:G:H4'	1:A:429:U:O5'	2.14	0.48
1:A:538:G:C4	1:A:539:A:C8	3.01	0.48
1:A:558:G:H2'	1:A:559:A:H2	1.78	0.48
1:A:736:C:H2'	1:A:737:A:C8	2.49	0.48
1:A:1250:A:H4'	9:I:68:GLY:N	2.28	0.48
1:A:865:A:H5'	1:A:1078:U:O4	2.13	0.48
4:D:53:ASP:HB3	4:D:57:ARG:NH2	2.29	0.48
5:E:144:THR:O	5:E:148:VAL:HG23	2.13	0.48
18:R:17:SER:OG	18:R:54:ARG:HB2	2.13	0.48
1:A:409:G:C2'	1:A:409:G:N3	2.76	0.48
1:A:524:G:H2'	1:A:525:C:C6	2.49	0.48
1:A:840:C:H5''	1:A:841:U:OP1	2.14	0.48
2:B:95:GLN:C	2:B:96:ARG:HD2	2.34	0.48
3:C:191:THR:HG21	3:C:193:TYR:CZ	2.48	0.48
4:D:151:LYS:H	4:D:151:LYS:HD2	1.79	0.48
12:L:33:ARG:O	12:L:84:LEU:HD12	2.14	0.48
1:A:489:C:H2'	1:A:490:G:H8	1.78	0.48
3:C:91:LEU:HD21	3:C:99:VAL:H	1.79	0.48
1:A:1202:G:N2	14:N:43:CYS:HB3	2.29	0.48
1:A:1202:G:H21	14:N:43:CYS:HB3	1.77	0.47
16:P:81:ARG:O	16:P:83:GLU:N	2.47	0.47
1:A:1106:G:H5''	3:C:172:ARG:HG2	1.95	0.47
12:L:89:ARG:CZ	12:L:97:ARG:HD3	2.44	0.47
1:A:462:G:H1	16:P:82:GLN:HA	1.79	0.47
7:G:15:ASP:HB3	7:G:20:ASP:H	1.79	0.47
1:A:600:C:OP1	8:H:97:VAL:HG12	2.14	0.47
2:B:61:LEU:HD12	2:B:66:GLY:HA3	1.96	0.47
5:E:51:VAL:HB	5:E:52:PRO:HD3	1.96	0.47
7:G:155:ARG:CZ	7:G:155:ARG:HA	2.44	0.47
1:A:407:G:O2'	1:A:431:A:OP1	2.33	0.47
1:A:542:G:H2'	1:A:543:C:C6	2.50	0.47
5:E:80:ILE:CD1	5:E:91:LEU:HB2	2.43	0.47
8:H:104:ARG:HG2	8:H:104:ARG:HH11	1.80	0.47
12:L:125:PRO:HB2	12:L:127:GLU:HG2	1.96	0.47
19:S:9:VAL:HB	19:S:10:PHE:H	1.55	0.47
1:A:1432:G:O2'	1:A:1468:A:N6	2.47	0.47
1:A:928:G:C5'	1:A:1533:C:H5''	2.42	0.47
14:N:27:CYS:HB3	14:N:43:CYS:HB3	1.95	0.47
16:P:8:ARG:HB3	16:P:28:ARG:NH1	2.30	0.47
9:I:10:ARG:HD3	9:I:75:ASP:CB	2.44	0.47
19:S:71:LEU:H	19:S:71:LEU:HD12	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:187:C:OP1	20:T:82:SER:OG	2.32	0.47
1:A:371:G:O2'	1:A:372:C:H5'	2.15	0.47
3:C:11:ARG:C	3:C:13:GLY:N	2.68	0.47
15:O:87:ILE:HG22	15:O:88:ARG:NH1	2.29	0.47
20:T:47:GLY:O	20:T:49:ALA:N	2.47	0.47
1:A:1104:G:H2'	1:A:1105:A:C8	2.49	0.47
2:B:185:ILE:HG23	2:B:199:TYR:HB2	1.97	0.47
4:D:5:ILE:HA	4:D:115:ARG:HH22	1.80	0.47
13:M:77:ASN:O	13:M:81:LEU:HG	2.14	0.47
1:A:108:G:H5''	1:A:109:A:H5''	1.95	0.47
1:A:427:U:OP2	1:A:428:G:O2'	2.23	0.47
1:A:991:U:C4	1:A:1212:U:C2	3.03	0.47
1:A:1190:G:OP1	3:C:5:ILE:HG13	2.15	0.47
5:E:95:ALA:O	5:E:98:THR:OG1	2.29	0.47
8:H:26:VAL:HG13	8:H:59:LEU:HB2	1.96	0.47
9:I:125:TYR:CE2	9:I:128:ARG:NE	2.83	0.47
10:J:39:PRO:O	10:J:40:LEU:HB2	2.15	0.47
3:C:33:LEU:HD21	14:N:53:LEU:CD2	2.45	0.47
16:P:8:ARG:HB3	16:P:28:ARG:HH12	1.80	0.47
1:A:1241:G:H2'	1:A:1242:C:C6	2.50	0.47
1:A:1276:G:H2'	1:A:1277:C:O4'	2.16	0.47
1:A:1298:C:C4	7:G:114:ARG:HD3	2.50	0.47
1:A:901:A:C5	1:A:902:G:H1'	2.50	0.47
6:F:5:GLU:HB2	6:F:91:VAL:HG13	1.97	0.47
20:T:83:ARG:HB3	20:T:87:LYS:NZ	2.30	0.47
8:H:20:TYR:CE1	8:H:76:PRO:HD2	2.50	0.46
1:A:1300:G:O2'	1:A:1301:U:OP2	2.28	0.46
1:A:662:G:H2'	1:A:663:A:C8	2.50	0.46
1:A:739:C:OP1	15:O:2:PRO:HD3	2.15	0.46
1:A:1117:G:H5'	1:A:1118:C:OP2	2.15	0.46
1:A:1442:G:N2	1:A:1446:A:H2'	2.30	0.46
1:A:1399:C:C2	1:A:1502:A:N6	2.83	0.46
1:A:706:A:O2'	11:K:29:ILE:HD11	2.15	0.46
2:B:132:LYS:HA	2:B:135:GLN:HB2	1.96	0.46
10:J:45:ARG:HG2	10:J:47:PHE:CE2	2.51	0.46
12:L:40:VAL:HG21	12:L:77:LEU:O	2.16	0.46
21:V:2:GLY:C	21:V:4:GLY:H	2.19	0.46
1:A:812:C:H6	1:A:812:C:H2'	1.45	0.46
8:H:89:PRO:HA	8:H:92:ARG:NH1	2.30	0.46
1:A:1128:C:H5'	9:I:16:ARG:HH12	1.80	0.46
11:K:67:ASP:O	11:K:71:LYS:HG3	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:P:8:ARG:O	16:P:9:PHE:HD1	1.98	0.46
18:R:61:LYS:O	18:R:65:ILE:HG13	2.15	0.46
1:A:1128:C:H42	1:A:1143:G:H1	1.63	0.46
1:A:1302:U:OP1	13:M:13:LYS:NZ	2.41	0.46
1:A:501:C:H2'	1:A:502:G:H8	1.78	0.46
1:A:956:U:H2'	1:A:957:U:O4'	2.15	0.46
2:B:45:GLN:O	2:B:49:GLU:HG3	2.16	0.46
1:A:1298:C:H2'	7:G:114:ARG:HH12	1.81	0.46
11:K:48:ILE:HD13	11:K:48:ILE:HA	1.72	0.46
1:A:1338:G:C6	1:A:1339:A:C6	3.03	0.46
1:A:1531:A:O5'	1:A:1531:A:H8	1.99	0.46
1:A:68:G:C2	1:A:69:G:H1'	2.51	0.46
4:D:47:ARG:HH22	4:D:49:ARG:HG2	1.81	0.46
20:T:14:LYS:O	20:T:18:GLN:HG3	2.16	0.46
20:T:54:LYS:HG3	20:T:100:ILE:HD13	1.98	0.46
1:A:1119:C:H42	1:A:1154:G:H1	1.64	0.46
1:A:1474:G:H2'	1:A:1475:G:H8	1.81	0.46
1:A:479:C:H2'	1:A:480:U:O4'	2.15	0.46
1:A:980:C:H3'	1:A:981:U:C6	2.50	0.46
12:L:60:LEU:HD11	12:L:85:ILE:HD12	1.98	0.46
1:A:1417:G:H2'	1:A:1482:G:N2	2.31	0.46
1:A:353:A:H5'	1:A:353:A:C8	2.46	0.46
2:B:8:LYS:HB2	2:B:9:GLU:H	1.47	0.46
5:E:39:GLY:O	5:E:68:GLU:HA	2.15	0.46
7:G:116:ALA:HA	7:G:119:ARG:NH1	2.31	0.46
1:A:130:A:OP2	1:A:190(E):U:H2'	2.16	0.46
4:D:26:CYS:HA	4:D:31:CYS:HB2	1.96	0.46
11:K:124:LYS:HG3	11:K:125:PHE:CD1	2.51	0.46
1:A:927:G:N2	1:A:1532:U:OP1	2.44	0.46
2:B:130:ARG:HB3	2:B:131:PRO:HD2	1.97	0.46
3:C:156:ARG:HH21	3:C:161:GLU:HA	1.81	0.46
6:F:21:LEU:O	6:F:24:GLU:HB3	2.16	0.46
1:A:1347:G:OP2	9:I:107:ARG:HD2	2.16	0.46
1:A:1126:U:HO2'	1:A:1127:G:P	2.40	0.45
1:A:1128:C:O2'	1:A:1130:A:H8	1.99	0.45
1:A:338:A:H2'	1:A:339:C:C6	2.51	0.45
1:A:933:G:O6	7:G:3:ARG:NH2	2.49	0.45
2:B:189:ASP:OD1	2:B:205:ASP:HB3	2.16	0.45
3:C:112:SER:O	3:C:116:VAL:HG23	2.16	0.45
3:C:155:GLY:HA2	3:C:164:ARG:H	1.80	0.45
8:H:2:LEU:HA	8:H:2:LEU:HD23	1.75	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:J:9:ARG:HE	10:J:9:ARG:HB3	1.18	0.45
1:A:376:G:H5''	16:P:5:ARG:HB2	1.98	0.45
21:V:14:TRP:CZ3	21:V:15:ARG:HG3	2.51	0.45
1:A:1230:C:H1'	13:M:126:LYS:HA	1.98	0.45
1:A:38:G:H22	1:A:397:A:H5''	1.80	0.45
6:F:101:ALA:HA	18:R:28:GLU:HG3	1.97	0.45
7:G:88:PRO:HG3	7:G:149:ARG:HA	1.98	0.45
14:N:27:CYS:CB	14:N:43:CYS:CB	2.93	0.45
19:S:50:ALA:HA	19:S:58:VAL:O	2.16	0.45
1:A:1001:A:H2'	1:A:1002:G:C8	2.51	0.45
1:A:409:G:C6	1:A:433:C:H5'	2.51	0.45
1:A:1154:G:H2'	1:A:1155:G:H8	1.80	0.45
1:A:458:C:H2'	1:A:459:G:H8	1.82	0.45
1:A:500:G:H2'	1:A:501:C:C6	2.50	0.45
6:F:21:LEU:O	6:F:25:ILE:HG13	2.17	0.45
7:G:59:LEU:HD11	7:G:63:LYS:NZ	2.31	0.45
9:I:4:TYR:CZ	9:I:88:TYR:HD1	2.35	0.45
14:N:29:ARG:HE	14:N:31:ARG:HG3	1.81	0.45
14:N:42:ILE:C	14:N:44:LEU:N	2.69	0.45
17:Q:85:VAL:O	17:Q:89:LEU:N	2.33	0.45
20:T:14:LYS:HA	20:T:17:ARG:HG3	1.99	0.45
2:B:82:ARG:HB3	2:B:94:ASN:ND2	2.32	0.45
13:M:37:THR:HG22	13:M:55:ARG:NH1	2.31	0.45
16:P:3:LYS:HG3	16:P:24:ALA:HB2	1.98	0.45
1:A:1070:U:H2'	1:A:1071:C:C6	2.52	0.45
1:A:1442:G:H2'	1:A:1446:A:C8	2.51	0.45
1:A:1448:C:H2'	1:A:1449:C:H6	1.81	0.45
1:A:321:A:H2	1:A:332:G:H22	1.65	0.45
1:A:410:G:H2'	1:A:411:A:O4'	2.16	0.45
9:I:82:ALA:HB1	9:I:96:LEU:HD23	1.97	0.45
1:A:1480:G:H2'	1:A:1481:U:C6	2.52	0.45
1:A:35:G:H2'	1:A:36:C:C6	2.51	0.45
2:B:53:ARG:HG2	2:B:56:ARG:NH2	2.32	0.45
4:D:90:GLY:CA	4:D:204:ILE:HD11	2.47	0.45
7:G:29:LYS:HD3	7:G:29:LYS:HA	1.80	0.45
10:J:15:THR:O	10:J:19:SER:HB3	2.16	0.45
19:S:22:LEU:HD22	19:S:28:LYS:HD3	1.97	0.45
1:A:1303:C:N4	1:A:1304:G:C6	2.85	0.45
1:A:1417:G:H2'	1:A:1482:G:H22	1.81	0.45
1:A:460:A:H2'	1:A:461:C:H5'	1.98	0.45
1:A:463:A:C2	16:P:82:GLN:HB2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:974:A:P	14:N:29:ARG:HH21	2.40	0.45
20:T:10:LEU:HG	20:T:12:ALA:H	1.80	0.45
1:A:1399:C:H4'	1:A:1400:C:H5''	1.97	0.45
1:A:200:G:H2'	1:A:201:C:O4'	2.16	0.45
1:A:445:G:H2'	1:A:446:G:H8	1.82	0.45
5:E:99:GLY:O	5:E:117:ASP:HA	2.16	0.45
1:A:130:A:H5'	17:Q:63:ARG:HE	1.82	0.45
1:A:389:A:C6	1:A:390:C:H1'	2.52	0.45
1:A:563:A:H2'	1:A:567:G:C8	2.51	0.45
2:B:88:ALA:HB2	2:B:219:VAL:HG13	1.98	0.45
1:A:939:G:H5''	7:G:102:ARG:NH2	2.32	0.45
1:A:113:G:H1'	1:A:354:G:H5'	1.99	0.44
3:C:85:ARG:O	3:C:89:GLU:HB2	2.17	0.44
6:F:69:GLU:O	6:F:72:VAL:HG22	2.17	0.44
1:A:485:G:O2'	1:A:486:U:P	2.75	0.44
1:A:93:G:C6	1:A:95:U:C4	3.05	0.44
2:B:69:LEU:HD12	2:B:159:PRO:HG3	1.99	0.44
2:B:15:VAL:HG11	2:B:209:ARG:HB3	1.99	0.44
4:D:152:SER:HB3	4:D:155:LEU:HD12	2.00	0.44
6:F:75:LEU:HD13	6:F:79:LEU:HD12	1.99	0.44
7:G:143:ARG:NH1	7:G:144:MET:SD	2.91	0.44
3:C:164:ARG:HG2	3:C:165:THR:H	1.82	0.44
11:K:98:LEU:HD23	11:K:98:LEU:HA	1.82	0.44
12:L:28:LYS:HB2	12:L:33:ARG:HH22	1.82	0.44
13:M:46:LYS:HG3	13:M:47:ASP:OD2	2.17	0.44
1:A:1068:G:H8	1:A:1068:G:OP2	2.00	0.44
1:A:1286:A:H2'	1:A:1287:A:H4'	1.98	0.44
1:A:376:G:H2'	1:A:377:G:H8	1.82	0.44
5:E:82:VAL:HG21	5:E:138:ALA:HA	1.98	0.44
5:E:75:THR:C	5:E:76:ILE:HD12	2.37	0.44
12:L:75:HIS:HB2	12:L:77:LEU:HG	1.99	0.44
1:A:987:G:H1	1:A:1218:C:H42	1.64	0.44
1:A:1514:C:N4	1:A:1521:G:H1	2.15	0.44
8:H:92:ARG:HH11	8:H:92:ARG:HG2	1.82	0.44
10:J:6:ILE:HG23	10:J:98:ILE:HG22	2.00	0.44
12:L:27:LEU:HD13	12:L:64:TYR:HE1	1.83	0.44
1:A:223:U:OP1	20:T:68:LYS:NZ	2.50	0.44
1:A:586:C:C2'	1:A:587:G:H5'	2.48	0.44
1:A:73:C:H2'	1:A:74:C:C6	2.53	0.44
12:L:46:LYS:HG2	12:L:92:ASP:O	2.17	0.44
13:M:84:ILE:HG21	19:S:65:ASN:HB2	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1048:G:OP1	14:N:4:LYS:N	2.34	0.44
1:A:192:U:H2'	1:A:193:C:H6	1.83	0.44
8:H:121:ASP:OD1	8:H:121:ASP:N	2.47	0.44
8:H:20:TYR:HA	8:H:65:TYR:CZ	2.53	0.44
10:J:74:ILE:HG22	10:J:75:ILE:H	1.83	0.44
15:O:25:THR:HG21	15:O:70:LEU:HD23	2.00	0.44
18:R:53:ARG:NH1	18:R:60:GLY:N	2.66	0.44
1:A:298:A:H2'	1:A:299:G:O4'	2.18	0.44
1:A:839:U:H5'	1:A:840:C:H5	1.82	0.44
2:B:217:ARG:HA	2:B:220:ASP:HB2	1.99	0.44
4:D:7:PRO:HG2	4:D:10:ARG:HD2	1.99	0.44
9:I:114:TYR:CD1	10:J:60:ARG:HB2	2.53	0.44
1:A:881:G:P	12:L:12:ARG:HH22	2.40	0.44
1:A:974:A:P	14:N:41:ARG:HH12	2.40	0.44
1:A:1223:C:P	19:S:78:ARG:HH12	2.40	0.44
1:A:1132:C:H2'	1:A:1133:G:H8	1.83	0.44
1:A:452:A:O2'	1:A:453:A:H8	2.00	0.44
1:A:573:A:O2'	1:A:574:A:H5'	2.17	0.44
4:D:150:GLU:OE1	4:D:153:ARG:NH1	2.51	0.44
6:F:13:ASN:O	6:F:14:LEU:HD23	2.17	0.44
17:Q:97:SER:HB2	17:Q:98:LEU:H	1.67	0.44
1:A:1392:G:N2	1:A:1502:A:H8	2.11	0.43
1:A:197:A:N6	1:A:221:C:H5'	2.32	0.43
1:A:419:C:C4	1:A:425:G:C6	3.06	0.43
1:A:794:A:H2'	1:A:795:C:C6	2.53	0.43
2:B:218:ALA:O	2:B:222:ILE:HG13	2.18	0.43
7:G:113:GLU:O	7:G:119:ARG:HD3	2.18	0.43
14:N:23:ARG:CZ	14:N:30:ALA:HB2	2.48	0.43
1:A:927:G:N2	1:A:1391:U:H1'	2.33	0.43
1:A:1518:A:H2'	1:A:1519:A:C8	2.54	0.43
1:A:663:A:O2'	1:A:664:G:H5'	2.18	0.43
5:E:45:PHE:CE2	5:E:47:LYS:HD2	2.53	0.43
7:G:73:MET:HG2	7:G:90:GLU:HA	2.00	0.43
8:H:51:VAL:HG11	8:H:60:ARG:CZ	2.49	0.43
1:A:1349:A:OP2	9:I:118:LYS:NZ	2.51	0.43
12:L:33:ARG:HD2	12:L:33:ARG:HA	1.60	0.43
13:M:57:ARG:O	13:M:61:GLU:HG3	2.17	0.43
1:A:1126:U:H3	1:A:1127:G:N2	2.15	0.43
1:A:44:G:N2	1:A:399:G:C4	2.86	0.43
1:A:76:C:H2'	1:A:77:G:H8	1.83	0.43
3:C:113:ALA:HB2	3:C:183:ASP:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1297:C:OP2	13:M:44:ARG:NH2	2.51	0.43
13:M:88:ARG:HG3	13:M:98:VAL:CG1	2.48	0.43
18:R:17:SER:HB2	18:R:54:ARG:HE	1.82	0.43
1:A:1256:A:H2	1:A:1277:C:N4	2.15	0.43
1:A:975:A:O2'	14:N:32:SER:HA	2.19	0.43
11:K:32:ILE:O	11:K:40:ILE:N	2.50	0.43
12:L:115:LYS:O	12:L:117:ARG:N	2.52	0.43
16:P:21:VAL:HG21	16:P:59:TRP:CD1	2.54	0.43
1:A:1010:G:N1	1:A:1019:C:O2	2.47	0.43
1:A:190(A):C:N4	1:A:190(H):G:H1	2.17	0.43
1:A:363:A:OP1	12:L:33:ARG:HD2	2.18	0.43
1:A:598:U:H4'	8:H:94:TYR:CD2	2.54	0.43
1:A:701:C:H5''	1:A:703:G:O4'	2.18	0.43
6:F:53:ALA:C	6:F:55:ASP:H	2.21	0.43
9:I:53:VAL:HG11	9:I:92:TYR:CE2	2.53	0.43
12:L:54:LYS:HG2	12:L:75:HIS:CE1	2.54	0.43
13:M:125:ARG:HD2	13:M:125:ARG:C	2.39	0.43
1:A:1028:C:N4	1:A:1033:G:H1	2.13	0.43
1:A:707:C:H4'	11:K:20:TYR:CD1	2.53	0.43
12:L:117:ARG:NH2	12:L:124:LYS:HA	2.34	0.43
1:A:1327:C:H5''	21:V:20:LYS:HE2	2.00	0.43
1:A:1299:A:C5	1:A:1301:U:O2	2.71	0.43
1:A:259:G:OP1	20:T:83:ARG:HD3	2.18	0.43
1:A:826:C:H2'	1:A:827:U:H6	1.84	0.43
2:B:29:ALA:HA	2:B:32:ILE:HD12	1.99	0.43
17:Q:79:SER:OG	17:Q:80:GLY:N	2.52	0.43
1:A:243:A:C2	1:A:246:A:C8	3.06	0.43
1:A:646:U:H2'	1:A:647:C:H6	1.84	0.43
1:A:866:C:H2'	1:A:867:G:O4'	2.19	0.43
3:C:35:GLU:OE1	3:C:95:THR:HB	2.19	0.43
3:C:58:GLU:HB3	10:J:92:THR:CG2	2.43	0.43
17:Q:94:ASN:O	17:Q:98:LEU:HD13	2.19	0.43
1:A:622:A:C8	1:A:623:C:C6	3.07	0.43
5:E:76:ILE:HG22	5:E:78:HIS:H	1.84	0.43
10:J:60:ARG:HA	10:J:60:ARG:HD2	1.75	0.43
17:Q:63:ARG:HG2	17:Q:64:PRO:HD2	2.01	0.43
20:T:10:LEU:O	20:T:13:LEU:HG	2.18	0.43
1:A:1406:U:H2'	1:A:1407:C:O4'	2.19	0.43
1:A:738:C:H6	1:A:738:C:O5'	2.02	0.43
2:B:178:ARG:NE	2:B:196:LEU:O	2.49	0.43
5:E:53:LEU:H	5:E:53:LEU:HD12	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1060:C:O2	1:A:1198:G:C2	2.72	0.42
1:A:1370:G:H2'	1:A:1371:G:H8	1.83	0.42
1:A:660:G:C2	1:A:746:A:C2	3.06	0.42
1:A:410:G:P	1:A:431:A:H62	2.41	0.42
2:B:71:VAL:HA	2:B:93:VAL:HG23	2.01	0.42
12:L:76:ASN:N	12:L:76:ASN:OD1	2.52	0.42
1:A:1251:A:H4'	9:I:12:GLU:HG2	2.01	0.42
1:A:148:G:H2'	1:A:149:A:C8	2.55	0.42
1:A:193:C:H2'	1:A:194:C:H6	1.85	0.42
4:D:24:GLU:O	4:D:25:ARG:HB2	2.20	0.42
15:O:78:TYR:CZ	15:O:82:ILE:HD11	2.54	0.42
15:O:87:ILE:HG22	15:O:88:ARG:HH12	1.84	0.42
1:A:1148:U:O3'	9:I:14:VAL:HG11	2.20	0.42
1:A:1495:U:H2'	1:A:1496:C:H6	1.84	0.42
1:A:262:A:C6	1:A:263:A:C6	3.07	0.42
2:B:82:ARG:HB2	2:B:82:ARG:HE	1.72	0.42
2:B:166:ASP:OD2	2:B:169:LYS:HB2	2.20	0.42
3:C:73:PRO:HA	3:C:76:VAL:HG23	2.01	0.42
8:H:12:ARG:NH1	8:H:25:ASP:O	2.51	0.42
8:H:65:TYR:HA	8:H:79:VAL:HG23	2.02	0.42
14:N:12:ARG:HD3	14:N:12:ARG:HA	1.79	0.42
1:A:1216:G:H5''	14:N:5:ALA:CB	2.48	0.42
17:Q:59:ILE:HD13	17:Q:59:ILE:HA	1.65	0.42
18:R:47:THR:HG22	18:R:48:GLY:H	1.84	0.42
20:T:56:MET:HE3	20:T:88:VAL:HG11	2.00	0.42
1:A:1262:C:H42	1:A:1273:G:H1	1.67	0.42
1:A:1287:A:H2'	1:A:1288:A:C8	2.55	0.42
1:A:1541:U:H6	1:A:1541:U:O5'	2.02	0.42
1:A:480:U:H5'	1:A:481:G:OP2	2.20	0.42
1:A:857:C:H2'	1:A:858:G:O4'	2.19	0.42
3:C:10:PHE:CE2	3:C:178:LEU:HD13	2.54	0.42
7:G:95:ARG:HG2	7:G:99:LEU:HD12	2.02	0.42
9:I:70:LYS:HA	9:I:73:GLN:NE2	2.35	0.42
9:I:85:LEU:HA	9:I:85:LEU:HD12	1.91	0.42
1:A:1014:A:H5'	19:S:14:HIS:CD2	2.54	0.42
1:A:359:U:H2'	1:A:360:A:C8	2.55	0.42
1:A:938:A:N6	1:A:939:G:C6	2.88	0.42
2:B:21:ARG:HB2	2:B:22:LYS:H	1.56	0.42
7:G:148:ASN:OD1	7:G:148:ASN:N	2.51	0.42
10:J:8:LEU:HB3	10:J:16:LEU:HD22	2.02	0.42
12:L:102:ARG:HB2	12:L:120:TYR:HA	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:974:A:P	14:N:29:ARG:NH2	2.93	0.42
1:A:1021:G:H2'	1:A:1022:G:C8	2.51	0.42
1:A:1070:U:H2'	1:A:1071:C:H6	1.85	0.42
1:A:1095:U:OP1	1:A:1108:G:N2	2.42	0.42
1:A:1345:U:OP1	9:I:120:ARG:NH1	2.48	0.42
1:A:1483:A:H2'	1:A:1484:C:O4'	2.20	0.42
1:A:962:C:H2'	1:A:963:G:O4'	2.19	0.42
12:L:90:VAL:HG12	12:L:93:LEU:H	1.84	0.42
1:A:1192:C:H2'	1:A:1193:G:O4'	2.20	0.42
1:A:109:A:C6	1:A:326:G:C6	3.08	0.42
1:A:376:G:H2'	1:A:377:G:C8	2.54	0.42
7:G:50:ILE:HG22	7:G:51:GLN:HE21	1.85	0.42
1:A:1525:G:OP1	11:K:120:ARG:NH2	2.52	0.42
1:A:797:C:OP1	11:K:124:LYS:HE2	2.20	0.42
13:M:15:VAL:HG23	13:M:43:THR:O	2.19	0.42
16:P:74:LEU:O	16:P:79:VAL:HG23	2.20	0.42
1:A:1007:C:N4	1:A:1022:G:H1	2.17	0.42
1:A:1279:A:O2'	1:A:1282:C:N4	2.52	0.42
1:A:1323:G:H2'	1:A:1324:A:C8	2.55	0.42
1:A:1502:A:H2	1:A:1505:G:N1	2.02	0.42
3:C:6:HIS:CD2	3:C:8:ILE:H	2.33	0.42
3:C:95:THR:O	3:C:97:LYS:HE2	2.19	0.42
18:R:16:PRO:HB2	18:R:54:ARG:HH21	1.85	0.42
18:R:48:GLY:O	18:R:74:ARG:NH2	2.53	0.42
20:T:53:LEU:HB2	20:T:100:ILE:HG23	2.02	0.42
1:A:1167:A:H8	1:A:1167:A:OP1	2.02	0.41
1:A:266:G:H5'	1:A:268:C:H41	1.85	0.41
1:A:664:G:OP1	18:R:64:ARG:HD2	2.20	0.41
9:I:92:TYR:HB3	9:I:96:LEU:HD13	2.00	0.41
11:K:108:ILE:O	18:R:87:ARG:N	2.52	0.41
1:A:1004:A:O2'	1:A:1038:C:O2	2.17	0.41
1:A:949:A:C2	1:A:1233:G:N3	2.88	0.41
1:A:1248:A:H2'	1:A:1249:C:C6	2.55	0.41
1:A:578:C:H2'	1:A:579:G:O4'	2.20	0.41
1:A:61:G:H2'	1:A:62:U:O4'	2.21	0.41
1:A:945:G:C2	1:A:946:A:C8	3.07	0.41
2:B:22:LYS:HE2	2:B:35:GLU:OE2	2.20	0.41
2:B:80:ILE:HG21	2:B:212:GLN:HA	2.02	0.41
14:N:11:LYS:HB2	14:N:11:LYS:HE3	1.77	0.41
1:A:1010:G:H2'	1:A:1011:G:C8	2.56	0.41
1:A:252:U:H2'	1:A:253:U:C5	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:899:C:H2'	1:A:900:A:C8	2.54	0.41
1:A:929:G:H1	1:A:1388:C:N4	2.14	0.41
3:C:188:LEU:HD22	3:C:188:LEU:HA	1.76	0.41
3:C:96:GLY:O	3:C:97:LYS:HD3	2.20	0.41
5:E:34:VAL:HG12	5:E:62:ALA:HB1	2.02	0.41
5:E:76:ILE:HG23	5:E:77:PRO:HD2	2.02	0.41
9:I:118:LYS:O	9:I:119:ALA:HB3	2.20	0.41
12:L:112:ASP:O	12:L:114:LYS:HE2	2.20	0.41
15:O:5:LYS:HE3	15:O:5:LYS:HB2	1.40	0.41
1:A:1048:G:P	14:N:4:LYS:HG2	2.60	0.41
1:A:115:G:H8	1:A:115:G:O5'	2.03	0.41
1:A:309:G:H2'	1:A:310:G:H8	1.86	0.41
4:D:191:ARG:O	4:D:191:ARG:HD2	2.20	0.41
6:F:44:GLY:HA2	6:F:59:TYR:CE1	2.55	0.41
6:F:61:LEU:HD23	6:F:63:TYR:OH	2.20	0.41
9:I:48:GLU:N	9:I:49:PRO:HD2	2.34	0.41
11:K:92:GLU:O	11:K:96:ARG:HD3	2.21	0.41
12:L:5:PRO:HG2	12:L:10:LEU:HD21	2.01	0.41
17:Q:84:LEU:HD23	17:Q:84:LEU:HA	1.90	0.41
18:R:17:SER:OG	18:R:54:ARG:O	2.35	0.41
1:A:1212:U:H4'	1:A:1213:A:O5'	2.21	0.41
1:A:1236:A:OP1	21:V:3:LYS:N	2.46	0.41
1:A:266:G:C5'	1:A:268:C:H41	2.33	0.41
1:A:407:G:O2'	1:A:431:A:OP2	2.36	0.41
5:E:41:VAL:HG22	5:E:113:ALA:HA	2.03	0.41
5:E:50:GLU:HB3	5:E:53:LEU:HD13	2.02	0.41
8:H:103:VAL:CG2	8:H:136:GLU:HG2	2.50	0.41
8:H:13:ILE:O	8:H:17:THR:HG23	2.21	0.41
8:H:48:TYR:HA	8:H:60:ARG:O	2.21	0.41
12:L:69:TYR:HE2	12:L:71:PRO:HA	1.85	0.41
13:M:4:ILE:HD13	13:M:56:LEU:HD22	2.01	0.41
20:T:39:LYS:HD3	20:T:55:ILE:HD13	2.03	0.41
1:A:66:G:H4'	1:A:173:U:C5	2.55	0.41
1:A:142:G:O2'	1:A:196:A:N1	2.40	0.41
1:A:812:C:OP1	1:A:903:G:H1'	2.20	0.41
5:E:139:LEU:HA	5:E:139:LEU:HD23	1.75	0.41
7:G:65:ALA:O	7:G:69:VAL:HG23	2.21	0.41
8:H:104:ARG:NH2	8:H:138:TRP:CH2	2.88	0.41
1:A:1157:A:C6	1:A:1180:A:C6	3.08	0.41
1:A:1256:A:H5''	1:A:1257:U:OP1	2.21	0.41
1:A:264:U:H2'	1:A:265:G:O4'	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:519:C:H2'	1:A:520:A:C8	2.55	0.41
1:A:77:G:O2'	1:A:78:G:H5'	2.20	0.41
2:B:44:LEU:HD23	2:B:44:LEU:HA	1.67	0.41
1:A:933:G:OP2	7:G:3:ARG:HB3	2.20	0.41
11:K:126:ARG:HB3	11:K:127:LYS:H	1.46	0.41
11:K:16:SER:HA	11:K:79:SER:O	2.20	0.41
11:K:54:ARG:HG2	11:K:54:ARG:H	1.46	0.41
11:K:98:LEU:O	11:K:101:SER:OG	2.30	0.41
12:L:7:ILE:HA	12:L:7:ILE:HD13	1.77	0.41
21:V:6:ARG:HG3	21:V:6:ARG:H	1.51	0.41
1:A:1201:A:H4'	1:A:1202:G:O5'	2.21	0.41
1:A:1268:A:N3	1:A:1326:C:O2'	2.53	0.41
1:A:1417:G:C6	1:A:1482:G:C6	3.08	0.41
1:A:413:G:OP2	1:A:417:C:N4	2.53	0.41
1:A:411:A:N6	1:A:418:C:H42	2.18	0.41
1:A:801:U:H2'	1:A:802:A:C8	2.56	0.41
3:C:29:TYR:OH	14:N:54:PRO:HG2	2.21	0.41
4:D:170:VAL:HG22	4:D:171:GLY:H	1.85	0.41
13:M:3:ARG:HE	13:M:7:VAL:HA	1.85	0.41
18:R:46:GLU:H	18:R:46:GLU:CD	2.24	0.41
20:T:36:LEU:HA	20:T:36:LEU:HD23	1.78	0.41
1:A:1095:U:H2'	1:A:1096:C:C6	2.56	0.41
1:A:484:G:H4'	1:A:485:G:O5'	2.21	0.41
5:E:11:ILE:HG21	5:E:105:VAL:HG22	2.03	0.41
5:E:31:LEU:HD22	5:E:43:LEU:HD21	2.02	0.41
9:I:56:LEU:HB3	9:I:57:GLY:H	1.76	0.41
14:N:21:TYR:HE1	14:N:23:ARG:NE	2.18	0.41
14:N:42:ILE:O	14:N:44:LEU:N	2.54	0.41
16:P:19:ILE:HG22	16:P:36:ILE:HG13	2.03	0.41
18:R:58:LEU:HA	18:R:58:LEU:HD23	1.78	0.41
1:A:1305:G:N2	1:A:1331:G:H1'	2.36	0.41
1:A:1505:G:H3'	1:A:1505:G:C8	2.55	0.41
3:C:120:VAL:O	3:C:124:ILE:HG13	2.21	0.41
9:I:53:VAL:HG11	9:I:92:TYR:CZ	2.55	0.41
13:M:34:LEU:HD13	13:M:34:LEU:HA	1.93	0.41
1:A:570:G:H1'	1:A:820:U:C4	2.56	0.41
11:K:51:LYS:C	11:K:51:LYS:HD3	2.41	0.41
17:Q:27:PHE:O	17:Q:36:ILE:HG12	2.21	0.41
17:Q:5:VAL:HG22	17:Q:60:ILE:HG12	2.02	0.41
18:R:88:LYS:HB3	18:R:88:LYS:HE3	1.94	0.41
20:T:20:LEU:HD23	20:T:20:LEU:HA	1.79	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1167:A:C6	1:A:1168:A:C6	3.09	0.40
1:A:1250:A:H5'	9:I:68:GLY:O	2.21	0.40
1:A:691:G:H2'	1:A:692:U:C6	2.56	0.40
1:A:760:G:H1	17:Q:105:ALA:CB	2.34	0.40
1:A:1193:G:O2'	5:E:25:ARG:NH2	2.54	0.40
7:G:78:ARG:HG2	7:G:80:VAL:HG13	2.02	0.40
10:J:27:ALA:HB2	10:J:85:LEU:HD21	2.03	0.40
1:A:1007:C:N3	1:A:1022:G:N2	2.62	0.40
1:A:177:C:H2'	1:A:178:C:H6	1.86	0.40
1:A:324:G:N2	1:A:327:A:C8	2.90	0.40
1:A:500:G:H2'	1:A:501:C:H6	1.86	0.40
1:A:547:A:H4'	1:A:548:G:O5'	2.21	0.40
5:E:11:ILE:CG2	5:E:105:VAL:HG22	2.51	0.40
8:H:14:ARG:HB3	8:H:14:ARG:HH11	1.86	0.40
9:I:19:LEU:HD23	9:I:19:LEU:HA	1.78	0.40
9:I:4:TYR:CE1	9:I:88:TYR:HD1	2.39	0.40
1:A:1481:U:H2'	1:A:1482:G:C8	2.57	0.40
1:A:45:U:H2'	1:A:46:G:C8	2.56	0.40
1:A:853:G:C2	1:A:854:G:C8	3.08	0.40
1:A:967:C:OP1	1:A:969:A:H5'	2.21	0.40
1:A:974:A:OP2	14:N:29:ARG:NH2	2.54	0.40
1:A:974:A:H4'	1:A:975:A:H3'	2.02	0.40
2:B:52:GLU:O	2:B:56:ARG:HB2	2.20	0.40
4:D:122:ARG:NH2	4:D:134:ASP:OD1	2.54	0.40
4:D:102:ASP:HB3	4:D:136:PRO:HB3	2.03	0.40
4:D:90:GLY:HA2	4:D:204:ILE:HD11	2.03	0.40
1:A:878:G:H5'	8:H:89:PRO:HG2	2.02	0.40
1:A:1448:C:H2'	1:A:1449:C:C6	2.56	0.40
1:A:26:A:O2'	4:D:209:ARG:NH1	2.52	0.40
1:A:392:G:H2'	1:A:393:A:C8	2.55	0.40
1:A:938:A:N6	1:A:939:G:C5	2.89	0.40
1:A:939:G:H2'	1:A:940:C:C6	2.57	0.40
4:D:97:LEU:HD23	4:D:97:LEU:HA	1.88	0.40
5:E:131:ILE:HD13	5:E:131:ILE:HA	1.85	0.40
5:E:15:ARG:HD3	5:E:26:PHE:CG	2.57	0.40
1:A:921:U:O2	5:E:19:MET:HB2	2.22	0.40
7:G:23:VAL:O	7:G:27:ILE:HG13	2.22	0.40
2:B:195:ASP:O	8:H:68:ARG:NH2	2.54	0.40
9:I:29:ASN:HB3	9:I:64:THR:HA	2.02	0.40
10:J:76:ASN:O	10:J:76:ASN:ND2	2.55	0.40
11:K:66:LEU:HA	11:K:66:LEU:HD23	1.84	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:M:14:ARG:HG2	13:M:15:VAL:N	2.36	0.40
13:M:3:ARG:HA	13:M:8:GLU:O	2.21	0.40
18:R:52:PRO:HB2	18:R:54:ARG:CD	2.48	0.40
18:R:68:LYS:HB2	18:R:68:LYS:HE2	1.95	0.40
19:S:81:ARG:HB3	19:S:81:ARG:HE	1.62	0.40
20:T:84:LEU:HD23	20:T:84:LEU:HA	1.93	0.40
1:A:390:C:H6	1:A:390:C:O5'	2.04	0.40
1:A:427:U:C4	1:A:428:G:C6	3.10	0.40
1:A:502:G:C6	1:A:503:C:C4	3.10	0.40
1:A:775:G:H2'	1:A:776:G:O4'	2.21	0.40
1:A:946:A:H2'	1:A:947:G:C8	2.57	0.40
2:B:101:MET:O	2:B:105:PHE:HD1	2.05	0.40
7:G:44:TYR:O	7:G:47:CYS:HB2	2.22	0.40
14:N:39:LEU:HA	14:N:39:LEU:HD23	1.81	0.40
17:Q:59:ILE:HG23	17:Q:71:PHE:HB3	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	232/256 (91%)	190 (82%)	36 (16%)	6 (3%)	5	32
3	C	204/239 (85%)	163 (80%)	28 (14%)	13 (6%)	1	13
4	D	206/209 (99%)	181 (88%)	19 (9%)	6 (3%)	4	30
5	E	148/162 (91%)	135 (91%)	9 (6%)	4 (3%)	5	31
6	F	99/101 (98%)	89 (90%)	9 (9%)	1 (1%)	15	52
7	G	153/156 (98%)	135 (88%)	16 (10%)	2 (1%)	12	46
8	H	136/138 (99%)	124 (91%)	9 (7%)	3 (2%)	6	35
9	I	125/128 (98%)	101 (81%)	17 (14%)	7 (6%)	2	15

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
10	J	96/105 (91%)	69 (72%)	17 (18%)	10 (10%)	0	6
11	K	117/129 (91%)	99 (85%)	16 (14%)	2 (2%)	9	40
12	L	122/132 (92%)	96 (79%)	17 (14%)	9 (7%)	1	10
13	M	123/126 (98%)	100 (81%)	20 (16%)	3 (2%)	6	34
14	N	58/61 (95%)	46 (79%)	11 (19%)	1 (2%)	9	40
15	O	86/89 (97%)	79 (92%)	5 (6%)	2 (2%)	6	34
16	P	81/88 (92%)	70 (86%)	10 (12%)	1 (1%)	13	48
17	Q	102/105 (97%)	86 (84%)	14 (14%)	2 (2%)	7	37
18	R	71/88 (81%)	59 (83%)	10 (14%)	2 (3%)	5	31
19	S	78/93 (84%)	60 (77%)	15 (19%)	3 (4%)	3	24
20	T	97/106 (92%)	82 (84%)	8 (8%)	7 (7%)	1	10
21	V	22/27 (82%)	14 (64%)	7 (32%)	1 (4%)	2	20
All	All	2356/2538 (93%)	1978 (84%)	293 (12%)	85 (4%)	3	25

All (85) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	17	PHE
3	C	189	ALA
3	C	206	GLU
8	H	91	ARG
9	I	29	ASN
9	I	33	PHE
9	I	101	PHE
9	I	121	ARG
10	J	34	VAL
10	J	90	LEU
12	L	27	LEU
12	L	28	LYS
12	L	47	LYS
15	O	71	GLN
19	S	9	VAL
20	T	73	HIS
3	C	5	ILE
3	C	12	LEU
3	C	62	ASP
3	C	154	SER
3	C	179	ARG

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Mol	Chain	Res	Type
4	D	23	GLY
4	D	25	ARG
5	E	16	THR
7	G	4	ARG
7	G	147	ALA
9	I	38	GLN
9	I	56	LEU
10	J	39	PRO
10	J	72	VAL
10	J	86	MET
12	L	41	ARG
12	L	73	GLU
16	P	82	GLN
19	S	25	LYS
20	T	96	GLY
20	T	99	LEU
2	B	8	LYS
2	B	78	GLN
2	B	220	ASP
2	B	221	LEU
3	C	81	GLY
3	C	91	LEU
5	E	38	GLN
10	J	76	ASN
12	L	48	PRO
12	L	62	SER
13	M	86	CYS
13	M	122	LYS
14	N	24	CYS
15	O	85	LEU
17	Q	97	SER
18	R	21	LYS
18	R	87	ARG
20	T	102	GLY
21	V	3	LYS
3	C	168	ALA
4	D	4	TYR
5	E	78	HIS
9	I	127	LYS
12	L	13	LYS
13	M	13	LYS
19	S	30	LEU

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Mol	Chain	Res	Type
20	T	95	ALA
3	C	60	ALA
4	D	9	CYS
10	J	26	ALA
10	J	85	LEU
10	J	89	ASP
11	K	117	ASN
17	Q	79	SER
20	T	97	ALA
3	C	205	GLY
4	D	42	GLN
6	F	54	LYS
8	H	66	GLY
12	L	88	GLY
2	B	130	ARG
20	T	98	PRO
4	D	5	ILE
5	E	11	ILE
11	K	118	GLY
3	C	66	VAL
8	H	137	VAL
10	J	40	LEU

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	
2	B	202/220 (92%)	174 (86%)	28 (14%)	3	18
3	C	160/188 (85%)	139 (87%)	21 (13%)	4	20
4	D	180/181 (99%)	162 (90%)	18 (10%)	7	31
5	E	115/123 (94%)	103 (90%)	12 (10%)	7	29
6	F	90/90 (100%)	76 (84%)	14 (16%)	2	14
7	G	126/127 (99%)	118 (94%)	8 (6%)	18	50
8	H	119/119 (100%)	104 (87%)	15 (13%)	4	21

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
9	I	98/99 (99%)	87 (89%)	11 (11%)	6	25
10	J	87/92 (95%)	74 (85%)	13 (15%)	3	16
11	K	90/99 (91%)	83 (92%)	7 (8%)	12	41
12	L	104/109 (95%)	93 (89%)	11 (11%)	6	28
13	M	100/101 (99%)	89 (89%)	11 (11%)	6	27
14	N	49/50 (98%)	40 (82%)	9 (18%)	1	7
15	O	79/80 (99%)	70 (89%)	9 (11%)	5	24
16	P	72/74 (97%)	65 (90%)	7 (10%)	8	32
17	Q	96/97 (99%)	87 (91%)	9 (9%)	8	33
18	R	64/77 (83%)	59 (92%)	5 (8%)	12	41
19	S	71/80 (89%)	62 (87%)	9 (13%)	4	20
20	T	75/82 (92%)	69 (92%)	6 (8%)	12	40
21	V	19/22 (86%)	16 (84%)	3 (16%)	2	14
All	All	1996/2110 (95%)	1770 (89%)	226 (11%)	6	25

All (226) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	8	LYS
2	B	9	GLU
2	B	15	VAL
2	B	21	ARG
2	B	22	LYS
2	B	24	TRP
2	B	59	GLU
2	B	69	LEU
2	B	76	GLN
2	B	93	VAL
2	B	114	ARG
2	B	117	GLU
2	B	129	GLU
2	B	134	GLU
2	B	135	GLN
2	B	143	GLU
2	B	153	ARG
2	B	162	ILE
2	B	163	PHE

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Mol	Chain	Res	Type
2	B	170	GLU
2	B	172	ILE
2	B	178	ARG
2	B	185	ILE
2	B	189	ASP
2	B	190	THR
2	B	196	LEU
2	B	231	GLU
2	B	236	TYR
3	C	4	LYS
3	C	26	LYS
3	C	31	HIS
3	C	32	LEU
3	C	34	LEU
3	C	62	ASP
3	C	72	LYS
3	C	76	VAL
3	C	85	ARG
3	C	91	LEU
3	C	107	GLN
3	C	112	SER
3	C	128	PHE
3	C	162	GLN
3	C	166	GLU
3	C	167	TRP
3	C	179	ARG
3	C	188	LEU
3	C	191	THR
3	C	192	THR
3	C	204	LEU
4	D	3	ARG
4	D	8	VAL
4	D	9	CYS
4	D	15	GLU
4	D	64	LEU
4	D	78	LEU
4	D	93	PHE
4	D	96	LEU
4	D	112	VAL
4	D	122	ARG
4	D	127	THR
4	D	131	ARG

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Mol	Chain	Res	Type
4	D	151	LYS
4	D	157	LEU
4	D	161	ASN
4	D	175	SER
4	D	176	LEU
4	D	190	ASP
5	E	9	LYS
5	E	12	LEU
5	E	15	ARG
5	E	16	THR
5	E	31	LEU
5	E	41	VAL
5	E	47	LYS
5	E	80	ILE
5	E	83	GLU
5	E	89	ILE
5	E	120	THR
5	E	150	ARG
6	F	3	ARG
6	F	10	LEU
6	F	19	LEU
6	F	28	ARG
6	F	47	ARG
6	F	54	LYS
6	F	55	ASP
6	F	70	ASP
6	F	72	VAL
6	F	73	ASN
6	F	74	ASP
6	F	79	LEU
6	F	82	ARG
6	F	91	VAL
7	G	24	THR
7	G	79	ARG
7	G	80	VAL
7	G	92	SER
7	G	96	GLN
7	G	113	GLU
7	G	115	ARG
7	G	140	ASP
8	H	24	THR
8	H	29	SER

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Mol	Chain	Res	Type
8	H	39	LEU
8	H	49	GLU
8	H	50	ARG
8	H	65	TYR
8	H	85	ARG
8	H	88	LYS
8	H	91	ARG
8	H	95	VAL
8	H	104	ARG
8	H	113	SER
8	H	119	LEU
8	H	127	LEU
8	H	133	LEU
9	I	12	GLU
9	I	27	THR
9	I	28	VAL
9	I	34	ASN
9	I	38	GLN
9	I	58	ARG
9	I	104	ARG
9	I	111	ARG
9	I	113	LYS
9	I	121	ARG
9	I	127	LYS
10	J	3	LYS
10	J	4	ILE
10	J	9	ARG
10	J	60	ARG
10	J	66	ARG
10	J	71	LEU
10	J	73	ASP
10	J	75	ILE
10	J	82	ILE
10	J	84	GLN
10	J	86	MET
10	J	87	THR
10	J	90	LEU
11	K	29	ILE
11	K	34	ASP
11	K	51	LYS
11	K	54	ARG
11	K	79	SER

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Mol	Chain	Res	Type
11	K	106	LYS
11	K	127	LYS
12	L	13	LYS
12	L	20	LYS
12	L	28	LYS
12	L	33	ARG
12	L	42	THR
12	L	62	SER
12	L	66	VAL
12	L	78	GLN
12	L	79	GLU
12	L	97	ARG
12	L	126	LYS
13	M	9	ILE
13	M	13	LYS
13	M	66	LEU
13	M	70	LEU
13	M	88	ARG
13	M	98	VAL
13	M	106	ASN
13	M	110	ARG
13	M	115	LYS
13	M	125	ARG
13	M	126	LYS
14	N	7	ILE
14	N	8	GLU
14	N	9	LYS
14	N	12	ARG
14	N	22	THR
14	N	24	CYS
14	N	27	CYS
14	N	33	VAL
14	N	41	ARG
15	O	7	GLU
15	O	10	LYS
15	O	34	LEU
15	O	39	LEU
15	O	40	SER
15	O	64	ARG
15	O	70	LEU
15	O	81	LEU
15	O	88	ARG

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Mol	Chain	Res	Type
16	P	1	MET
16	P	2	VAL
16	P	45	THR
16	P	67	THR
16	P	80	PHE
16	P	81	ARG
16	P	82	GLN
17	Q	9	VAL
17	Q	52	LYS
17	Q	55	ASP
17	Q	59	ILE
17	Q	60	ILE
17	Q	68	ARG
17	Q	78	GLU
17	Q	97	SER
17	Q	100	LYS
18	R	39	VAL
18	R	42	ARG
18	R	47	THR
18	R	54	ARG
18	R	88	LYS
19	S	6	LYS
19	S	7	LYS
19	S	9	VAL
19	S	12	ASP
19	S	13	ASP
19	S	61	TYR
19	S	62	ILE
19	S	63	THR
19	S	65	ASN
20	T	13	LEU
20	T	35	THR
20	T	42	GLN
20	T	75	ASN
20	T	82	SER
20	T	84	LEU
21	V	6	ARG
21	V	8	THR
21	V	17	THR

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

Mol	Chain	Res	Type
2	B	78	GLN
3	C	6	HIS
3	C	118	GLN
7	G	37	ASN
12	L	75	HIS
15	O	37	ASN
16	P	82	GLN
19	S	14	HIS
20	T	16	HIS

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	1506/1507 (99%)	285 (18%)	46 (3%)
22	X	10/17 (58%)	2 (20%)	0
23	W	3/6 (50%)	0	0
All	All	1519/1530 (99%)	287 (18%)	46 (3%)

All (287) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	7	G
1	A	9	G
1	A	26	A
1	A	32	A
1	A	39	G
1	A	47	C
1	A	48	C
1	A	49	U
1	A	50	A
1	A	51	A
1	A	59	A
1	A	60	A
1	A	61	G
1	A	101	A
1	A	108	G
1	A	115	G
1	A	116	A
1	A	120	A
1	A	121	C
1	A	129(A)	G
1	A	130	A

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Mol	Chain	Res	Type
1	A	131	C
1	A	163	C
1	A	182	U
1	A	183	G
1	A	189	G
1	A	190(A)	C
1	A	190(B)	C
1	A	190(C)	C
1	A	190(D)	U
1	A	190(E)	U
1	A	190(F)	G
1	A	190(G)	G
1	A	195	A
1	A	197	A
1	A	202	U
1	A	203	U
1	A	204	U
1	A	216	G
1	A	220	G
1	A	244	U
1	A	247	G
1	A	251	G
1	A	252	U
1	A	258	G
1	A	266	G
1	A	267	C
1	A	279	A
1	A	289	G
1	A	300	A
1	A	321	A
1	A	328	C
1	A	329	A
1	A	330	C
1	A	332	G
1	A	345	C
1	A	347	G
1	A	351	G
1	A	352	C
1	A	353	A
1	A	354	G
1	A	356	A
1	A	367	U

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Mol	Chain	Res	Type
1	A	372	C
1	A	373	A
1	A	379	C
1	A	384	G
1	A	388	G
1	A	390	C
1	A	397	A
1	A	398	C
1	A	406	G
1	A	407	G
1	A	408	A
1	A	409	G
1	A	410	G
1	A	412	A
1	A	413	G
1	A	414	A
1	A	415	A
1	A	416	G
1	A	418	C
1	A	419	C
1	A	420	U
1	A	421	U
1	A	422	C
1	A	423	G
1	A	424	G
1	A	425	G
1	A	429	U
1	A	430	A
1	A	431	A
1	A	433	C
1	A	434	U
1	A	435	C
1	A	439	A
1	A	452	A
1	A	460	A
1	A	461	C
1	A	462	G
1	A	463	A
1	A	474	G
1	A	475	G
1	A	476	G
1	A	477	G

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Mol	Chain	Res	Type
1	A	478	A
1	A	485	G
1	A	486	U
1	A	497	A
1	A	498	U
1	A	511	C
1	A	517	G
1	A	518	C
1	A	521	G
1	A	527	G
1	A	531	U
1	A	532	A
1	A	533	A
1	A	534	U
1	A	545	C
1	A	547	A
1	A	550	G
1	A	559	A
1	A	560	U
1	A	562	C
1	A	572	A
1	A	573	A
1	A	575	G
1	A	576	G
1	A	579	G
1	A	582	U
1	A	596	C
1	A	616	G
1	A	653	A
1	A	665	A
1	A	687	A
1	A	695	A
1	A	701	C
1	A	702	A
1	A	703	G
1	A	721	G
1	A	731	G
1	A	748	C
1	A	749	C
1	A	755	G
1	A	777	A
1	A	782	A

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Mol	Chain	Res	Type
1	A	787	A
1	A	792	A
1	A	793	U
1	A	794	A
1	A	813	U
1	A	815	A
1	A	817	C
1	A	828	A
1	A	836	G
1	A	839	U
1	A	840	C
1	A	841	U
1	A	848	C
1	A	859	A
1	A	876	G
1	A	902	G
1	A	914	A
1	A	926	G
1	A	927	G
1	A	934	C
1	A	935	A
1	A	936	C
1	A	945	G
1	A	960	U
1	A	961	U
1	A	966	G
1	A	969	A
1	A	971	G
1	A	972	C
1	A	974	A
1	A	975	A
1	A	976	G
1	A	977	A
1	A	983	A
1	A	991	U
1	A	992	U
1	A	993	G
1	A	1005	A
1	A	1007	C
1	A	1024	G
1	A	1025	U
1	A	1029	C

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Mol	Chain	Res	Type
1	A	1030	C
1	A	1031	G
1	A	1032	G
1	A	1050	G
1	A	1053	G
1	A	1054	C
1	A	1055	A
1	A	1064	G
1	A	1065	U
1	A	1066	C
1	A	1068	G
1	A	1094	G
1	A	1095	U
1	A	1096	C
1	A	1101	A
1	A	1102	A
1	A	1117	G
1	A	1124	G
1	A	1125	U
1	A	1126	U
1	A	1127	G
1	A	1129	C
1	A	1130	A
1	A	1131	G
1	A	1137	C
1	A	1139	G
1	A	1145	C
1	A	1146	A
1	A	1152	A
1	A	1157	A
1	A	1159	U
1	A	1171	G
1	A	1183	A
1	A	1184	G
1	A	1196	U
1	A	1197	G
1	A	1201	A
1	A	1202	G
1	A	1211	U
1	A	1212	U
1	A	1213	A
1	A	1214	C

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Mol	Chain	Res	Type
1	A	1225	A
1	A	1227	A
1	A	1238	A
1	A	1256	A
1	A	1257	U
1	A	1258	G
1	A	1270	C
1	A	1280	A
1	A	1281	U
1	A	1282	C
1	A	1286	A
1	A	1287	A
1	A	1300	G
1	A	1301	U
1	A	1302	U
1	A	1305	G
1	A	1319	A
1	A	1340	A
1	A	1346	A
1	A	1347	G
1	A	1348	U
1	A	1353	G
1	A	1360	A
1	A	1361	G
1	A	1362	C
1	A	1363	A
1	A	1364	U
1	A	1368	G
1	A	1370	G
1	A	1397	C
1	A	1398	A
1	A	1442	G
1	A	1443	G
1	A	1446	A
1	A	1447	G
1	A	1452	C
1	A	1487	G
1	A	1492	A
1	A	1494	G
1	A	1499	A
1	A	1502	A
1	A	1503	A

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Mol	Chain	Res	Type
1	A	1504	G
1	A	1505	G
1	A	1506	U
1	A	1517	G
1	A	1519	A
1	A	1520	G
1	A	1521	G
1	A	1529	G
1	A	1530	G
1	A	1533	C
1	A	1534	A
1	A	1539	C
22	X	32	C
22	X	33	U

All (46) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A	60	A
1	A	108	G
1	A	115	G
1	A	129(A)	G
1	A	181	G
1	A	243	A
1	A	244	U
1	A	250	A
1	A	251	G
1	A	266	G
1	A	328	C
1	A	372	C
1	A	407	G
1	A	428	G
1	A	429	U
1	A	484	G
1	A	485	G
1	A	533	A
1	A	559	A
1	A	575	G
1	A	701	C
1	A	812	C
1	A	840	C
1	A	960	U

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Mol	Chain	Res	Type
1	A	965	A
1	A	975	A
1	A	1049	U
1	A	1065	U
1	A	1067	A
1	A	1101	A
1	A	1126	U
1	A	1145	C
1	A	1182	G
1	A	1183	A
1	A	1201	A
1	A	1212	U
1	A	1281	U
1	A	1285	A
1	A	1300	G
1	A	1346	A
1	A	1347	G
1	A	1397	C
1	A	1498	U
1	A	1503	A
1	A	1505	G
1	A	1532	U

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 79 ligands modelled in this entry, 78 are monoatomic - leaving 1 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	$\# Z > 2$
24	PAR	A	1601	-	45,45,45	1.27	6 (13%)	64,67,67	1.67	13 (20%)

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
24	PAR	A	1601	-	-	6/18/94/94	0/4/4/4

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
24	A	1601	PAR	C64-C54	3.80	1.57	1.52
24	A	1601	PAR	C13-C23	2.78	1.56	1.52
24	A	1601	PAR	O54-C14	2.48	1.48	1.41
24	A	1601	PAR	C52-C42	2.42	1.57	1.52
24	A	1601	PAR	O43-C13	2.11	1.45	1.41
24	A	1601	PAR	O33-C14	2.02	1.47	1.41

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
24	A	1601	PAR	O33-C14-C24	6.71	119.77	108.22
24	A	1601	PAR	C34-C24-N24	-3.46	103.97	111.05
24	A	1601	PAR	O52-C13-C23	3.22	114.62	107.96
24	A	1601	PAR	C14-O33-C33	-2.96	110.64	117.96
24	A	1601	PAR	O34-C34-C44	-2.83	103.81	110.35
24	A	1601	PAR	O11-C11-O51	2.78	118.45	110.67
24	A	1601	PAR	O52-C13-O43	-2.77	108.43	111.43
24	A	1601	PAR	C11-O51-C51	2.46	118.51	113.69
24	A	1601	PAR	O54-C54-C44	2.44	114.12	109.69
24	A	1601	PAR	O51-C51-C61	2.28	112.10	106.44
24	A	1601	PAR	C13-O52-C52	-2.27	112.35	117.96
24	A	1601	PAR	C13-C23-C33	2.25	104.81	102.10
24	A	1601	PAR	O34-C34-C24	-2.02	106.59	110.22

There are no chirality outliers.

All (6) torsion outliers are listed below:

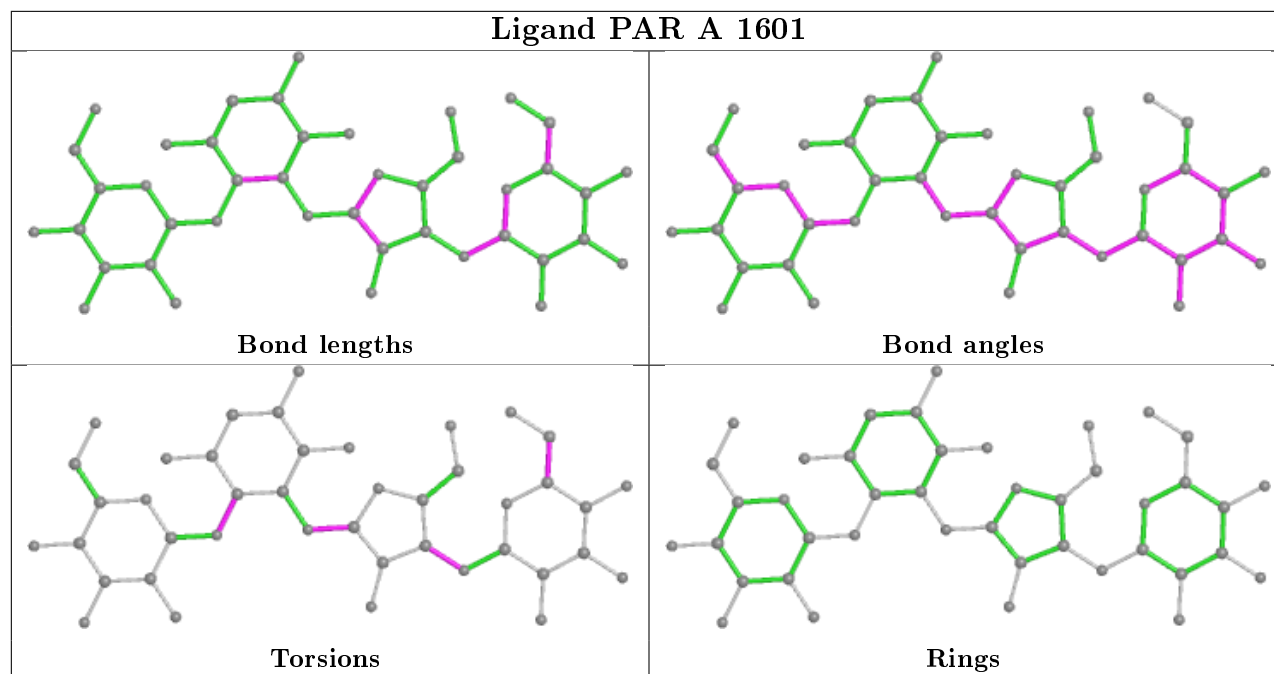
Mol	Chain	Res	Type	Atoms
24	A	1601	PAR	O54-C54-C64-N64
24	A	1601	PAR	C52-C42-O11-C11
24	A	1601	PAR	O43-C13-O52-C52
24	A	1601	PAR	C23-C13-O52-C52
24	A	1601	PAR	C23-C33-O33-C14
24	A	1601	PAR	C44-C54-C64-N64

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
24	A	1601	PAR	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	1507/1507 (100%)	0.06	27 (1%) 68 65	90, 129, 200, 303	0
2	B	234/256 (91%)	0.44	19 (8%) 12 14	125, 164, 195, 207	0
3	C	206/239 (86%)	0.36	16 (7%) 13 15	130, 155, 205, 213	0
4	D	208/209 (99%)	0.47	13 (6%) 20 20	119, 138, 166, 173	0
5	E	150/162 (92%)	0.01	2 (1%) 77 73	102, 119, 134, 148	0
6	F	101/101 (100%)	0.28	4 (3%) 38 36	131, 154, 174, 183	0
7	G	155/156 (99%)	0.29	7 (4%) 33 32	131, 156, 178, 183	0
8	H	138/138 (100%)	0.22	4 (2%) 51 49	95, 115, 132, 145	0
9	I	127/128 (99%)	0.76	20 (15%) 2 3	117, 167, 187, 192	0
10	J	98/105 (93%)	1.45	30 (30%) 0 0	126, 202, 253, 257	0
11	K	119/129 (92%)	0.63	12 (10%) 7 9	109, 134, 154, 176	0
12	L	124/132 (93%)	0.53	8 (6%) 18 19	92, 128, 145, 159	0
13	M	125/126 (99%)	1.31	32 (25%) 0 0	130, 162, 177, 185	0
14	N	60/61 (98%)	0.39	2 (3%) 46 44	136, 147, 178, 189	0
15	O	88/89 (98%)	0.60	7 (7%) 12 14	106, 130, 155, 177	0
16	P	83/88 (94%)	0.68	3 (3%) 42 41	105, 117, 141, 183	0
17	Q	104/105 (99%)	0.81	9 (8%) 10 13	96, 120, 142, 150	0
18	R	73/88 (82%)	0.45	6 (8%) 11 14	121, 138, 173, 184	0
19	S	80/93 (86%)	0.94	15 (18%) 1 2	152, 187, 204, 208	0
20	T	99/106 (93%)	0.52	6 (6%) 21 22	114, 132, 160, 162	0
21	V	24/27 (88%)	1.03	5 (20%) 1 1	132, 151, 161, 170	0
22	X	10/17 (58%)	0.49	1 (10%) 7 9	151, 181, 199, 205	0
23	W	4/6 (66%)	0.08	0 100 100	140, 140, 143, 144	0
All	All	3917/4068 (96%)	0.36	248 (6%) 20 20	90, 138, 197, 303	0

All (248) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	416	G	13.6
13	M	123	ALA	12.9
13	M	124	PRO	11.2
17	Q	103	GLY	10.8
1	A	422	C	9.5
1	A	414	A	7.9
1	A	415	A	7.9
17	Q	104	LYS	7.4
10	J	74	ILE	7.3
17	Q	102	GLY	7.1
13	M	122	LYS	7.0
1	A	413	G	6.7
11	K	129	SER	6.3
13	M	126	LYS	6.1
10	J	73	ASP	6.0
1	A	1534	A	5.9
17	Q	105	ALA	5.7
1	A	423	G	5.6
11	K	128	ALA	5.6
10	J	34	VAL	5.5
3	C	196	LEU	5.4
9	I	15	ALA	5.2
13	M	7	VAL	5.2
10	J	72	VAL	5.1
15	O	21	ASP	5.1
4	D	42	GLN	5.0
1	A	421	U	5.0
13	M	121	LYS	4.8
13	M	125	ARG	4.8
1	A	1533	C	4.8
9	I	7	THR	4.8
10	J	75	ILE	4.7
18	R	88	LYS	4.7
2	B	238	LEU	4.7
13	M	120	LYS	4.7
15	O	22	THR	4.6
19	S	32	LYS	4.5
4	D	47	ARG	4.3
9	I	8	GLY	4.3
10	J	8	LEU	4.3
19	S	30	LEU	4.2
10	J	70	ARG	4.2

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Mol	Chain	Res	Type	RSRZ
9	I	128	ARG	4.2
1	A	1006	C	4.1
10	J	4	ILE	4.1
10	J	35	SER	4.0
13	M	11	ARG	4.0
21	V	24	ARG	3.9
4	D	37	PRO	3.9
10	J	36	GLY	3.8
10	J	26	ALA	3.7
1	A	1002	G	3.7
1	A	1034	G	3.7
1	A	1038	C	3.7
4	D	181	MET	3.6
9	I	65	VAL	3.6
19	S	72	GLY	3.6
9	I	64	THR	3.5
11	K	127	LYS	3.5
4	D	41	GLY	3.5
22	X	33	U	3.5
11	K	12	ARG	3.5
2	B	239	VAL	3.4
10	J	71	LEU	3.4
13	M	118	ALA	3.4
13	M	12	ASN	3.4
3	C	102	ASN	3.4
19	S	34	TRP	3.4
11	K	77	MET	3.4
7	G	81	GLY	3.3
2	B	125	PRO	3.3
13	M	6	GLY	3.3
7	G	82	GLY	3.3
3	C	103	VAL	3.3
3	C	101	LEU	3.3
9	I	36	TYR	3.3
18	R	17	SER	3.2
13	M	4	ILE	3.2
15	O	42	HIS	3.2
2	B	16	HIS	3.2
19	S	50	ALA	3.2
4	D	8	VAL	3.1
17	Q	101	ARG	3.1
15	O	17	ARG	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	1039	C	3.1
9	I	102	LEU	3.1
19	S	69	HIS	3.1
13	M	119	GLY	3.0
4	D	45	GLN	3.0
1	A	412	A	3.0
10	J	80	LYS	3.0
21	V	18	TYR	3.0
10	J	64	GLU	2.9
10	J	32	ALA	2.9
13	M	45	VAL	2.9
13	M	2	ALA	2.9
1	A	1003	G	2.9
2	B	132	LYS	2.9
15	O	49	ASP	2.9
1	A	1001	A	2.8
11	K	11	LYS	2.8
7	G	124	LEU	2.8
1	A	1040	U	2.8
5	E	45	PHE	2.8
13	M	117	VAL	2.8
18	R	29	PHE	2.8
13	M	8	GLU	2.8
2	B	131	PRO	2.8
13	M	10	PRO	2.8
19	S	29	ARG	2.8
10	J	77	PRO	2.8
11	K	13	GLN	2.8
19	S	2	PRO	2.8
19	S	36	ARG	2.8
10	J	83	GLU	2.7
4	D	38	TYR	2.7
13	M	60	VAL	2.7
20	T	104	LEU	2.7
9	I	37	PHE	2.7
12	L	50	SER	2.7
10	J	37	PRO	2.7
19	S	71	LEU	2.7
1	A	432	A	2.7
21	V	21	TYR	2.7
2	B	130	ARG	2.7
19	S	31	ILE	2.7

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Mol	Chain	Res	Type	RSRZ
19	S	3	ARG	2.6
8	H	129	VAL	2.6
9	I	61	ALA	2.6
2	B	206	ASP	2.6
2	B	192	SER	2.6
11	K	42	TRP	2.6
10	J	100	THR	2.6
4	D	13	ARG	2.6
13	M	3	ARG	2.6
2	B	31	TYR	2.6
3	C	42	LEU	2.6
19	S	81	ARG	2.6
1	A	202	U	2.5
9	I	26	VAL	2.5
20	T	64	ASP	2.5
21	V	5	ASP	2.5
1	A	1519	A	2.5
3	C	100	ALA	2.5
3	C	161	GLU	2.5
3	C	197	GLY	2.5
3	C	54	ARG	2.5
10	J	3	LYS	2.5
4	D	43	HIS	2.5
16	P	29	ASP	2.5
10	J	96	ILE	2.5
2	B	202	PRO	2.5
20	T	100	ILE	2.5
9	I	60	ASP	2.5
2	B	135	GLN	2.5
2	B	35	GLU	2.5
17	Q	44	ALA	2.4
10	J	5	ARG	2.4
1	A	1539	C	2.4
6	F	61	LEU	2.4
11	K	14	VAL	2.4
10	J	63	PHE	2.4
13	M	27	LYS	2.4
13	M	102	ARG	2.4
1	A	1007	C	2.4
2	B	190	THR	2.4
10	J	45	ARG	2.4
2	B	230	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
13	M	24	GLY	2.4
12	L	127	GLU	2.4
13	M	87	TYR	2.4
16	P	19	ILE	2.4
21	V	22	ARG	2.4
4	D	46	LYS	2.3
13	M	30	ALA	2.3
3	C	193	TYR	2.3
5	E	17	ALA	2.3
20	T	103	GLY	2.3
9	I	16	ARG	2.3
9	I	93	ARG	2.3
10	J	33	GLN	2.3
7	G	156	TRP	2.3
3	C	53	ALA	2.3
3	C	201	TYR	2.3
13	M	59	TYR	2.3
12	L	112	ASP	2.3
9	I	63	ILE	2.3
19	S	18	LYS	2.3
4	D	36	ARG	2.3
17	Q	98	LEU	2.3
1	A	1005	A	2.3
17	Q	19	VAL	2.3
12	L	28	LYS	2.2
14	N	29	ARG	2.2
3	C	52	LEU	2.2
3	C	156	ARG	2.2
13	M	5	ALA	2.2
13	M	13	LYS	2.2
9	I	33	PHE	2.2
17	Q	58	GLU	2.2
10	J	69	ASN	2.2
8	H	1	MET	2.2
12	L	31	PRO	2.2
10	J	38	ILE	2.2
12	L	96	VAL	2.2
9	I	88	TYR	2.2
11	K	50	TYR	2.2
2	B	22	LYS	2.2
6	F	14	LEU	2.2
7	G	103	TRP	2.1

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Mol	Chain	Res	Type	RSRZ
10	J	18	ALA	2.1
18	R	16	PRO	2.1
7	G	134	ALA	2.1
9	I	25	LYS	2.1
1	A	1540	U	2.1
8	H	97	VAL	2.1
14	N	30	ALA	2.1
18	R	18	ARG	2.1
18	R	43	PHE	2.1
4	D	30	LYS	2.1
1	A	1531	A	2.1
1	A	1008	C	2.1
11	K	40	ILE	2.1
2	B	72	GLY	2.1
6	F	67	MET	2.1
3	C	68	VAL	2.1
15	O	52	SER	2.1
2	B	138	LEU	2.1
13	M	19	LEU	2.1
19	S	74	PHE	2.1
6	F	4	TYR	2.1
9	I	46	ALA	2.1
10	J	61	GLU	2.1
16	P	39	TYR	2.1
20	T	36	LEU	2.1
10	J	23	ILE	2.1
15	O	23	GLY	2.1
20	T	56	MET	2.1
11	K	126	ARG	2.1
13	M	111	LYS	2.1
9	I	81	ILE	2.0
8	H	128	GLY	2.0
13	M	113	PRO	2.0
3	C	195	VAL	2.0
2	B	142	LEU	2.0
7	G	128	ALA	2.0
12	L	95	GLY	2.0
12	L	32	PHE	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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(Å ²)	Q<0.9
25	MG	A	1610	1/1	0.11	0.37	116,116,116,116	0
25	MG	A	1667	1/1	0.17	0.49	156,156,156,156	0
25	MG	A	1624	1/1	0.19	0.96	109,109,109,109	0
25	MG	A	1671	1/1	0.50	1.07	117,117,117,117	0
25	MG	A	1604	1/1	0.52	0.47	116,116,116,116	0
25	MG	A	1664	1/1	0.56	0.57	136,136,136,136	0
25	MG	A	1614	1/1	0.57	1.08	99,99,99,99	0
25	MG	E	201	1/1	0.57	1.06	83,83,83,83	0
25	MG	A	1602	1/1	0.58	0.48	112,112,112,112	0
25	MG	A	1629	1/1	0.59	0.86	98,98,98,98	0
25	MG	A	1662	1/1	0.59	0.99	99,99,99,99	0
25	MG	A	1605	1/1	0.66	0.15	123,123,123,123	0
25	MG	A	1643	1/1	0.66	0.50	91,91,91,91	0
25	MG	A	1647	1/1	0.69	0.34	134,134,134,134	0
25	MG	A	1606	1/1	0.70	1.19	95,95,95,95	0
25	MG	A	1672	1/1	0.71	0.19	116,116,116,116	0
25	MG	A	1615	1/1	0.73	0.38	143,143,143,143	0
25	MG	A	1618	1/1	0.73	0.33	100,100,100,100	0
25	MG	A	1644	1/1	0.73	0.43	100,100,100,100	0
25	MG	A	1619	1/1	0.74	0.55	118,118,118,118	0
25	MG	A	1632	1/1	0.77	1.25	122,122,122,122	0
25	MG	A	1675	1/1	0.78	0.58	105,105,105,105	0
25	MG	A	1674	1/1	0.78	0.36	118,118,118,118	0
25	MG	A	1663	1/1	0.78	0.65	99,99,99,99	0
25	MG	A	1609	1/1	0.79	0.45	110,110,110,110	0
25	MG	A	1666	1/1	0.81	0.70	115,115,115,115	0
25	MG	A	1636	1/1	0.83	0.42	111,111,111,111	0
25	MG	A	1638	1/1	0.84	0.20	109,109,109,109	0
25	MG	A	1665	1/1	0.84	0.58	112,112,112,112	0
25	MG	A	1621	1/1	0.85	0.33	104,104,104,104	0
25	MG	A	1627	1/1	0.85	0.37	101,101,101,101	0

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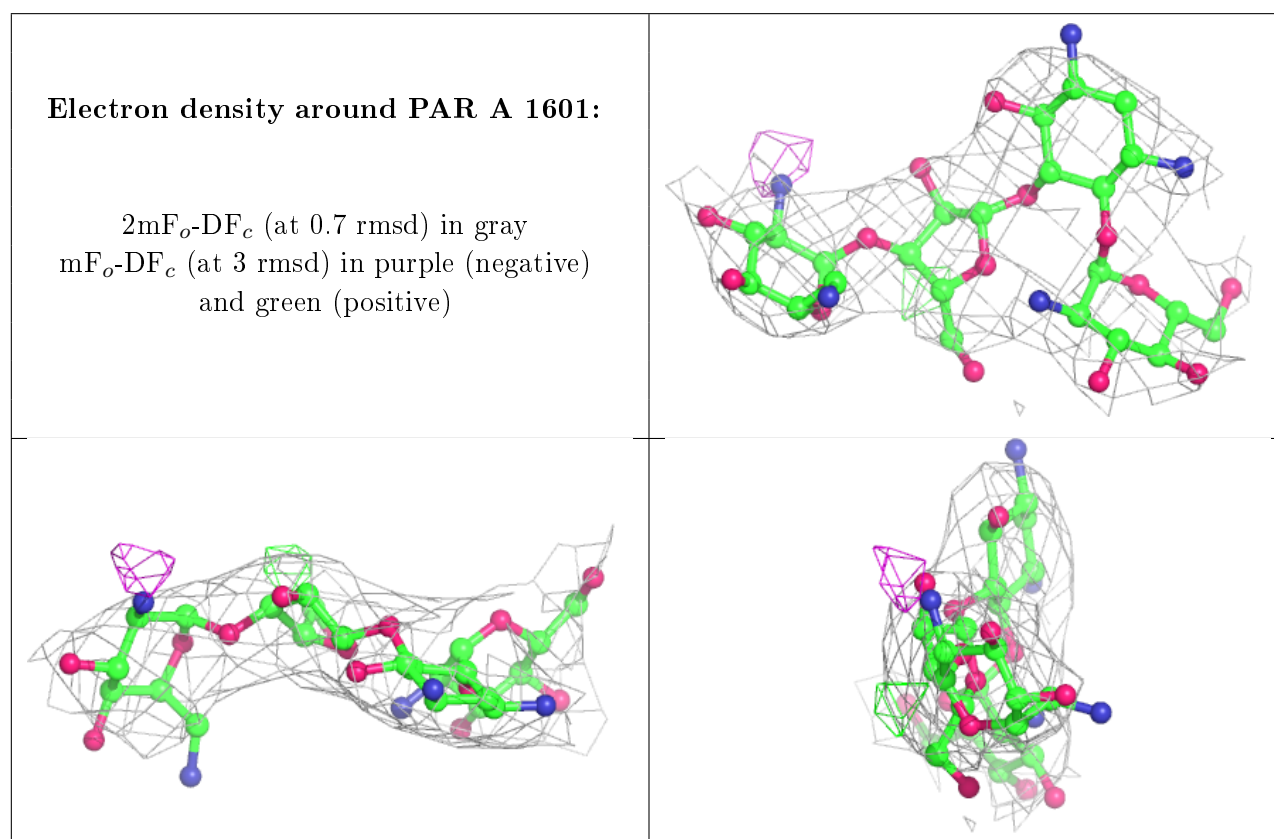
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
25	MG	A	1613	1/1	0.86	0.72	116,116,116,116	0
25	MG	A	1657	1/1	0.87	0.18	90,90,90,90	0
25	MG	J	201	1/1	0.87	0.28	115,115,115,115	0
25	MG	A	1661	1/1	0.87	0.40	113,113,113,113	0
25	MG	A	1630	1/1	0.87	1.41	108,108,108,108	0
25	MG	A	1603	1/1	0.88	1.35	118,118,118,118	0
25	MG	A	1625	1/1	0.89	0.56	90,90,90,90	0
25	MG	A	1642	1/1	0.90	0.41	127,127,127,127	0
25	MG	A	1616	1/1	0.91	0.50	82,82,82,82	0
25	MG	A	1670	1/1	0.92	0.36	122,122,122,122	0
25	MG	A	1651	1/1	0.92	1.12	91,91,91,91	0
24	PAR	A	1601	42/42	0.92	0.23	118,122,126,127	0
25	MG	A	1607	1/1	0.92	0.24	125,125,125,125	0
25	MG	A	1608	1/1	0.92	0.20	116,116,116,116	0
25	MG	A	1655	1/1	0.92	1.49	108,108,108,108	0
25	MG	A	1641	1/1	0.92	0.56	139,139,139,139	0
25	MG	A	1652	1/1	0.93	0.20	114,114,114,114	0
25	MG	A	1617	1/1	0.93	0.41	90,90,90,90	0
25	MG	A	1654	1/1	0.93	0.25	94,94,94,94	0
25	MG	A	1659	1/1	0.94	0.25	112,112,112,112	0
25	MG	A	1677	1/1	0.94	0.13	130,130,130,130	0
25	MG	A	1626	1/1	0.94	0.15	114,114,114,114	0
25	MG	A	1623	1/1	0.94	0.31	86,86,86,86	0
25	MG	A	1658	1/1	0.94	0.38	113,113,113,113	0
25	MG	A	1628	1/1	0.95	0.67	108,108,108,108	0
25	MG	A	1639	1/1	0.95	0.69	113,113,113,113	0
25	MG	A	1673	1/1	0.95	0.31	107,107,107,107	0
25	MG	A	1637	1/1	0.95	0.31	82,82,82,82	0
25	MG	A	1648	1/1	0.95	0.30	122,122,122,122	0
25	MG	A	1611	1/1	0.95	0.42	104,104,104,104	0
25	MG	A	1646	1/1	0.96	0.24	100,100,100,100	0
25	MG	A	1660	1/1	0.96	0.20	136,136,136,136	0
25	MG	A	1633	1/1	0.97	0.37	118,118,118,118	0
25	MG	A	1640	1/1	0.97	0.41	100,100,100,100	0
25	MG	A	1631	1/1	0.97	0.49	110,110,110,110	0
25	MG	A	1634	1/1	0.97	0.42	113,113,113,113	0
25	MG	A	1669	1/1	0.97	0.71	110,110,110,110	0
25	MG	A	1653	1/1	0.97	0.62	121,121,121,121	0
25	MG	A	1668	1/1	0.98	0.29	135,135,135,135	0
25	MG	A	1656	1/1	0.98	0.23	127,127,127,127	0
25	MG	A	1645	1/1	0.98	0.22	92,92,92,92	0
25	MG	A	1676	1/1	0.98	0.39	97,97,97,97	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
25	MG	A	1635	1/1	0.98	0.31	90,90,90,90	0
25	MG	A	1612	1/1	0.98	0.63	114,114,114,114	0
25	MG	A	1650	1/1	0.98	0.60	96,96,96,96	0
25	MG	A	1620	1/1	0.98	0.44	81,81,81,81	0
25	MG	A	1649	1/1	0.98	0.21	92,92,92,92	0
25	MG	A	1622	1/1	0.99	0.43	106,106,106,106	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers ⓘ

There are no such residues in this entry.