



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 12:33 PM UTC

PDB ID : 6MPI / pdb_00006mpi
Title : Structure of the *Thermus thermophilus* 30S ribosomal subunit complexed with a 2-thiocytidine (s2C32) and inosine (I34) modified anticodon stem loop (ASL) of *Escherichia coli* transfer RNA Arginine 1 (TRNAARG1) 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 : 2018-10-07
Resolution : 3.33 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

The following versions of software and data (see [references](#) ①) were used in the production of this report:

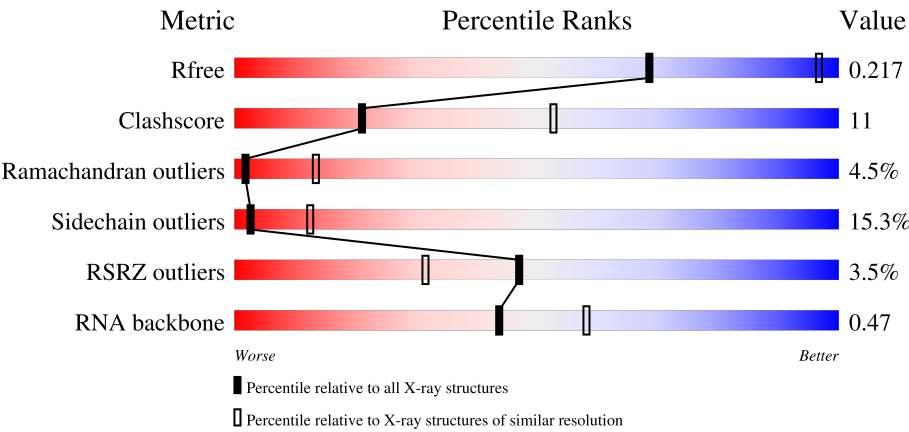
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.33 Å.

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



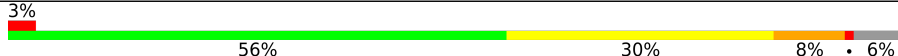
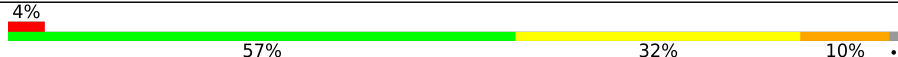
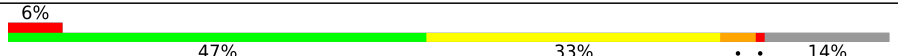
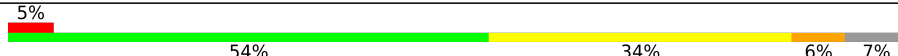
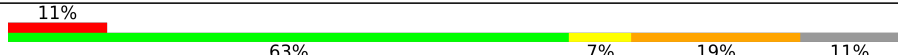
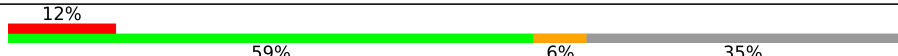
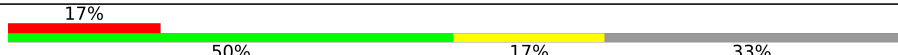
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	1434 (3.38-3.30)
Clashscore	190562	1479 (3.38-3.30)
Ramachandran outliers	187476	1456 (3.38-3.30)
Sidechain outliers	187428	1455 (3.38-3.30)
RSRZ outliers	180081	1434 (3.38-3.30)
RNA backbone	3983	1083 (3.68-3.00)

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

Mol	Chain	Length	Quality of chain
1	A	1507	2% 56% 33% 9% .
2	B	256	2% 48% 36% 7% 9%
3	C	239	3% 50% 27% 8% . 14%
4	D	209	4% 62% 33% 5%

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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
24	MG	A	1634	-	-	-	X
24	MG	A	1653	-	-	-	X
24	MG	A	1680	-	-	-	X

2 Entry composition

There are 26 unique types of molecules in this entry. The entry contains 52062 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 Arg1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
22	X	11	Total	C	N	O	P S	0	0	0
			232	105	43	73	10 1			

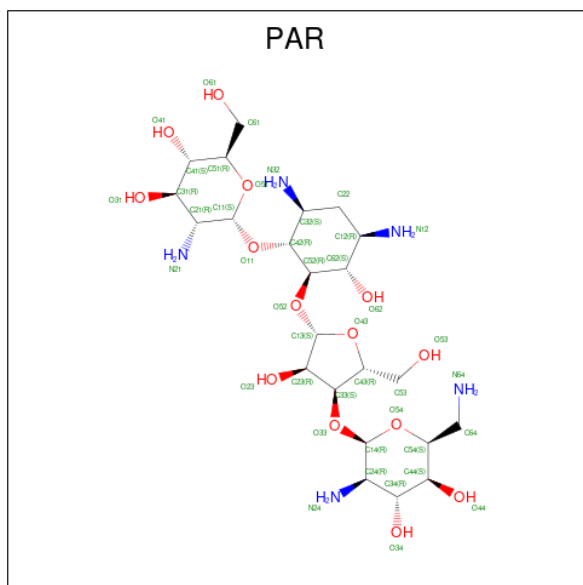
- Molecule 23 is a RNA chain called mRNA A-site fragment.

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 MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
24	A	93	Total	Mg	0	0
			93	93		
24	J	1	Total	Mg	0	0
			1	1		
24	N	1	Total	Mg	0	0
			1	1		
24	W	1	Total	Mg	0	0
			1	1		

- Molecule 25 is PAROMOMYCIN (CCD ID: PAR) (formula: C₂₃H₄₅N₅O₁₄).



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

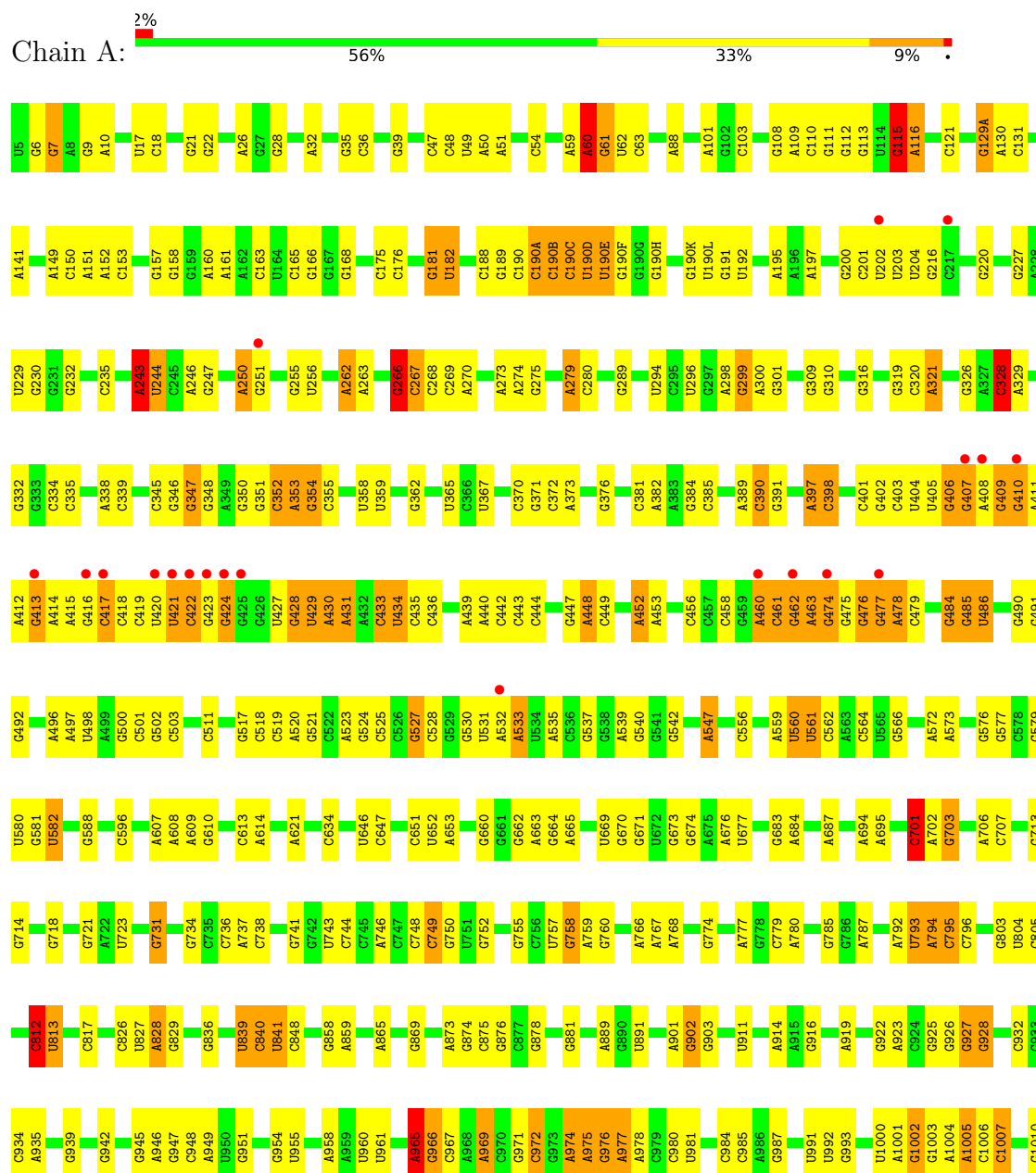
- Molecule 26 is ZINC ION (CCD ID: ZN) (formula: Zn).

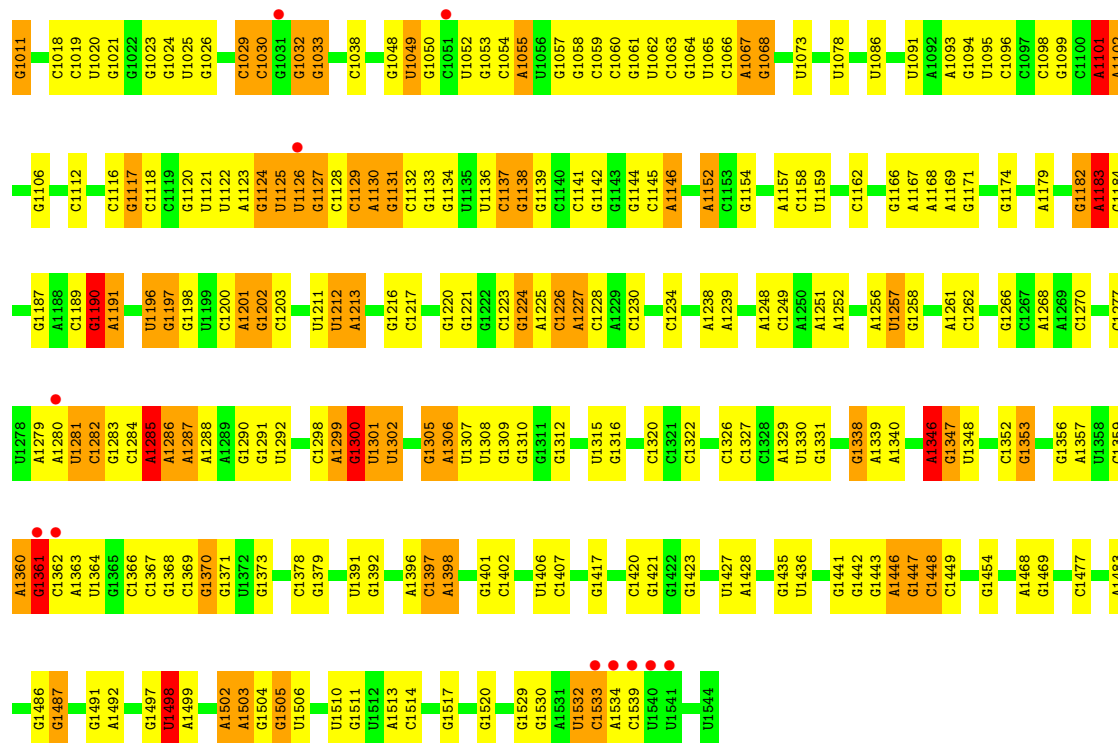
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
26	D	1	Total 1	Zn 1	0	0

3 Residue-property plots

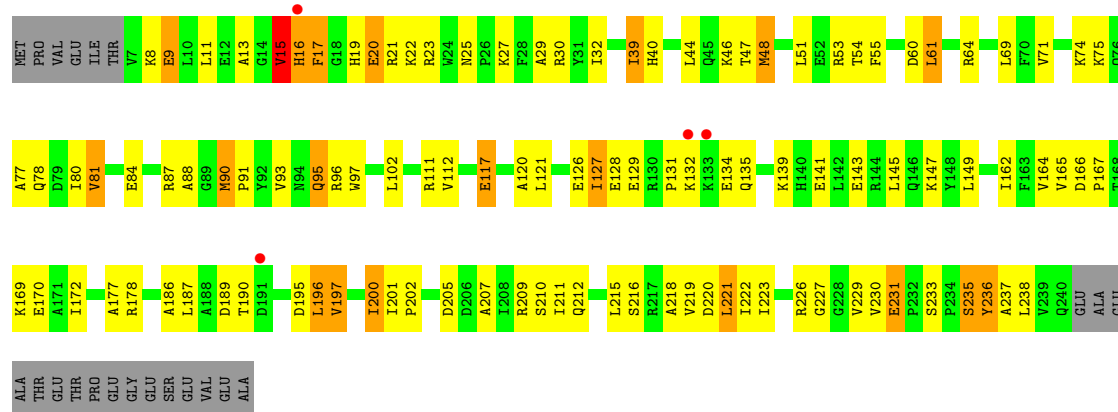
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

• Molecule 1: 16S rRNA

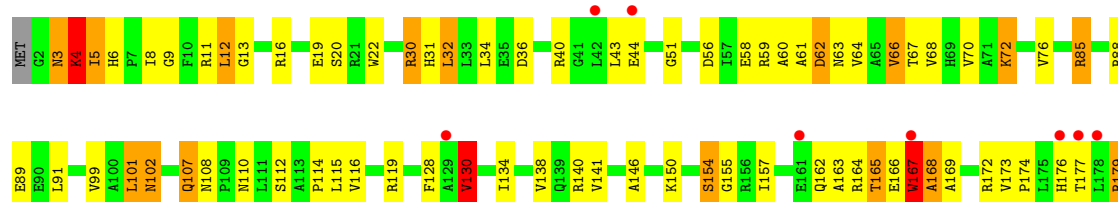


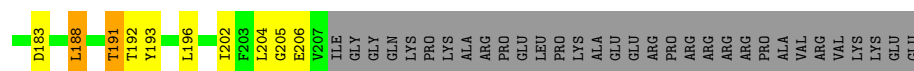


• Molecule 2: 30S ribosomal protein S2



• Molecule 3: 30S ribosomal protein S3





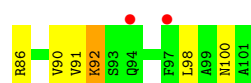
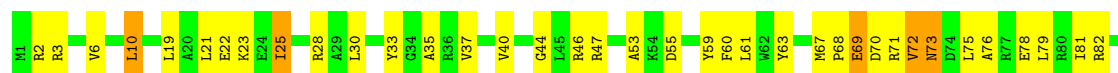
• Molecule 4: 30S ribosomal protein S4



• Molecule 5: 30S ribosomal protein S5



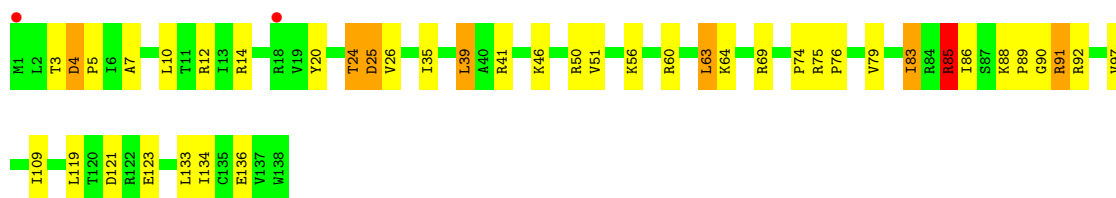
• Molecule 6: 30S ribosomal protein S6



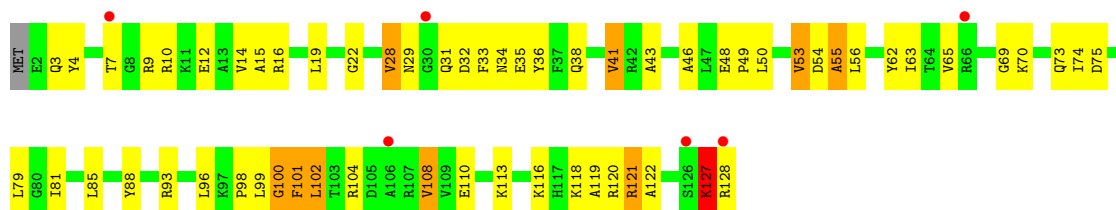
• Molecule 7: 30S ribosomal protein S7



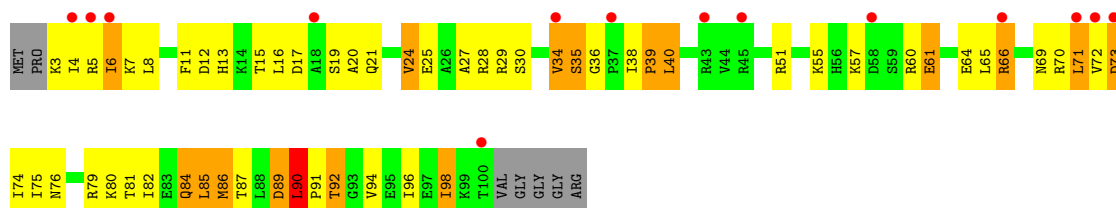
• Molecule 8: 30S ribosomal protein S8



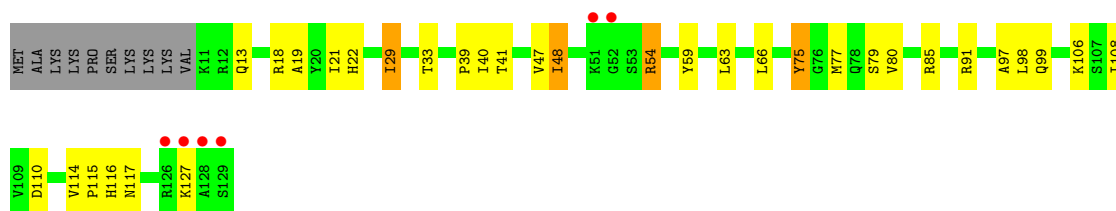
• Molecule 9: 30S ribosomal protein S9



• Molecule 10: 30S ribosomal protein S10



• Molecule 11: 30S ribosomal protein S11

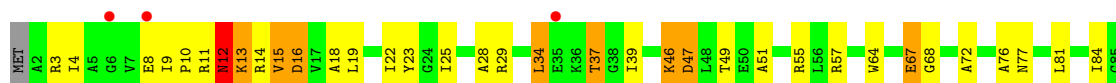


• Molecule 12: 30S ribosomal protein S12

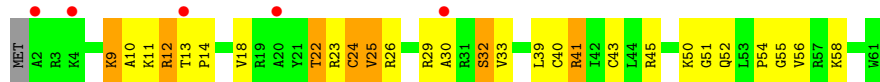




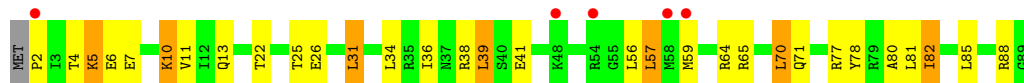
- Molecule 13: 30S ribosomal protein S13



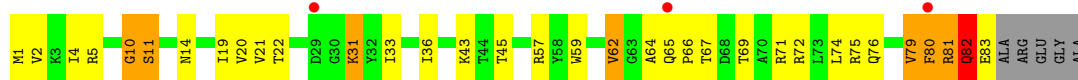
- Molecule 14: 30S ribosomal protein S14 type Z



- Molecule 15: 30S ribosomal protein S15



- Molecule 16: 30S ribosomal protein S16

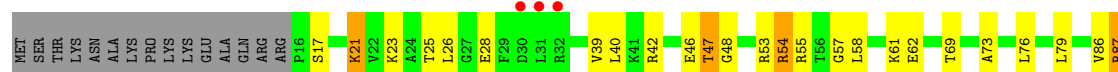


- Molecule 17: 30S ribosomal protein S17



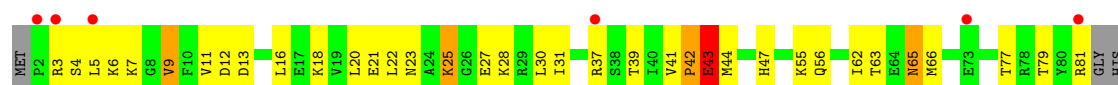
- Molecule 18: 30S ribosomal protein S18





K88

- Molecule 19: 30S ribosomal protein S19

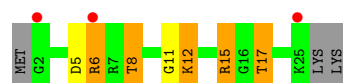


GLY
LYS
GLU
ALA
LYS
ALA
LYS
THR
LYS
LYS

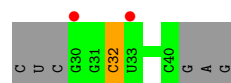
- Molecule 20: 30S ribosomal protein S20



- Molecule 21: 30S ribosomal protein Thx



- Molecule 22: tRNA ASL Escherichia coli Arg1



- Molecule 23: mRNA A-site fragment



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	403.96Å 403.96Å 176.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.77 – 3.33 49.77 – 3.33	Depositor EDS
% Data completeness (in resolution range)	86.7 (49.77-3.33) 83.0 (49.77-3.33)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.34 (at 3.33Å)	Xtriage
Refinement program	PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.190 , 0.216 0.193 , 0.217	Depositor DCC
R_{free} test set	1739 reflections (0.82%)	wwPDB-VP
Wilson B-factor (Å ²)	115.8	Xtriage
Anisotropy	0.101	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 87.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	52062	wwPDB-VP
Average B, all atoms (Å ²)	124.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.97% 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: MG, ZN, RSP, PAR

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	0.42	1/36244 (0.0%)	0.63	26/56567 (0.0%)
2	B	0.51	0/1935	0.93	5/2609 (0.2%)
3	C	0.58	0/1636	0.94	2/2205 (0.1%)
4	D	0.55	0/1733	0.95	2/2318 (0.1%)
5	E	0.69	0/1162	1.13	6/1564 (0.4%)
6	F	0.54	0/856	0.95	2/1154 (0.2%)
7	G	0.49	0/1276	0.87	2/1709 (0.1%)
8	H	0.68	0/1136	1.08	7/1527 (0.5%)
9	I	0.52	0/1029	0.92	1/1378 (0.1%)
10	J	0.58	0/805	0.96	0/1082
11	K	0.60	0/900	0.96	1/1213 (0.1%)
12	L	0.66	0/986	1.13	7/1320 (0.5%)
13	M	0.56	0/1008	1.00	4/1347 (0.3%)
14	N	0.59	0/501	1.00	0/664
15	O	0.59	0/745	0.89	0/992
16	P	0.69	0/716	1.02	1/963 (0.1%)
17	Q	0.65	0/870	1.02	3/1159 (0.3%)
18	R	0.54	0/603	0.91	0/799
19	S	0.47	0/661	0.94	2/890 (0.2%)
20	T	0.65	0/764	1.02	3/1006 (0.3%)
21	V	0.56	0/212	0.90	1/277 (0.4%)
22	X	0.65	0/235	0.64	0/360
23	W	0.36	0/91	0.53	0/140
All	All	0.48	1/56104 (0.0%)	0.76	75/83243 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
8	H	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1361	G	O3'-P	5.50	1.69	1.61

All (75) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	F	37	VAL	N-CA-C	9.54	116.42	106.21
20	T	13	LEU	N-CA-C	-8.48	102.69	113.72
2	B	16	HIS	CB-CA-C	-8.45	106.83	116.63
13	M	15	VAL	N-CA-C	8.33	119.10	110.36
12	L	93	LEU	CA-C-N	-8.13	112.34	120.31
12	L	93	LEU	C-N-CA	-8.13	112.34	120.31
12	L	47	LYS	CA-C-N	-7.89	109.97	119.84
12	L	47	LYS	C-N-CA	-7.89	109.97	119.84
8	H	134	ILE	N-CA-C	7.74	117.80	110.53
2	B	90	MET	CA-C-N	7.28	127.44	120.31
2	B	90	MET	C-N-CA	7.28	127.44	120.31
13	M	98	VAL	N-CA-C	-7.06	107.00	113.71
8	H	4	ASP	CA-C-N	6.96	126.44	118.85
8	H	4	ASP	C-N-CA	6.96	126.44	118.85
5	E	48	ALA	CA-C-N	6.86	128.41	119.84
5	E	48	ALA	C-N-CA	6.86	128.41	119.84
1	A	758	G	N1-C6-O6	6.81	140.33	119.90
1	A	965	A	C2'-C3'-O3'	6.79	119.68	109.50
7	G	87	VAL	N-CA-C	6.78	115.36	107.77
20	T	97	ALA	CA-C-N	6.77	128.31	119.84
20	T	97	ALA	C-N-CA	6.77	128.31	119.84
1	A	1498	U	C2'-C3'-O3'	6.75	119.62	109.50
5	E	23	GLY	N-CA-C	6.50	121.34	110.56
12	L	44	THR	CA-C-N	6.36	127.79	120.98
12	L	44	THR	C-N-CA	6.36	127.79	120.98
13	M	67	GLU	CB-CA-C	-6.26	109.36	116.63
1	A	1285	A	P-O3'-C3'	6.22	129.53	120.20
1	A	1498	U	P-O3'-C3'	6.19	129.48	120.20
8	H	56	LYS	CA-C-N	6.06	127.68	120.23
8	H	56	LYS	C-N-CA	6.06	127.68	120.23
1	A	60	A	P-O3'-C3'	6.01	129.21	120.20
1	A	243	A	P-O3'-C3'	5.98	129.17	120.20
2	B	16	HIS	N-CA-C	5.91	118.06	108.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	L	29	GLY	N-CA-C	-5.82	104.98	113.48
1	A	1503	A	P-O3'-C3'	5.72	128.78	120.20
1	A	1346	A	C2'-C3'-O3'	5.70	118.05	109.50
3	C	167	TRP	N-CA-C	5.68	113.44	108.78
21	V	17	THR	N-CA-C	5.63	117.17	109.18
9	I	22	GLY	N-CA-C	5.60	122.36	111.66
1	A	1346	A	P-O3'-C3'	5.59	128.58	120.20
1	A	115	G	C2'-C3'-O3'	5.58	117.87	109.50
5	E	105	VAL	CA-C-N	-5.54	113.31	119.19
5	E	105	VAL	C-N-CA	-5.54	113.31	119.19
19	S	41	VAL	CA-C-N	5.52	126.74	119.84
19	S	41	VAL	C-N-CA	5.52	126.74	119.84
3	C	167	TRP	CB-CA-C	-5.51	109.13	117.07
1	A	129(A)	G	C4'-C3'-O3'	5.50	117.64	109.40
16	P	11	SER	N-CA-C	-5.47	100.80	108.96
1	A	1299	A	N9-C1'-C2'	5.44	120.16	112.00
1	A	812	C	P-O3'-C3'	5.39	128.28	120.20
1	A	328	C	C2'-C3'-O3'	5.38	117.57	109.50
1	A	1101	A	P-O3'-C3'	5.38	128.27	120.20
17	Q	29	HIS	CA-C-N	5.33	124.78	119.24
17	Q	29	HIS	C-N-CA	5.33	124.78	119.24
17	Q	39	SER	N-CA-C	5.31	116.87	108.96
1	A	1300	G	C2'-C3'-O3'	5.30	117.46	109.50
1	A	965	A	P-O3'-C3'	5.30	128.15	120.20
1	A	1190	G	P-O3'-C3'	5.30	128.15	120.20
1	A	701	C	C4'-C3'-O3'	5.29	117.34	109.40
4	D	188	LEU	CA-C-N	5.29	125.21	119.76
4	D	188	LEU	C-N-CA	5.29	125.21	119.76
5	E	41	VAL	CB-CA-C	-5.25	103.33	110.90
11	K	75	TYR	N-CA-C	-5.23	106.58	113.17
6	F	37	VAL	CB-CA-C	-5.20	107.39	113.22
1	A	1183	A	P-O3'-C3'	5.16	127.94	120.20
1	A	1190	G	C2'-C3'-O3'	5.13	117.20	109.50
8	H	85	ARG	CA-C-N	-5.11	117.96	122.97
8	H	85	ARG	C-N-CA	-5.11	117.96	122.97
1	A	243	A	C4'-C3'-O3'	5.11	117.06	109.40
7	G	147	ALA	N-CA-C	-5.08	99.99	110.80
1	A	266	G	C2'-C3'-O3'	5.07	117.11	109.50
2	B	13	ALA	N-CA-C	-5.07	107.06	113.20
1	A	1503	A	C2'-C3'-O3'	5.06	117.09	109.50
1	A	1067	A	P-O3'-C3'	5.03	127.75	120.20
13	M	25	ILE	N-CA-C	5.03	115.14	108.11

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
8	H	90	GLY	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	439	0
2	B	1900	0	1951	65	0
3	C	1612	0	1677	58	0
4	D	1703	0	1763	44	0
5	E	1146	0	1207	39	0
6	F	843	0	857	29	0
7	G	1257	0	1296	31	0
8	H	1116	0	1177	26	0
9	I	1011	0	1043	46	0
10	J	792	0	835	43	0
11	K	885	0	904	22	0
12	L	970	0	1057	42	0
13	M	997	0	1072	37	0
14	N	492	0	529	30	0
15	O	734	0	771	14	0
16	P	700	0	720	29	0
17	Q	857	0	930	23	0
18	R	597	0	668	17	0
19	S	647	0	673	12	0
20	T	762	0	856	20	0
21	V	208	0	221	6	0
22	X	232	0	121	0	0
23	W	82	0	45	1	0
24	A	93	0	0	0	0
24	J	1	0	0	0	0
24	N	1	0	0	0	0
24	W	1	0	0	0	0
25	A	42	0	45	1	0
26	D	1	0	0	0	0
All	All	52062	0	36764	958	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 (958) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:77:ALA:HB2	2:B:211:ILE:HD13	1.52	0.91
1:A:410:G:OP1	1:A:431:A:N6	2.07	0.87
8:H:86:ILE:HD11	8:H:136:GLU:HG3	1.57	0.86
4:D:125:HIS:HD1	4:D:152:SER:HG	1.20	0.85
1:A:542:G:OP1	4:D:10:ARG:NH2	2.10	0.84
7:G:9:VAL:HG13	7:G:94:ARG:HH21	1.42	0.83
1:A:1003:G:N2	1:A:1038:C:O2	2.12	0.83
1:A:448:A:OP2	1:A:485:G:N2	2.11	0.82
1:A:1029:C:N4	1:A:1032:G:O6	2.12	0.82
1:A:581:G:N7	1:A:758:G:C6	2.49	0.81
7:G:146:GLU:HA	7:G:149:ARG:HB2	1.63	0.81
10:J:90:LEU:H	10:J:91:PRO:HD2	1.46	0.80
1:A:664:G:H22	1:A:741:G:H1	1.30	0.78
1:A:836:G:OP1	18:R:61:LYS:NZ	2.16	0.78
13:M:10:PRO:HB2	13:M:18:ALA:HB1	1.66	0.78
10:J:57:LYS:O	10:J:60:ARG:NH1	2.16	0.77
13:M:11:ARG:O	13:M:13:LYS:N	2.17	0.77
12:L:47:LYS:HB3	12:L:48:PRO:HD3	1.65	0.76
1:A:1502:A:H2	1:A:1505:G:H1	1.33	0.76
12:L:56:ALA:HB2	12:L:70:ILE:HD11	1.68	0.76
1:A:1191:A:H5''	3:C:4:LYS:HE2	1.68	0.75
3:C:8:ILE:HG23	3:C:16:ARG:HG2	1.68	0.75
1:A:1305:G:HO2'	1:A:1306:A:H8	1.34	0.75
1:A:758:G:O5'	1:A:758:G:H8	1.68	0.75
1:A:1285:A:H4'	1:A:1286:A:O5'	1.86	0.75
3:C:154:SER:OG	3:C:155:GLY:N	2.17	0.75
1:A:1128:C:H4'	9:I:16:ARG:HH22	1.50	0.75
2:B:93:VAL:HG21	2:B:97:TRP:HD1	1.51	0.75
2:B:93:VAL:HG21	2:B:97:TRP:CD1	2.21	0.75
1:A:1118:C:OP1	9:I:104:ARG:NH1	2.20	0.74
9:I:10:ARG:HD3	9:I:75:ASP:HB3	1.69	0.74
1:A:1004:A:O2'	1:A:1038:C:O2	2.05	0.74
1:A:476:G:C2	1:A:478:A:H5'	2.23	0.73
1:A:1532:U:O2'	1:A:1533:C:O5'	2.07	0.73
1:A:1315:U:O2'	1:A:1360:A:N3	2.20	0.73
18:R:17:SER:OG	18:R:54:ARG:O	2.05	0.73
20:T:75:ASN:OD1	20:T:75:ASN:N	2.21	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1230:C:H1'	13:M:126:LYS:HA	1.69	0.72
9:I:121:ARG:NH1	9:I:122:ALA:O	2.23	0.72
12:L:60:LEU:HD11	12:L:85:ILE:HD12	1.72	0.72
8:H:69:ARG:NH1	8:H:75:ARG:O	2.23	0.72
9:I:100:GLY:O	9:I:102:LEU:N	2.23	0.71
1:A:192:U:H4'	20:T:57:ARG:HD2	1.72	0.71
1:A:477:G:H8	1:A:478:A:H4'	1.55	0.71
1:A:975:A:H4'	1:A:976:G:O5'	1.89	0.71
1:A:1442:G:N2	1:A:1447:G:N7	2.38	0.70
1:A:1129:C:OP1	9:I:62:TYR:OH	2.08	0.70
4:D:22:LYS:O	4:D:24:GLU:N	2.24	0.69
2:B:197:VAL:HB	2:B:200:ILE:HG23	1.72	0.69
1:A:413:G:OP2	1:A:417:C:N4	2.25	0.69
1:A:1310:G:OP2	13:M:88:ARG:NH2	2.26	0.69
2:B:74:LYS:NZ	2:B:205:ASP:OD2	2.25	0.69
9:I:9:ARG:HG2	9:I:14:VAL:HG22	1.75	0.69
1:A:503:C:OP2	12:L:116:SER:OG	2.10	0.69
1:A:1423:G:N2	1:A:1477:C:O2	2.25	0.69
8:H:41:ARG:NH1	8:H:123:GLU:OE2	2.26	0.69
1:A:1392:G:H21	1:A:1502:A:H8	1.38	0.69
1:A:839:U:OP2	1:A:840:C:N4	2.22	0.68
3:C:102:ASN:OD1	3:C:102:ASN:N	2.27	0.68
1:A:474:G:O6	16:P:81:ARG:NH2	2.26	0.68
1:A:1086:U:H3	1:A:1099:G:H22	1.41	0.68
19:S:13:ASP:HA	19:S:16:LEU:HB3	1.76	0.68
13:M:49:THR:HG22	13:M:51:ALA:H	1.58	0.68
15:O:39:LEU:HD13	15:O:56:LEU:HB2	1.76	0.68
1:A:582:U:OP1	15:O:64:ARG:NH2	2.27	0.68
1:A:1305:G:O2'	1:A:1306:A:H8	1.76	0.68
9:I:46:ALA:HB2	9:I:74:ILE:HG23	1.75	0.67
10:J:27:ALA:HB2	10:J:85:LEU:HD21	1.75	0.67
2:B:229:VAL:HG12	2:B:231:GLU:HG3	1.75	0.67
7:G:50:ILE:HG21	7:G:58:PRO:HA	1.77	0.67
2:B:75:LYS:HE2	2:B:96:ARG:HH12	1.58	0.67
17:Q:55:ASP:N	17:Q:55:ASP:OD1	2.26	0.67
8:H:24:THR:HG22	8:H:63:LEU:HD21	1.77	0.66
8:H:91:ARG:HG3	12:L:7:ILE:HG13	1.77	0.66
2:B:16:HIS:CE1	2:B:210:SER:HB3	2.31	0.66
14:N:50:LYS:HB3	14:N:52:GLN:HG3	1.78	0.66
1:A:26:A:O2'	4:D:209:ARG:NH1	2.29	0.66
1:A:462:G:C2	16:P:82:GLN:HB3	2.31	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:58:GLU:HB3	10:J:92:THR:HG21	1.78	0.66
1:A:376:G:H5''	16:P:5:ARG:HD2	1.78	0.65
16:P:81:ARG:C	16:P:83:GLU:H	2.02	0.65
1:A:141:A:H1'	1:A:182:U:O2	1.95	0.65
2:B:60:ASP:HB3	2:B:64:ARG:HH21	1.61	0.65
4:D:149:ALA:HB3	4:D:152:SER:HB2	1.77	0.65
1:A:456:C:H42	1:A:476:G:H1'	1.61	0.65
1:A:1162:C:O2	1:A:1174:G:N2	2.20	0.65
4:D:7:PRO:HG2	4:D:10:ARG:HD2	1.78	0.65
12:L:27:LEU:O	12:L:29:GLY:N	2.29	0.65
1:A:477:G:C8	1:A:478:A:H4'	2.32	0.65
16:P:22:THR:HA	16:P:33:ILE:HG13	1.78	0.65
5:E:79:GLU:HG3	5:E:93:PRO:HD3	1.77	0.65
16:P:81:ARG:HA	16:P:81:ARG:HH11	1.62	0.64
1:A:353:A:H5'	1:A:353:A:H8	1.61	0.64
2:B:102:LEU:HD21	2:B:162:ILE:HD11	1.78	0.64
3:C:32:LEU:HD21	3:C:59:ARG:HD2	1.79	0.64
1:A:406:G:C8	1:A:496:A:C5	2.86	0.64
1:A:462:G:N2	16:P:82:GLN:HB3	2.13	0.64
1:A:1144:G:H21	1:A:1146:A:H62	1.46	0.64
3:C:11:ARG:O	3:C:13:GLY:N	2.30	0.64
7:G:18:TYR:CD2	7:G:59:LEU:HB2	2.33	0.64
1:A:275:G:H5'	17:Q:14:LYS:HE2	1.78	0.64
3:C:130:VAL:HG21	3:C:157:ILE:HG23	1.80	0.64
10:J:89:ASP:HB2	10:J:91:PRO:HD2	1.80	0.63
1:A:113:G:H1'	1:A:354:G:H5'	1.79	0.63
1:A:1435:G:H2'	1:A:1436:U:C6	2.33	0.63
1:A:35:G:H2'	1:A:36:C:C6	2.34	0.63
3:C:6:HIS:HD2	3:C:9:GLY:H	1.44	0.63
13:M:16:ASP:OD1	13:M:16:ASP:N	2.23	0.63
1:A:1073:U:OP2	5:E:57:LYS:NZ	2.31	0.63
5:E:43:LEU:HB2	5:E:136:MET:HE2	1.80	0.63
13:M:12:ASN:HB2	13:M:46:LYS:HD2	1.79	0.63
1:A:255:G:O3'	17:Q:17:LYS:NZ	2.31	0.63
1:A:1356:G:H2'	1:A:1357:A:C8	2.33	0.63
4:D:176:LEU:HA	4:D:183:GLY:HA2	1.81	0.62
2:B:15:VAL:HG11	2:B:209:ARG:HB3	1.80	0.62
16:P:71:ARG:HD3	16:P:75:ARG:NH2	2.15	0.62
8:H:46:LYS:HG3	8:H:64:LYS:HG3	1.79	0.62
1:A:713:G:H2'	1:A:714:G:C8	2.35	0.62
11:K:40:ILE:HG22	11:K:41:THR:HG23	1.80	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:A:H2'	1:A:326:G:N2	2.15	0.61
2:B:87:ARG:NH2	2:B:220:ASP:OD1	2.34	0.61
1:A:181:G:H4'	1:A:182:U:H5'	1.83	0.61
1:A:537:G:OP1	12:L:113:ARG:NH2	2.34	0.61
16:P:81:ARG:HA	16:P:81:ARG:NH1	2.16	0.61
1:A:1189:C:OP1	10:J:51:ARG:NH2	2.33	0.61
5:E:81:GLU:HG2	5:E:90:VAL:HG22	1.81	0.61
7:G:18:TYR:HD2	7:G:59:LEU:HB2	1.65	0.61
5:E:50:GLU:HG3	5:E:52:PRO:HD2	1.83	0.60
10:J:60:ARG:O	10:J:61:GLU:HB3	2.01	0.60
5:E:43:LEU:HD11	5:E:132:ALA:HB1	1.84	0.60
1:A:581:G:C8	1:A:758:G:C6	2.90	0.60
1:A:1302:U:OP1	13:M:13:LYS:NZ	2.34	0.60
1:A:1327:C:OP2	21:V:12:LYS:NZ	2.33	0.60
1:A:190(D):U:O2'	1:A:190(E):U:H5'	2.02	0.59
1:A:582:U:P	15:O:64:ARG:HH22	2.25	0.59
3:C:36:ASP:OD2	3:C:59:ARG:NH2	2.36	0.59
1:A:975:A:N6	1:A:1366:C:O2'	2.35	0.59
1:A:1298:C:C4	7:G:114:ARG:HD3	2.37	0.59
18:R:53:ARG:HD2	18:R:58:LEU:O	2.03	0.59
2:B:235:SER:O	2:B:237:ALA:N	2.35	0.59
1:A:760:G:H22	17:Q:105:ALA:HA	1.68	0.59
1:A:1281:U:H5'	1:A:1282:C:H5	1.68	0.59
1:A:477:G:H2'	1:A:479:C:OP2	2.03	0.59
1:A:958:A:N3	1:A:985:C:O2'	2.35	0.59
1:A:1029:C:N3	1:A:1033:G:N2	2.51	0.59
1:A:109:A:H2'	1:A:326:G:H21	1.67	0.58
5:E:33:VAL:HG11	5:E:109:ILE:HA	1.84	0.58
12:L:27:LEU:C	12:L:29:GLY:H	2.12	0.58
13:M:34:LEU:HD12	13:M:39:ILE:HB	1.85	0.58
1:A:795:C:H5''	1:A:796:C:OP2	2.02	0.58
1:A:975:A:H5'	1:A:975:A:C8	2.38	0.58
4:D:151:LYS:H	4:D:151:LYS:HD2	1.67	0.58
19:S:42:PRO:O	19:S:44:MET:N	2.32	0.58
1:A:1392:G:N2	1:A:1502:A:H8	2.00	0.58
10:J:65:LEU:HB2	14:N:56:VAL:HG22	1.86	0.58
1:A:460:A:N6	1:A:463:A:N1	2.52	0.58
3:C:157:ILE:HD12	3:C:164:ARG:HB3	1.85	0.58
1:A:235:C:H5'	17:Q:70:ARG:HG2	1.84	0.58
10:J:24:VAL:HG12	10:J:28:ARG:HD2	1.85	0.58
13:M:37:THR:HG22	13:M:55:ARG:HD2	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:243:A:H4'	1:A:244:U:O5'	2.04	0.58
1:A:530:G:O6	23:W:3:U:O2'	2.15	0.58
1:A:1227:A:C2	13:M:117:VAL:HG21	2.38	0.58
1:A:669:U:H2'	1:A:670:G:C8	2.39	0.58
1:A:1129:C:H1'	1:A:1132:C:H5	1.68	0.58
1:A:1190:G:H5'	3:C:176:HIS:CE1	2.39	0.58
1:A:409:G:N1	1:A:433:C:H5'	2.19	0.57
1:A:1057:G:H2'	1:A:1058:G:O4'	2.04	0.57
9:I:116:LYS:HD2	9:I:122:ALA:HA	1.86	0.57
12:L:32:PHE:HB3	12:L:84:LEU:HD11	1.87	0.57
1:A:757:U:H2'	1:A:758:G:O4'	2.04	0.57
3:C:107:GLN:CD	3:C:107:GLN:H	2.11	0.57
6:F:76:ALA:HA	6:F:79:LEU:HB2	1.85	0.57
1:A:1366:C:O2'	10:J:60:ARG:NH2	2.38	0.57
2:B:88:ALA:HB2	2:B:219:VAL:HG13	1.87	0.57
12:L:70:ILE:HD13	12:L:77:LEU:HD12	1.85	0.57
9:I:93:ARG:HB3	9:I:93:ARG:NH1	2.20	0.57
1:A:299:G:H2'	1:A:300:A:C8	2.40	0.57
1:A:1124:G:O2'	1:A:1126:U:O4	2.22	0.57
3:C:164:ARG:HG2	3:C:165:THR:H	1.69	0.57
6:F:71:ARG:O	6:F:73:ASN:N	2.38	0.57
16:P:21:VAL:HG21	16:P:59:TRP:CG	2.39	0.57
1:A:1001:A:H2'	1:A:1002:G:C8	2.40	0.57
4:D:162:LEU:HD22	4:D:181:MET:HG2	1.87	0.57
1:A:328:C:H2'	1:A:328:C:O2	2.04	0.56
20:T:33:ILE:HD13	20:T:63:ILE:HG12	1.87	0.56
1:A:1141:C:H2'	1:A:1142:G:H8	1.70	0.56
1:A:1316:G:H4'	14:N:18:VAL:HG11	1.87	0.56
1:A:1510:U:H2'	1:A:1511:G:C8	2.40	0.56
4:D:108:LEU:HD22	4:D:176:LEU:HB2	1.87	0.56
1:A:409:G:H5'	1:A:430:A:N7	2.21	0.56
1:A:1101:A:H4'	1:A:1102:A:O5'	2.05	0.56
13:M:84:ILE:HG21	19:S:65:ASN:HB2	1.87	0.56
1:A:731:G:OP1	1:A:766:A:H1'	2.05	0.56
4:D:162:LEU:HD23	4:D:178:VAL:HG13	1.88	0.56
1:A:463:A:H1'	1:A:474:G:N2	2.20	0.56
20:T:44:ALA:C	20:T:46:GLU:H	2.14	0.56
1:A:792:A:H4'	1:A:793:U:H5''	1.88	0.56
16:P:81:ARG:O	16:P:83:GLU:N	2.38	0.56
1:A:409:G:C6	1:A:433:C:H5'	2.41	0.56
2:B:80:ILE:HD12	2:B:80:ILE:H	1.71	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:189:ASP:OD1	2:B:189:ASP:N	2.33	0.56
3:C:88:ARG:HG3	3:C:101:LEU:HD12	1.86	0.56
1:A:669:U:H2'	1:A:670:G:H8	1.70	0.56
6:F:2:ARG:NE	6:F:69:GLU:HG2	2.20	0.56
6:F:100:ASN:ND2	18:R:26:LEU:O	2.38	0.56
20:T:67:ALA:HA	20:T:73:HIS:H	1.71	0.56
1:A:1288:A:N3	1:A:1352:C:O2'	2.35	0.55
3:C:3:ASN:OD1	3:C:3:ASN:N	2.39	0.55
8:H:10:LEU:HD22	8:H:83:ILE:HD11	1.87	0.55
2:B:97:TRP:HZ2	2:B:102:LEU:HD13	1.71	0.55
10:J:79:ARG:HG3	10:J:82:ILE:HD12	1.87	0.55
20:T:63:ILE:HD13	20:T:80:ARG:HB3	1.88	0.55
3:C:63:ASN:HA	3:C:99:VAL:HG12	1.88	0.55
4:D:12:CYS:HB3	4:D:33:MET:HG2	1.88	0.55
11:K:21:ILE:HD11	11:K:98:LEU:HD12	1.87	0.55
15:O:26:GLU:OE1	15:O:77:ARG:HD2	2.06	0.55
21:V:15:ARG:O	21:V:17:THR:HG23	2.06	0.55
1:A:453:A:H62	1:A:479:C:H5	1.53	0.55
1:A:1502:A:H2	1:A:1505:G:N1	2.03	0.55
2:B:121:LEU:HD22	2:B:126:GLU:HB2	1.89	0.55
4:D:140:VAL:HG11	4:D:146:ILE:HD11	1.89	0.55
6:F:46:ARG:HB2	6:F:60:PHE:CE1	2.41	0.55
19:S:55:LYS:HG2	19:S:56:GLN:HG2	1.88	0.55
1:A:346:G:H2'	1:A:347:G:O4'	2.06	0.55
1:A:320:C:H2'	1:A:321:A:C8	2.42	0.55
1:A:928:G:H5'	1:A:1533:C:H5'	1.88	0.55
1:A:1126:U:H3	1:A:1127:G:N2	2.05	0.55
20:T:39:LYS:HD3	20:T:55:ILE:HD13	1.89	0.55
2:B:75:LYS:HA	2:B:78:GLN:HB2	1.89	0.54
8:H:89:PRO:HA	8:H:92:ARG:NH1	2.22	0.54
1:A:662:G:H2'	1:A:663:A:C8	2.41	0.54
1:A:1305:G:C8	1:A:1305:G:OP2	2.60	0.54
1:A:1360:A:H3'	1:A:1361:G:N7	2.23	0.54
1:A:878:G:H5'	8:H:89:PRO:HG2	1.89	0.54
6:F:98:LEU:HD12	18:R:28:GLU:HB3	1.89	0.54
16:P:19:ILE:HG22	16:P:36:ILE:HG13	1.90	0.54
1:A:405:U:OP1	1:A:406:G:O2'	2.24	0.54
8:H:20:TYR:CE1	8:H:76:PRO:HD2	2.43	0.54
14:N:12:ARG:C	14:N:14:PRO:HD3	2.33	0.54
5:E:151:LEU:HD23	8:H:79:VAL:HA	1.90	0.54
17:Q:28:PRO:HA	17:Q:35:VAL:HA	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:V:5:ASP:O	21:V:8:THR:OG1	2.25	0.54
6:F:61:LEU:HB3	6:F:63:TYR:HE1	1.72	0.54
1:A:267:C:H2'	1:A:268:C:C6	2.43	0.54
1:A:1268:A:N3	1:A:1326:C:O2'	2.38	0.54
12:L:89:ARG:HH22	12:L:97:ARG:HH11	1.55	0.54
1:A:676:A:H1'	11:K:115:PRO:HB3	1.90	0.54
1:A:1396:A:H2	5:E:19:MET:HG3	1.73	0.54
2:B:51:LEU:HD21	2:B:201:ILE:HG23	1.89	0.54
1:A:1116:C:O2'	9:I:108:VAL:HG21	2.08	0.54
1:A:448:A:H62	1:A:486:U:H3	1.56	0.54
1:A:1397:C:H4'	1:A:1398:A:OP2	2.08	0.54
1:A:1420:C:H2'	1:A:1421:G:C8	2.43	0.54
12:L:126:LYS:H	12:L:126:LYS:HD2	1.73	0.54
5:E:84:PHE:HB3	5:E:134:ALA:HB2	1.89	0.53
1:A:310:G:H5'	16:P:31:LYS:HB2	1.90	0.53
1:A:547:A:OP2	4:D:2:GLY:N	2.41	0.53
4:D:173:TRP:CD1	4:D:189:PRO:HD3	2.44	0.53
5:E:145:LYS:O	5:E:149:GLU:HG2	2.08	0.53
9:I:10:ARG:NH1	9:I:75:ASP:OD2	2.41	0.53
12:L:27:LEU:C	12:L:29:GLY:N	2.66	0.53
1:A:490:G:H2'	1:A:491:G:H8	1.72	0.53
1:A:677:U:H3	1:A:713:G:H22	1.56	0.53
1:A:409:G:N2	1:A:431:A:O2'	2.41	0.53
7:G:78:ARG:HB2	7:G:156:TRP:HZ3	1.74	0.53
8:H:7:ALA:HA	8:H:85:ARG:HG3	1.90	0.53
1:A:1136:U:H5''	1:A:1137:C:H5	1.74	0.53
1:A:1112:C:O2	3:C:179:ARG:HD2	2.08	0.53
8:H:4:ASP:OD2	8:H:85:ARG:NH1	2.35	0.53
14:N:24:CYS:O	14:N:26:ARG:N	2.41	0.53
1:A:1427:U:H2'	1:A:1428:A:C8	2.44	0.53
1:A:1427:U:H2'	1:A:1428:A:H8	1.74	0.53
1:A:462:G:H1	16:P:82:GLN:HA	1.74	0.53
1:A:701:C:H5''	1:A:703:G:O4'	2.08	0.53
3:C:141:VAL:HG11	3:C:202:ILE:HG12	1.91	0.53
1:A:1048:G:O3'	1:A:1049:U:H3'	2.09	0.53
3:C:68:VAL:HG12	3:C:70:VAL:HG23	1.91	0.53
5:E:80:ILE:HD13	5:E:138:ALA:HB1	1.89	0.53
6:F:23:LYS:HE3	6:F:61:LEU:HD21	1.91	0.53
12:L:25:PRO:C	12:L:27:LEU:H	2.16	0.53
1:A:115:G:H1'	1:A:116:A:N7	2.24	0.53
1:A:1052:U:O2'	1:A:1055:A:OP2	2.22	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:46:ARG:HB2	6:F:60:PHE:HE1	1.74	0.53
10:J:24:VAL:O	10:J:28:ARG:HB2	2.09	0.53
1:A:103:C:OP1	20:T:17:ARG:NH1	2.43	0.52
1:A:517:G:N1	1:A:533:A:OP2	2.30	0.52
2:B:19:HIS:NE2	2:B:205:ASP:OD1	2.42	0.52
4:D:175:SER:O	4:D:175:SER:OG	2.25	0.52
10:J:29:ARG:HD2	10:J:84:GLN:HE22	1.75	0.52
1:A:279:A:OP2	17:Q:95:TYR:OH	2.22	0.52
1:A:409:G:N2	1:A:433:C:OP1	2.43	0.52
1:A:421:U:O2'	1:A:422:C:H5'	2.09	0.52
12:L:89:ARG:NH2	12:L:97:ARG:HH11	2.08	0.52
13:M:11:ARG:C	13:M:13:LYS:H	2.18	0.52
1:A:527:G:O2'	1:A:535:A:N1	2.41	0.52
1:A:1443:G:H5''	1:A:1446:A:H5''	1.91	0.52
8:H:92:ARG:HG2	8:H:92:ARG:HH11	1.74	0.52
16:P:21:VAL:HG12	16:P:33:ILE:HB	1.92	0.52
1:A:409:G:N1	1:A:433:C:OP1	2.42	0.52
6:F:44:GLY:HA2	6:F:59:TYR:CZ	2.44	0.52
13:M:8:GLU:OE1	13:M:22:ILE:HA	2.09	0.52
1:A:316:G:OP2	1:A:351:G:O2'	2.28	0.52
1:A:975:A:H5'	1:A:975:A:H8	1.74	0.52
1:A:976:G:N7	1:A:1359:C:H1'	2.25	0.52
5:E:12:LEU:HD13	5:E:31:LEU:HB2	1.91	0.52
9:I:43:ALA:HA	9:I:74:ILE:HD13	1.92	0.52
13:M:90:LEU:HD21	13:M:94:ARG:NH2	2.24	0.52
17:Q:79:SER:OG	17:Q:80:GLY:N	2.42	0.52
13:M:47:ASP:OD1	13:M:47:ASP:N	2.42	0.52
1:A:417:C:H6	1:A:417:C:H3'	1.75	0.52
7:G:89:MET:HE1	7:G:156:TRP:HB3	1.92	0.52
10:J:39:PRO:HA	10:J:70:ARG:HH11	1.73	0.52
1:A:974:A:OP2	14:N:29:ARG:NH2	2.43	0.52
1:A:1366:C:H2'	1:A:1367:C:H6	1.75	0.52
10:J:90:LEU:H	10:J:91:PRO:CD	2.19	0.52
12:L:82:VAL:N	12:L:106:ASP:OD2	2.32	0.51
13:M:77:ASN:O	13:M:81:LEU:HG	2.10	0.51
16:P:11:SER:HB3	16:P:14:ASN:HB3	1.93	0.51
1:A:1187:G:O5'	9:I:113:LYS:NZ	2.44	0.51
2:B:51:LEU:HD22	2:B:55:PHE:HE2	1.76	0.51
10:J:79:ARG:HA	10:J:82:ILE:HG13	1.92	0.51
1:A:1120:G:H2'	1:A:1121:U:C6	2.46	0.51
3:C:116:VAL:HG21	3:C:202:ILE:HD11	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:P:69:THR:O	16:P:72:ARG:HB2	2.10	0.51
1:A:501:C:H2'	1:A:502:G:C8	2.46	0.51
15:O:36:ILE:HG12	15:O:59:MET:HE2	1.93	0.51
1:A:1366:C:H2'	1:A:1367:C:C6	2.46	0.51
3:C:6:HIS:CD2	3:C:9:GLY:H	2.27	0.51
9:I:49:PRO:O	9:I:53:VAL:HG23	2.11	0.51
1:A:581:G:N7	1:A:758:G:O6	2.43	0.51
2:B:132:LYS:HD3	2:B:135:GLN:HG3	1.93	0.51
12:L:46:LYS:HE2	12:L:47:LYS:HB2	1.92	0.51
13:M:49:THR:HG22	13:M:51:ALA:N	2.26	0.51
19:S:28:LYS:HG2	19:S:31:ILE:HD11	1.93	0.51
1:A:298:A:H2'	1:A:299:G:O4'	2.11	0.51
1:A:523:A:C2	12:L:91:LYS:HB3	2.46	0.51
1:A:673:G:H2'	1:A:674:G:C8	2.46	0.51
1:A:519:C:H2'	1:A:520:A:C8	2.46	0.50
5:E:51:VAL:HB	5:E:52:PRO:HD3	1.91	0.50
13:M:3:ARG:O	13:M:57:ARG:NH2	2.44	0.50
1:A:706:A:H1'	11:K:29:ILE:HD11	1.92	0.50
2:B:80:ILE:HG21	2:B:212:GLN:HA	1.93	0.50
9:I:53:VAL:C	9:I:55:ALA:H	2.19	0.50
9:I:118:LYS:O	9:I:119:ALA:HB3	2.12	0.50
15:O:78:TYR:CZ	15:O:82:ILE:HD11	2.46	0.50
1:A:458:C:H42	1:A:463:A:H3'	1.75	0.50
1:A:485:G:O2'	1:A:486:U:P	2.70	0.50
1:A:939:G:H5''	7:G:102:ARG:HH22	1.75	0.50
11:K:48:ILE:HG13	11:K:63:LEU:HB3	1.91	0.50
1:A:858:G:O6	1:A:869:G:C8	2.64	0.50
1:A:1133:G:H2'	1:A:1134:G:H8	1.76	0.50
1:A:1196:U:OP1	1:A:1197:G:H5'	2.12	0.50
4:D:24:GLU:OE1	4:D:25:ARG:N	2.43	0.50
19:S:22:LEU:HD22	19:S:28:LYS:HD3	1.94	0.50
1:A:1068:G:H8	1:A:1068:G:OP2	1.94	0.50
10:J:82:ILE:HA	10:J:85:LEU:HD12	1.93	0.50
16:P:4:ILE:HG13	16:P:64:ALA:HB1	1.94	0.50
1:A:581:G:N7	1:A:758:G:C5	2.80	0.50
1:A:707:C:OP1	11:K:85:ARG:NH1	2.43	0.50
1:A:1486:G:H2'	1:A:1487:G:O4'	2.12	0.50
3:C:108:ASN:C	3:C:110:ASN:H	2.20	0.50
12:L:47:LYS:HB3	12:L:48:PRO:CD	2.38	0.50
1:A:267:C:H2'	1:A:268:C:H6	1.76	0.50
1:A:581:G:O6	1:A:758:G:C8	2.65	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:R:58:LEU:HD13	18:R:62:GLU:HB3	1.94	0.50
1:A:255:G:H2'	1:A:256:U:C6	2.47	0.49
1:A:406:G:N3	1:A:406:G:H2'	2.27	0.49
1:A:1329:A:P	13:M:28:ALA:HB3	2.52	0.49
6:F:92:LYS:HB2	6:F:92:LYS:NZ	2.27	0.49
7:G:114:ARG:O	7:G:119:ARG:NH1	2.45	0.49
2:B:236:TYR:C	2:B:238:LEU:H	2.20	0.49
1:A:294:U:OP1	1:A:610:G:O2'	2.26	0.49
1:A:1141:C:H2'	1:A:1142:G:C8	2.46	0.49
6:F:33:TYR:HA	6:F:71:ARG:NH2	2.27	0.49
8:H:86:ILE:HG21	8:H:133:LEU:HD22	1.94	0.49
1:A:1062:U:H2'	1:A:1063:C:C6	2.47	0.49
1:A:1128:C:H5'	9:I:16:ARG:HH12	1.76	0.49
10:J:16:LEU:O	10:J:20:ALA:N	2.29	0.49
12:L:65:GLU:OE1	12:L:65:GLU:N	2.45	0.49
1:A:1443:G:H4'	1:A:1446:A:H5''	1.95	0.49
3:C:150:LYS:HG3	3:C:169:ALA:HB2	1.94	0.49
4:D:90:GLY:CA	4:D:204:ILE:HD11	2.43	0.49
5:E:5:ASP:OD1	5:E:5:ASP:N	2.45	0.49
7:G:108:ALA:O	7:G:111:ARG:HB2	2.12	0.49
9:I:36:TYR:OH	9:I:73:GLN:NE2	2.31	0.49
1:A:232:G:H1'	1:A:262:A:N1	2.28	0.49
1:A:404:U:H2'	1:A:405:U:H6	1.77	0.49
4:D:102:ASP:OD1	4:D:103:ASN:N	2.45	0.49
1:A:165:C:H2'	1:A:166:G:C8	2.48	0.49
1:A:362:G:N2	1:A:365:U:OP2	2.45	0.49
1:A:401:C:O2'	1:A:621:A:N3	2.34	0.49
1:A:447:G:H1	1:A:485:G:HO2'	1.61	0.49
1:A:560:U:H4'	1:A:561:U:H5''	1.93	0.49
1:A:1000:U:H2'	1:A:1001:A:O4'	2.13	0.49
1:A:1329:A:H5'	13:M:29:ARG:HD2	1.94	0.49
5:E:76:ILE:HD13	5:E:118:ILE:HD12	1.95	0.49
1:A:21:G:H2'	1:A:22:G:C8	2.48	0.49
1:A:353:A:H5'	1:A:353:A:C8	2.45	0.49
1:A:718:G:C8	11:K:116:HIS:HB3	2.48	0.49
2:B:60:ASP:HB3	2:B:64:ARG:NH2	2.26	0.49
3:C:6:HIS:CD2	3:C:8:ILE:HB	2.47	0.49
14:N:26:ARG:HE	14:N:43:CYS:HB3	1.77	0.49
20:T:50:GLU:HA	20:T:100:ILE:HG22	1.95	0.49
1:A:1182:G:H4'	1:A:1183:A:H3'	1.94	0.49
2:B:186:ALA:HB3	2:B:197:VAL:HG11	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:167:TRP:HE3	3:C:168:ALA:H	1.61	0.49
4:D:62:GLN:OE1	4:D:65:ARG:NH1	2.46	0.49
9:I:15:ALA:HA	9:I:65:VAL:HG12	1.95	0.49
10:J:34:VAL:HG12	10:J:35:SER:H	1.76	0.49
12:L:24:VAL:HG13	12:L:98:TYR:HE2	1.78	0.49
1:A:262:A:C6	1:A:263:A:C6	3.00	0.49
1:A:269:C:H2'	1:A:270:A:C8	2.48	0.49
1:A:939:G:H5''	7:G:102:ARG:NH2	2.28	0.49
1:A:1346:A:H5'	1:A:1348:U:H1'	1.94	0.49
1:A:1137:C:H5'	1:A:1138:G:C4	2.48	0.48
1:A:1261:A:H3'	1:A:1262:C:H6	1.78	0.48
5:E:6:PHE:CE1	5:E:36:ASP:HB3	2.48	0.48
1:A:409:G:H22	1:A:433:C:P	2.36	0.48
1:A:1179:A:H5''	9:I:102:LEU:O	2.12	0.48
2:B:77:ALA:O	2:B:81:VAL:HG23	2.13	0.48
1:A:358:U:H2'	1:A:359:U:H6	1.77	0.48
1:A:424:G:N3	1:A:424:G:H3'	2.28	0.48
1:A:951:G:OP2	13:M:102:ARG:NH2	2.46	0.48
1:A:1005:A:H3'	1:A:1006:C:H6	1.78	0.48
1:A:1370:G:C2	1:A:1371:G:N7	2.81	0.48
5:E:11:ILE:HD11	5:E:108:ALA:HB3	1.94	0.48
14:N:9:LYS:C	14:N:11:LYS:N	2.71	0.48
18:R:47:THR:HG22	18:R:48:GLY:H	1.79	0.48
1:A:411:A:C2	1:A:417:C:H1'	2.48	0.48
10:J:51:ARG:HG3	10:J:60:ARG:O	2.14	0.48
15:O:25:THR:HG21	15:O:70:LEU:HD23	1.94	0.48
2:B:17:PHE:HB3	2:B:44:LEU:HD11	1.94	0.48
9:I:70:LYS:O	9:I:74:ILE:HG13	2.14	0.48
17:Q:59:ILE:HG23	17:Q:71:PHE:HB3	1.95	0.48
19:S:21:GLU:C	19:S:23:ASN:H	2.20	0.48
1:A:409:G:H2'	1:A:409:G:N3	2.29	0.48
1:A:484:G:H4'	1:A:485:G:O5'	2.13	0.48
1:A:539:A:H2'	1:A:540:G:C8	2.48	0.48
2:B:223:ILE:O	2:B:226:ARG:N	2.40	0.48
1:A:165:C:H2'	1:A:166:G:H8	1.78	0.48
1:A:1064:G:N2	1:A:1190:G:O2'	2.47	0.48
1:A:1228:C:H4'	13:M:116:THR:HA	1.95	0.48
1:A:1298:C:C5	7:G:114:ARG:HD3	2.47	0.48
6:F:6:VAL:HG22	6:F:90:VAL:HG22	1.96	0.48
14:N:22:THR:O	14:N:23:ARG:HG3	2.13	0.48
1:A:175:C:H2'	1:A:176:C:H6	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:485:G:O2'	1:A:486:U:OP2	2.32	0.48
1:A:803:G:H2'	1:A:804:U:O4'	2.13	0.48
10:J:8:LEU:HD23	10:J:94:VAL:HG11	1.96	0.48
10:J:38:ILE:HD12	10:J:71:LEU:HB3	1.95	0.48
1:A:190:C:H4'	1:A:190(A):C:OP1	2.12	0.48
1:A:266:G:H5''	1:A:268:C:H41	1.79	0.48
5:E:11:ILE:HB	5:E:31:LEU:HB3	1.96	0.48
12:L:27:LEU:HG	12:L:28:LYS:H	1.78	0.48
5:E:144:THR:O	5:E:148:VAL:HG23	2.14	0.48
1:A:7:G:H5'	1:A:298:A:O4'	2.14	0.47
1:A:403:C:O3'	4:D:122:ARG:HD2	2.14	0.47
1:A:812:C:H4'	1:A:813:U:H5'	1.96	0.47
1:A:1369:C:H2'	1:A:1370:G:O4'	2.14	0.47
2:B:162:ILE:HD12	2:B:177:ALA:HB2	1.96	0.47
4:D:173:TRP:HB2	4:D:187:ARG:O	2.14	0.47
10:J:61:GLU:OE1	14:N:45:ARG:NH1	2.43	0.47
1:A:263:A:OP2	20:T:79:ARG:NH1	2.47	0.47
1:A:381:C:H2'	1:A:382:A:O4'	2.14	0.47
1:A:397:A:H3'	1:A:397:A:N3	2.29	0.47
1:A:1116:C:H2'	1:A:1117:G:H5''	1.97	0.47
5:E:39:GLY:O	5:E:68:GLU:HA	2.14	0.47
6:F:10:LEU:HD12	6:F:59:TYR:HB3	1.95	0.47
13:M:67:GLU:HB3	13:M:68:GLY:H	1.45	0.47
1:A:200:G:H2'	1:A:201:C:O4'	2.14	0.47
1:A:409:G:OP1	1:A:429:U:H1'	2.14	0.47
1:A:1223:C:OP1	1:A:1224:G:H3'	2.13	0.47
1:A:6:G:H4'	1:A:298:A:H4'	1.96	0.47
1:A:1401:G:C2	1:A:1402:C:H1'	2.50	0.47
2:B:187:LEU:HD12	2:B:201:ILE:O	2.14	0.47
4:D:187:ARG:HH12	4:D:188:LEU:HD12	1.79	0.47
5:E:79:GLU:HG3	5:E:93:PRO:CD	2.42	0.47
18:R:73:ALA:HB3	18:R:79:LEU:HD12	1.95	0.47
1:A:227:G:O2'	16:P:62:VAL:HG11	2.13	0.47
1:A:403:C:OP1	4:D:136:PRO:HD2	2.14	0.47
1:A:418:C:H2'	1:A:419:C:C6	2.49	0.47
1:A:840:C:H5''	1:A:841:U:OP1	2.15	0.47
3:C:30:ARG:NH2	3:C:31:HIS:HE1	2.13	0.47
3:C:155:GLY:HA2	3:C:164:ARG:H	1.78	0.47
6:F:25:ILE:HG21	6:F:82:ARG:HD2	1.96	0.47
1:A:348:G:H2'	1:A:348:G:N3	2.30	0.47
11:K:59:TYR:CE2	11:K:63:LEU:HD11	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:407:G:O2'	1:A:431:A:OP1	2.27	0.47
1:A:502:G:H2'	1:A:503:C:O4'	2.15	0.47
1:A:965:A:O2'	1:A:966:G:OP2	2.23	0.47
1:A:1133:G:H2'	1:A:1134:G:C8	2.49	0.47
1:A:1191:A:OP2	3:C:3:ASN:ND2	2.47	0.47
1:A:1346:A:H5''	9:I:120:ARG:HH12	1.79	0.47
1:A:1417:G:O2'	1:A:1483:A:N6	2.48	0.47
8:H:20:TYR:CE2	8:H:75:ARG:HD2	2.50	0.47
8:H:25:ASP:N	8:H:25:ASP:OD1	2.48	0.47
11:K:22:HIS:O	11:K:22:HIS:ND1	2.48	0.47
11:K:77:MET:HE3	11:K:77:MET:HB3	1.75	0.47
20:T:68:LYS:HD2	20:T:68:LYS:HA	1.73	0.47
1:A:153:C:H42	1:A:168:G:H1	1.62	0.47
1:A:1221:G:OP2	19:S:37:ARG:NH2	2.47	0.47
3:C:119:ARG:HG2	3:C:140:ARG:HH12	1.79	0.47
10:J:5:ARG:HA	10:J:73:ASP:OD1	2.15	0.47
1:A:235:C:H1'	17:Q:61:GLU:OE1	2.15	0.47
1:A:1122:U:O4	1:A:1123:A:N6	2.48	0.47
3:C:12:LEU:HD11	14:N:51:GLY:HA2	1.96	0.47
5:E:66:MET:HE3	5:E:66:MET:HB3	1.76	0.47
10:J:12:ASP:HB3	10:J:15:THR:HG22	1.97	0.47
1:A:191:G:O2'	20:T:102:GLY:O	2.27	0.46
1:A:409:G:O2'	1:A:410:G:OP2	2.32	0.46
1:A:1010:G:N1	1:A:1019:C:O2	2.46	0.46
2:B:90:MET:HE3	2:B:222:ILE:HG21	1.97	0.46
12:L:37:CYS:HB2	12:L:81:SER:O	2.15	0.46
19:S:11:VAL:HG12	19:S:16:LEU:HB2	1.97	0.46
1:A:110:C:H2'	1:A:111:G:O4'	2.15	0.46
1:A:418:C:H6	1:A:419:C:C5	2.32	0.46
1:A:736:C:H2'	1:A:737:A:C8	2.50	0.46
2:B:95:GLN:O	2:B:96:ARG:HD2	2.16	0.46
9:I:127:LYS:HD3	13:M:126:LYS:NZ	2.30	0.46
10:J:6:ILE:HG23	10:J:98:ILE:HG22	1.96	0.46
1:A:149:A:H2'	1:A:150:C:C6	2.50	0.46
1:A:338:A:H2'	1:A:339:C:C6	2.49	0.46
1:A:560:U:H5'	1:A:566:G:C2	2.50	0.46
1:A:1281:U:H4'	1:A:1282:C:OP2	2.14	0.46
1:A:10:A:OP2	5:E:126:ARG:HD3	2.15	0.46
1:A:189:G:N2	1:A:190(K):G:N3	2.63	0.46
1:A:250:A:O5'	1:A:250:A:H8	1.99	0.46
1:A:273:A:N6	1:A:274:A:C6	2.84	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1347:G:N2	1:A:1373:G:H2'	2.29	0.46
2:B:19:HIS:HD2	2:B:189:ASP:HB2	1.79	0.46
2:B:25:ASN:O	2:B:27:LYS:N	2.48	0.46
6:F:22:GLU:OE1	6:F:82:ARG:HD3	2.15	0.46
10:J:64:GLU:OE2	10:J:66:ARG:NH1	2.48	0.46
13:M:91:ARG:HA	13:M:91:ARG:HD2	1.73	0.46
20:T:20:LEU:HA	20:T:20:LEU:HD23	1.71	0.46
21:V:6:ARG:HB3	21:V:15:ARG:NH1	2.31	0.46
1:A:749:C:H2'	1:A:750:G:H8	1.81	0.46
3:C:112:SER:O	3:C:116:VAL:HG23	2.15	0.46
1:A:1212:U:H1'	1:A:1213:A:OP2	2.15	0.46
2:B:196:LEU:HD13	2:B:196:LEU:HA	1.68	0.46
3:C:22:TRP:CE3	3:C:32:LEU:HD22	2.51	0.46
9:I:110:GLU:HG2	9:I:119:ALA:HB1	1.97	0.46
13:M:23:TYR:HB3	13:M:67:GLU:HA	1.98	0.46
16:P:81:ARG:C	16:P:83:GLU:N	2.72	0.46
17:Q:59:ILE:HD13	17:Q:59:ILE:HA	1.60	0.46
1:A:1190:G:H5'	3:C:176:HIS:NE2	2.30	0.46
5:E:82:VAL:HG21	5:E:138:ALA:HA	1.97	0.46
12:L:45:PRO:HB3	12:L:92:ASP:HB3	1.97	0.46
1:A:358:U:H2'	1:A:359:U:C6	2.51	0.46
1:A:556:C:OP2	12:L:20:LYS:NZ	2.39	0.46
1:A:737:A:H2'	1:A:738:C:C6	2.50	0.46
1:A:1239:A:C4	1:A:1298:C:N4	2.84	0.46
1:A:1298:C:H4'	1:A:1299:A:O4'	2.16	0.46
18:R:46:GLU:CD	18:R:46:GLU:H	2.24	0.46
1:A:160:A:H2'	1:A:161:A:O4'	2.16	0.46
1:A:1018:C:H2'	1:A:1019:C:O4'	2.16	0.46
1:A:1126:U:HO2'	1:A:1127:G:P	2.38	0.46
1:A:1226:C:H4'	1:A:1227:A:OP1	2.15	0.46
2:B:61:LEU:HD22	2:B:61:LEU:HA	1.80	0.46
3:C:19:GLU:HG2	3:C:40:ARG:HH21	1.80	0.46
3:C:64:VAL:HG23	3:C:99:VAL:HG11	1.97	0.46
11:K:110:ASP:HB2	18:R:88:LYS:HE2	1.97	0.46
13:M:19:LEU:HD23	13:M:19:LEU:HA	1.79	0.46
16:P:57:ARG:HG2	16:P:79:VAL:HG13	1.98	0.46
1:A:409:G:N3	1:A:409:G:C2'	2.79	0.46
1:A:646:U:H2'	1:A:647:C:C6	2.51	0.46
1:A:1279:A:O2'	1:A:1282:C:N4	2.49	0.46
1:A:1300:G:O2'	1:A:1301:U:P	2.74	0.46
2:B:47:THR:HA	2:B:202:PRO:HG2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:40:ARG:HB3	3:C:44:GLU:OE2	2.16	0.46
3:C:130:VAL:HG12	3:C:134:ILE:HD11	1.98	0.46
8:H:4:ASP:OD2	8:H:85:ARG:HD2	2.15	0.46
1:A:977:A:H2'	1:A:978:A:H5''	1.97	0.45
1:A:1283:G:H2'	1:A:1284:C:C6	2.51	0.45
1:A:1505:G:C8	1:A:1505:G:H3'	2.51	0.45
5:E:21:ALA:O	5:E:23:GLY:N	2.49	0.45
1:A:1190:G:O2'	1:A:1191:A:P	2.74	0.45
16:P:74:LEU:O	16:P:79:VAL:HG23	2.16	0.45
1:A:407:G:H1	1:A:496:A:HO2'	1.58	0.45
1:A:1406:U:H2'	1:A:1407:C:O4'	2.16	0.45
1:A:1497:G:H2'	1:A:1498:U:H5'	1.97	0.45
4:D:111:ALA:HB3	4:D:117:ALA:HB2	1.98	0.45
1:A:17:U:H2'	1:A:18:C:C6	2.51	0.45
1:A:582:U:OP2	1:A:758:G:N2	2.42	0.45
1:A:1131:G:H2'	1:A:1132:C:C6	2.52	0.45
2:B:141:GLU:O	2:B:145:LEU:HG	2.16	0.45
5:E:80:ILE:HD12	5:E:91:LEU:HB2	1.98	0.45
15:O:77:ARG:O	15:O:80:ALA:HB3	2.17	0.45
1:A:478:A:O2'	1:A:479:C:H5'	2.17	0.45
5:E:75:THR:OG1	5:E:76:ILE:N	2.50	0.45
1:A:443:C:H2'	1:A:444:C:H6	1.82	0.45
1:A:1030:C:O2'	1:A:1032:G:N1	2.48	0.45
1:A:1117:G:O3'	9:I:104:ARG:HD2	2.17	0.45
2:B:17:PHE:CG	2:B:17:PHE:O	2.69	0.45
7:G:9:VAL:HG13	7:G:94:ARG:NH2	2.21	0.45
14:N:29:ARG:HB3	14:N:40:CYS:HB3	1.99	0.45
15:O:5:LYS:HE3	15:O:5:LYS:HB2	1.66	0.45
1:A:410:G:O4'	1:A:410:G:P	2.75	0.45
1:A:1057:G:H5''	3:C:154:SER:HB2	1.99	0.45
1:A:1448:C:H2'	1:A:1449:C:H6	1.82	0.45
2:B:51:LEU:HA	2:B:54:THR:HB	1.98	0.45
10:J:17:ASP:O	10:J:21:GLN:HB2	2.17	0.45
11:K:33:THR:HA	11:K:39:PRO:HA	1.97	0.45
17:Q:6:LEU:HB3	17:Q:23:VAL:HG11	1.97	0.45
17:Q:99:SER:OG	17:Q:100:LYS:N	2.50	0.45
1:A:403:C:O2'	4:D:122:ARG:NH1	2.49	0.45
10:J:94:VAL:HG12	10:J:96:ILE:HG13	1.98	0.45
11:K:48:ILE:HG13	11:K:63:LEU:CB	2.46	0.45
17:Q:66:SER:OG	17:Q:69:LYS:HB3	2.17	0.45
1:A:462:G:N2	16:P:80:PHE:HB2	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:195:ASP:O	8:H:74:PRO:HG3	2.17	0.45
4:D:60:GLU:HG2	4:D:202:LEU:HD12	1.98	0.45
5:E:95:ALA:O	5:E:98:THR:OG1	2.26	0.45
9:I:28:VAL:HG12	9:I:63:ILE:HB	1.98	0.45
9:I:31:GLN:HB3	9:I:35:GLU:HB3	1.97	0.45
16:P:75:ARG:HH11	16:P:82:GLN:HE22	1.64	0.45
1:A:524:G:H2'	1:A:525:C:C6	2.52	0.45
1:A:881:G:P	12:L:12:ARG:HH22	2.39	0.45
1:A:1497:G:C2'	1:A:1498:U:H5'	2.47	0.45
7:G:15:ASP:HB3	7:G:20:ASP:H	1.82	0.45
10:J:39:PRO:HA	10:J:70:ARG:NH1	2.32	0.45
14:N:24:CYS:HB2	14:N:40:CYS:HB3	1.62	0.45
1:A:409:G:H3'	1:A:430:A:N6	2.32	0.44
1:A:609:A:C5	1:A:610:G:C8	3.06	0.44
1:A:1371:G:O3'	9:I:69:GLY:HA3	2.17	0.44
7:G:57:GLU:O	7:G:61:VAL:HG23	2.18	0.44
11:K:48:ILE:HA	11:K:48:ILE:HD13	1.67	0.44
17:Q:29:HIS:CG	17:Q:30:PRO:HD2	2.53	0.44
1:A:1106:G:H5''	3:C:172:ARG:HG2	1.99	0.44
1:A:1136:U:H5''	1:A:1137:C:C5	2.52	0.44
4:D:11:LEU:HD13	4:D:66:ARG:HD3	1.98	0.44
7:G:66:VAL:HG12	7:G:70:LYS:HE3	1.98	0.44
9:I:53:VAL:HG22	9:I:96:LEU:HD11	1.99	0.44
12:L:46:LYS:O	12:L:47:LYS:C	2.61	0.44
19:S:3:ARG:HB3	19:S:4:SER:H	1.70	0.44
1:A:175:C:H4'	20:T:25:ARG:HD2	1.99	0.44
1:A:1118:C:P	9:I:104:ARG:HH11	2.40	0.44
1:A:1125:U:H3	10:J:5:ARG:HH21	1.64	0.44
9:I:10:ARG:HD3	9:I:75:ASP:CB	2.41	0.44
9:I:99:LEU:HB2	9:I:101:PHE:HD2	1.82	0.44
10:J:39:PRO:O	10:J:40:LEU:HB2	2.17	0.44
14:N:11:LYS:HG2	14:N:13:THR:HB	1.98	0.44
17:Q:85:VAL:O	17:Q:89:LEU:N	2.50	0.44
1:A:229:U:H2'	1:A:230:G:O4'	2.17	0.44
1:A:334:C:H2'	1:A:335:C:H6	1.83	0.44
1:A:1305:G:H2'	1:A:1331:G:N2	2.32	0.44
1:A:1307:U:H2'	1:A:1308:U:C6	2.52	0.44
2:B:200:ILE:H	2:B:200:ILE:HG13	1.54	0.44
4:D:88:VAL:O	4:D:92:VAL:HG23	2.18	0.44
5:E:69:VAL:HA	5:E:70:PRO:HD3	1.85	0.44
6:F:44:GLY:HA2	6:F:59:TYR:CE2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:26:PHE:HB2	7:G:101:LEU:HD22	1.99	0.44
11:K:33:THR:HG22	11:K:39:PRO:HA	1.98	0.44
12:L:10:LEU:HD23	12:L:10:LEU:HA	1.71	0.44
17:Q:3:LYS:HB3	17:Q:61:GLU:HB3	1.98	0.44
19:S:43:GLU:CD	19:S:43:GLU:H	2.25	0.44
1:A:370:C:C2'	1:A:371:G:H5'	2.48	0.44
1:A:1130:A:C2	1:A:1146:A:C4	3.05	0.44
7:G:149:ARG:HG2	7:G:152:ALA:HB2	1.98	0.44
1:A:580:U:H2'	1:A:581:G:C1'	2.47	0.44
1:A:812:C:OP1	1:A:903:G:H1'	2.18	0.44
7:G:20:ASP:OD2	7:G:63:LYS:NZ	2.28	0.44
11:K:40:ILE:HG23	11:K:75:TYR:CE2	2.52	0.44
19:S:4:SER:O	19:S:5:LEU:HB2	2.18	0.44
1:A:428:G:O2'	1:A:429:U:OP2	2.29	0.44
1:A:1098:C:H2'	1:A:1099:G:O4'	2.17	0.44
1:A:1378:C:C5	1:A:1379:G:C8	3.06	0.44
3:C:179:ARG:O	3:C:179:ARG:HG2	2.18	0.44
20:T:83:ARG:HB3	20:T:87:LYS:NZ	2.33	0.44
1:A:309:G:H1'	1:A:608:A:C2	2.52	0.44
1:A:334:C:H2'	1:A:335:C:C6	2.53	0.44
1:A:945:G:C2	1:A:946:A:C8	3.06	0.44
1:A:972:C:O5'	10:J:57:LYS:HD3	2.17	0.44
1:A:1420:C:H2'	1:A:1421:G:H8	1.82	0.44
14:N:9:LYS:O	14:N:11:LYS:N	2.50	0.44
15:O:57:LEU:HD12	15:O:57:LEU:HA	1.77	0.44
21:V:6:ARG:H	21:V:6:ARG:HG3	1.40	0.44
1:A:560:U:H5'	1:A:566:G:N2	2.33	0.44
1:A:652:U:O4	1:A:752:G:O2'	2.27	0.44
1:A:804:U:H5''	1:A:805:C:OP2	2.17	0.44
1:A:1060:C:P	14:N:45:ARG:HH22	2.41	0.44
1:A:1308:U:H2'	1:A:1309:G:C8	2.53	0.44
4:D:61:LYS:HE2	4:D:206:PHE:CE2	2.53	0.44
15:O:2:PRO:C	15:O:38:ARG:HH12	2.25	0.44
2:B:20:GLU:O	2:B:39:ILE:HG23	2.18	0.43
2:B:30:ARG:O	2:B:46:LYS:NZ	2.36	0.43
2:B:90:MET:HA	2:B:91:PRO:HD3	1.71	0.43
5:E:144:THR:O	5:E:147:ASP:N	2.51	0.43
6:F:33:TYR:HB2	6:F:75:LEU:HD23	2.00	0.43
8:H:3:THR:O	8:H:5:PRO:HD3	2.18	0.43
11:K:18:ARG:HB2	11:K:33:THR:OG1	2.18	0.43
14:N:23:ARG:NH1	14:N:30:ALA:HB2	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:Q:24:GLU:HG3	17:Q:39:SER:HB3	2.00	0.43
1:A:520:A:OP2	12:L:51:ALA:HB1	2.18	0.43
1:A:1126:U:O2'	1:A:1127:G:OP1	2.28	0.43
3:C:112:SER:HB3	3:C:115:LEU:HD12	2.00	0.43
1:A:1091:U:O2	1:A:1093:A:C8	2.71	0.43
1:A:1256:A:H2	1:A:1277:C:N4	2.16	0.43
1:A:1468:A:H2'	1:A:1469:G:O4'	2.18	0.43
3:C:167:TRP:HB3	3:C:168:ALA:H	1.54	0.43
6:F:23:LYS:HB3	6:F:23:LYS:HE2	1.77	0.43
21:V:5:ASP:O	21:V:11:GLY:HA3	2.18	0.43
1:A:651:C:O2'	1:A:652:U:H5'	2.18	0.43
1:A:1123:A:H4'	10:J:36:GLY:HA3	2.00	0.43
1:A:1166:G:N2	1:A:1169:A:OP2	2.51	0.43
2:B:71:VAL:HG22	2:B:164:VAL:HA	2.00	0.43
3:C:85:ARG:HH22	3:C:88:ARG:HH11	1.67	0.43
7:G:111:ARG:NH1	7:G:113:GLU:OE2	2.51	0.43
7:G:115:ARG:H	7:G:115:ARG:HG3	1.43	0.43
12:L:43:VAL:HG12	12:L:93:LEU:HD22	2.00	0.43
18:R:76:LEU:HA	18:R:76:LEU:HD23	1.62	0.43
1:A:925:G:C2	1:A:927:G:C8	3.07	0.43
1:A:954:G:H2'	1:A:955:U:O4'	2.18	0.43
1:A:1152:A:H5''	10:J:13:HIS:ND1	2.33	0.43
1:A:1167:A:C6	1:A:1168:A:C6	3.06	0.43
5:E:6:PHE:HE1	5:E:36:ASP:HB3	1.83	0.43
5:E:71:LEU:HD11	5:E:114:GLY:HA3	2.01	0.43
8:H:39:LEU:HD12	8:H:39:LEU:HA	1.73	0.43
10:J:90:LEU:N	10:J:91:PRO:HD2	2.23	0.43
2:B:84:GLU:OE2	2:B:233:SER:OG	2.34	0.43
5:E:11:ILE:HD12	5:E:11:ILE:HG23	1.78	0.43
6:F:53:ALA:C	6:F:55:ASP:H	2.26	0.43
7:G:48:LYS:HE2	7:G:48:LYS:HB3	1.83	0.43
7:G:113:GLU:HG2	7:G:119:ARG:HG2	2.00	0.43
12:L:33:ARG:HA	12:L:33:ARG:HD2	1.35	0.43
17:Q:15:MET:HE1	17:Q:43:LEU:HD22	2.01	0.43
1:A:62:U:H2'	1:A:63:C:C6	2.54	0.43
1:A:490:G:H2'	1:A:491:G:C8	2.51	0.43
1:A:1057:G:H1	1:A:1203:C:H42	1.65	0.43
2:B:32:ILE:HG21	2:B:40:HIS:ND1	2.34	0.43
3:C:11:ARG:C	3:C:13:GLY:H	2.26	0.43
9:I:69:GLY:O	9:I:73:GLN:HG3	2.18	0.43
11:K:54:ARG:H	11:K:54:ARG:HG2	1.38	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:M:15:VAL:HA	13:M:18:ALA:HB3	2.00	0.43
13:M:87:TYR:O	13:M:91:ARG:HG2	2.19	0.43
20:T:43:LEU:HB3	20:T:52:ALA:HB2	2.01	0.43
1:A:390:C:H2'	1:A:391:G:C8	2.54	0.43
1:A:947:G:H2'	1:A:948:C:O4'	2.18	0.43
1:A:410:G:C2	1:A:411:A:H1'	2.53	0.43
3:C:130:VAL:CG2	3:C:157:ILE:HG23	2.47	0.43
1:A:865:A:H5'	1:A:1078:U:C5	2.54	0.43
1:A:1338:G:C6	1:A:1339:A:C6	3.07	0.43
1:A:1491:G:N7	25:A:1693:PAR:O53	2.45	0.43
6:F:55:ASP:CB	6:F:86:ARG:HH12	2.32	0.43
9:I:120:ARG:O	9:I:122:ALA:N	2.51	0.43
1:A:190(A):C:H42	1:A:190(H):G:H1	1.66	0.42
1:A:452:A:OP1	16:P:43:LYS:NZ	2.37	0.42
1:A:528:C:H5'	1:A:535:A:N6	2.34	0.42
1:A:1220:G:H2'	1:A:1221:G:C8	2.54	0.42
2:B:117:GLU:O	2:B:120:ALA:HB3	2.19	0.42
4:D:70:ILE:HD11	4:D:100:ARG:CZ	2.49	0.42
5:E:105:VAL:HB	5:E:106:PRO:HD3	2.01	0.42
7:G:51:GLN:OE1	7:G:58:PRO:HD3	2.18	0.42
12:L:92:ASP:O	12:L:94:PRO:HD3	2.18	0.42
13:M:122:LYS:HA	13:M:122:LYS:HD2	1.79	0.42
17:Q:10:VAL:HG13	17:Q:19:VAL:HB	2.01	0.42
17:Q:53:LEU:HD11	17:Q:85:VAL:HG11	2.01	0.42
1:A:60:A:H4'	1:A:61:G:O5'	2.19	0.42
1:A:190(B):C:H5'	1:A:190(C):C:OP2	2.18	0.42
1:A:407:G:HO2'	1:A:431:A:P	2.41	0.42
1:A:660:G:C2	1:A:746:A:C2	3.08	0.42
10:J:81:THR:O	10:J:85:LEU:HG	2.19	0.42
13:M:23:TYR:HB3	13:M:67:GLU:H	1.84	0.42
1:A:1283:G:H2'	1:A:1284:C:H6	1.85	0.42
6:F:67:MET:HE1	6:F:75:LEU:HG	2.01	0.42
1:A:407:G:O2'	1:A:431:A:P	2.77	0.42
1:A:1010:G:H2'	1:A:1011:G:C8	2.54	0.42
2:B:16:HIS:HB3	2:B:17:PHE:H	1.30	0.42
4:D:150:GLU:OE1	4:D:150:GLU:HA	2.19	0.42
5:E:11:ILE:HD13	5:E:11:ILE:HA	1.76	0.42
7:G:23:VAL:O	7:G:27:ILE:HG13	2.18	0.42
11:K:13:GLN:HA	11:K:75:TYR:O	2.20	0.42
11:K:66:LEU:HG	11:K:97:ALA:HB1	2.02	0.42
11:K:108:ILE:O	18:R:87:ARG:N	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:R:39:VAL:O	18:R:42:ARG:HB2	2.18	0.42
1:A:427:U:OP1	4:D:13:ARG:NH2	2.46	0.42
1:A:875:C:O2'	8:H:14:ARG:NH1	2.53	0.42
1:A:1256:A:H5''	1:A:1257:U:OP1	2.19	0.42
1:A:1286:A:C8	1:A:1287:A:H4'	2.55	0.42
1:A:1391:U:H2'	1:A:1392:G:C8	2.54	0.42
2:B:166:ASP:OD2	2:B:169:LYS:HB2	2.19	0.42
2:B:218:ALA:O	2:B:221:LEU:HB3	2.19	0.42
14:N:39:LEU:HD22	14:N:43:CYS:HB2	2.02	0.42
16:P:65:GLN:HA	16:P:66:PRO:HD2	1.82	0.42
1:A:151:A:H2'	1:A:152:A:O4'	2.19	0.42
1:A:376:G:OP2	16:P:67:THR:HG21	2.19	0.42
1:A:476:G:O6	1:A:479:C:N4	2.52	0.42
1:A:946:A:H2'	1:A:947:G:C8	2.54	0.42
1:A:949:A:OP2	13:M:106:ASN:ND2	2.52	0.42
1:A:1201:A:H4'	1:A:1202:G:O5'	2.19	0.42
1:A:1216:G:H2'	1:A:1217:C:H6	1.85	0.42
5:E:144:THR:HG22	5:E:147:ASP:OD1	2.20	0.42
8:H:51:VAL:HG21	8:H:60:ARG:NH1	2.34	0.42
12:L:47:LYS:CB	12:L:48:PRO:HD3	2.42	0.42
14:N:39:LEU:HD23	14:N:39:LEU:HA	1.89	0.42
18:R:40:LEU:HD23	18:R:40:LEU:HA	1.76	0.42
1:A:501:C:H2'	1:A:502:G:H8	1.82	0.42
1:A:901:A:C5	1:A:902:G:H1'	2.55	0.42
1:A:965:A:C2	1:A:969:A:C2	3.08	0.42
1:A:967:C:O3'	9:I:128:ARG:NH2	2.53	0.42
1:A:980:C:H3'	1:A:981:U:H6	1.84	0.42
1:A:1020:U:H2'	1:A:1021:G:C8	2.55	0.42
1:A:1316:G:H4'	14:N:18:VAL:CG1	2.49	0.42
3:C:20:SER:O	14:N:54:PRO:HB3	2.19	0.42
4:D:64:LEU:HG	4:D:198:VAL:HG11	2.01	0.42
10:J:7:LYS:HA	10:J:7:LYS:HD2	1.86	0.42
14:N:23:ARG:O	14:N:25:VAL:N	2.52	0.42
1:A:1118:C:H1'	1:A:1179:A:C4	2.54	0.42
4:D:50:ARG:HA	4:D:51:PRO:HD3	1.80	0.42
9:I:4:TYR:CE1	9:I:88:TYR:HD1	2.38	0.42
9:I:48:GLU:N	9:I:49:PRO:HD2	2.35	0.42
1:A:1190:G:OP1	3:C:5:ILE:HG13	2.20	0.42
4:D:187:ARG:NH1	4:D:188:LEU:H	2.17	0.42
7:G:40:ALA:HB1	9:I:41:VAL:HG21	2.02	0.42
1:A:28:G:O2'	1:A:296:U:OP1	2.33	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:U:OP1	1:A:385:C:O2'	2.36	0.42
1:A:1007:C:H1'	1:A:1023:G:N2	2.35	0.42
1:A:1129:C:H1'	1:A:1132:C:C5	2.51	0.42
2:B:71:VAL:O	2:B:165:VAL:HG23	2.19	0.42
2:B:128:GLU:HA	2:B:135:GLN:OE1	2.20	0.42
2:B:236:TYR:C	2:B:238:LEU:N	2.78	0.42
3:C:191:THR:HG21	3:C:193:TYR:CZ	2.54	0.42
6:F:25:ILE:HD12	6:F:82:ARG:HD2	2.01	0.42
12:L:7:ILE:O	12:L:11:VAL:HG23	2.20	0.42
1:A:462:G:C6	1:A:463:A:H2	2.37	0.41
1:A:500:G:C6	1:A:501:C:C4	3.08	0.41
1:A:975:A:O2'	14:N:32:SER:HA	2.20	0.41
1:A:987:G:O5'	1:A:987:G:H8	2.02	0.41
2:B:44:LEU:HD23	2:B:44:LEU:HA	1.84	0.41
1:A:355:C:H5'	1:A:389:A:OP2	2.20	0.41
1:A:1291:G:C6	1:A:1292:U:C4	3.08	0.41
1:A:1379:G:OP2	7:G:6:ARG:NE	2.52	0.41
2:B:48:MET:HE3	2:B:48:MET:HB3	1.67	0.41
2:B:80:ILE:HG13	2:B:238:LEU:HD23	2.01	0.41
9:I:81:ILE:O	9:I:85:LEU:HB2	2.19	0.41
10:J:11:PHE:HB3	14:N:55:GLY:HA3	2.02	0.41
10:J:24:VAL:O	10:J:28:ARG:N	2.53	0.41
15:O:25:THR:CG2	15:O:70:LEU:HD23	2.50	0.41
17:Q:86:GLU:O	17:Q:90:ILE:HG13	2.20	0.41
1:A:1513:A:H2'	1:A:1514:C:C6	2.54	0.41
9:I:50:LEU:HD23	9:I:50:LEU:HA	1.87	0.41
10:J:34:VAL:HG12	10:J:35:SER:N	2.34	0.41
12:L:7:ILE:HA	12:L:7:ILE:HD13	1.61	0.41
1:A:390:C:O5'	1:A:390:C:H6	2.03	0.41
1:A:401:C:H2'	1:A:402:G:C8	2.55	0.41
1:A:743:U:H2'	1:A:744:C:C6	2.55	0.41
4:D:70:ILE:HD11	4:D:100:ARG:NE	2.34	0.41
4:D:176:LEU:HA	4:D:183:GLY:CA	2.50	0.41
6:F:19:LEU:O	6:F:23:LYS:HG3	2.20	0.41
12:L:125:PRO:HG2	12:L:127:GLU:OE2	2.20	0.41
1:A:434:U:O2	1:A:434:U:H2'	2.20	0.41
1:A:443:C:H2'	1:A:444:C:C6	2.55	0.41
1:A:581:G:C5	1:A:758:G:C5	3.09	0.41
1:A:792:A:H1'	1:A:794:A:N7	2.35	0.41
1:A:922:G:C6	1:A:923:A:C6	3.08	0.41
3:C:114:PRO:HD3	3:C:183:ASP:OD2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:176:LEU:HD21	4:D:178:VAL:HG22	2.02	0.41
6:F:79:LEU:HD23	6:F:79:LEU:HA	1.84	0.41
7:G:152:ALA:O	7:G:155:ARG:NH2	2.53	0.41
7:G:155:ARG:NH1	7:G:155:ARG:HA	2.35	0.41
9:I:32:ASP:O	9:I:34:ASN:N	2.53	0.41
12:L:90:VAL:HG11	12:L:93:LEU:HG	2.02	0.41
16:P:10:GLY:HA3	16:P:14:ASN:O	2.20	0.41
18:R:23:LYS:HD2	18:R:57:GLY:O	2.20	0.41
1:A:580:U:H2'	1:A:581:G:O4'	2.20	0.41
6:F:33:TYR:HA	6:F:71:ARG:HH21	1.86	0.41
13:M:72:ALA:O	13:M:76:ALA:N	2.45	0.41
14:N:11:LYS:C	14:N:13:THR:H	2.28	0.41
20:T:61:SER:O	20:T:62:LEU:C	2.64	0.41
1:A:243:A:C2	1:A:246:A:C8	3.08	0.41
1:A:352:C:H5''	1:A:352:C:H6	1.85	0.41
1:A:397:A:H5'	1:A:398:C:P	2.60	0.41
1:A:1059:C:O3'	14:N:45:ARG:NH2	2.53	0.41
1:A:1060:C:H2'	1:A:1061:G:H8	1.86	0.41
1:A:1212:U:H4'	1:A:1213:A:O5'	2.20	0.41
1:A:1338:G:H2'	1:A:1339:A:C8	2.56	0.41
3:C:155:GLY:HA3	3:C:163:ALA:HB1	2.01	0.41
8:H:75:ARG:HA	8:H:76:PRO:HD3	1.81	0.41
9:I:19:LEU:HD23	9:I:19:LEU:HA	1.89	0.41
1:A:280:C:O2	17:Q:38:ARG:HG3	2.21	0.41
1:A:911:U:OP1	12:L:95:GLY:HA2	2.21	0.41
1:A:1168:A:H2'	1:A:1169:A:C8	2.56	0.41
1:A:1234:C:H1'	1:A:1364:U:O2	2.20	0.41
2:B:121:LEU:HD23	2:B:121:LEU:HA	1.87	0.41
2:B:127:ILE:HG13	2:B:128:GLU:OE2	2.21	0.41
2:B:164:VAL:HG11	2:B:167:PRO:HA	2.02	0.41
3:C:56:ASP:O	3:C:66:VAL:HA	2.21	0.41
4:D:154:ASN:HA	4:D:159:ARG:HE	1.86	0.41
5:E:31:LEU:HD23	5:E:31:LEU:HA	1.68	0.41
13:M:96:LEU:O	13:M:110:ARG:NH1	2.54	0.41
15:O:31:LEU:HD12	15:O:31:LEU:HA	1.85	0.41
1:A:273:A:N6	1:A:274:A:N6	2.69	0.41
1:A:460:A:H2'	1:A:461:C:H5'	2.03	0.41
1:A:673:G:N2	1:A:674:G:C2	2.89	0.41
1:A:683:G:H2'	1:A:684:A:C8	2.56	0.41
1:A:828:A:H2'	1:A:829:G:O4'	2.21	0.41
1:A:1329:A:H2'	1:A:1330:U:O4'	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:174:PRO:HB2	3:C:177:THR:HB	2.03	0.41
7:G:26:PHE:CE2	7:G:30:ILE:HD11	2.56	0.41
11:K:19:ALA:HB2	11:K:80:VAL:HG21	2.03	0.41
12:L:46:LYS:HB2	12:L:47:LYS:H	1.55	0.41
12:L:53:ARG:HG3	12:L:93:LEU:HD21	2.02	0.41
12:L:127:GLU:H	12:L:127:GLU:HG2	1.58	0.41
14:N:12:ARG:HA	14:N:12:ARG:HD3	1.52	0.41
16:P:71:ARG:HD3	16:P:75:ARG:HH21	1.85	0.41
20:T:43:LEU:HD23	20:T:43:LEU:HA	1.93	0.41
1:A:476:G:H2'	1:A:476:G:N3	2.36	0.41
1:A:1346:A:H5''	9:I:120:ARG:NH1	2.35	0.41
1:A:1448:C:H2'	1:A:1449:C:C6	2.56	0.41
3:C:61:ALA:O	3:C:63:ASN:N	2.48	0.41
3:C:188:LEU:HD22	3:C:188:LEU:HA	1.86	0.41
6:F:3:ARG:HE	6:F:3:ARG:HB3	1.72	0.41
10:J:39:PRO:O	10:J:69:ASN:O	2.39	0.41
13:M:29:ARG:NH2	13:M:64:TRP:CD1	2.89	0.41
18:R:40:LEU:HB3	18:R:79:LEU:HD11	2.03	0.41
20:T:57:ARG:NH1	20:T:102:GLY:HA3	2.36	0.41
20:T:92:LEU:HD23	20:T:92:LEU:HA	1.71	0.41
4:D:18:LYS:NZ	4:D:20:TYR:OH	2.31	0.40
1:A:613:C:H2'	1:A:614:A:C8	2.56	0.40
1:A:779:C:H2'	1:A:780:A:O4'	2.20	0.40
1:A:826:C:H2'	1:A:827:U:C6	2.56	0.40
1:A:1057:G:H5''	3:C:154:SER:CB	2.51	0.40
1:A:1288:A:H8	1:A:1288:A:O5'	2.03	0.40
2:B:84:GLU:OE1	2:B:216:SER:HA	2.21	0.40
3:C:11:ARG:C	3:C:13:GLY:N	2.78	0.40
5:E:79:GLU:HG2	5:E:92:LYS:HG3	2.03	0.40
6:F:68:PRO:HB2	6:F:71:ARG:HG2	2.03	0.40
8:H:4:ASP:CG	8:H:85:ARG:HH11	2.25	0.40
13:M:126:LYS:HB2	13:M:126:LYS:HE3	1.90	0.40
14:N:9:LYS:C	14:N:11:LYS:H	2.29	0.40
14:N:41:ARG:HG2	14:N:41:ARG:HH11	1.86	0.40
18:R:53:ARG:C	18:R:55:ARG:H	2.28	0.40
1:A:157:G:C2	1:A:158:G:C8	3.09	0.40
1:A:442:C:H42	1:A:492:G:H1	1.68	0.40
1:A:448:A:H2'	1:A:449:C:C6	2.56	0.40
1:A:1248:A:H2'	1:A:1249:C:C6	2.56	0.40
1:A:1251:A:H2'	1:A:1252:A:O4'	2.22	0.40
2:B:112:VAL:HG22	2:B:149:LEU:HD13	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:121:VAL:HG12	4:D:134:ASP:HA	2.03	0.40
9:I:98:PRO:HB2	9:I:99:LEU:HD22	2.04	0.40
12:L:124:LYS:HA	12:L:125:PRO:HD2	1.88	0.40
1:A:406:G:C8	1:A:496:A:C6	3.09	0.40
1:A:767:A:H2'	1:A:768:A:O4'	2.21	0.40
1:A:889:A:H5'	1:A:891:U:H1'	2.04	0.40
1:A:928:G:C5'	1:A:1533:C:H5'	2.51	0.40
4:D:78:LEU:HD23	4:D:78:LEU:HA	1.73	0.40
15:O:10:LYS:HE3	15:O:11:VAL:HA	2.04	0.40
1:A:112:G:O2'	1:A:113:G:H5'	2.22	0.40
1:A:1352:C:H2'	1:A:1353:G:C8	2.57	0.40
1:A:1441:G:H4'	1:A:1442:G:C5	2.57	0.40
3:C:70:VAL:HG12	3:C:72:LYS:H	1.86	0.40
4:D:120:LEU:HD23	4:D:125:HIS:CD2	2.56	0.40
6:F:30:LEU:HB3	6:F:35:ALA:HB3	2.04	0.40
8:H:7:ALA:CA	8:H:85:ARG:HG3	2.50	0.40
9:I:127:LYS:HD2	9:I:127:LYS:HA	1.78	0.40
10:J:25:GLU:C	10:J:27:ALA:H	2.29	0.40
14:N:9:LYS:O	14:N:9:LYS:HE2	2.21	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%)	184 (79%)	34 (15%)	14 (6%)	1	9
3	C	204/239 (85%)	166 (81%)	25 (12%)	13 (6%)	1	8
4	D	206/209 (99%)	182 (88%)	17 (8%)	7 (3%)	3	19
5	E	148/162 (91%)	134 (90%)	12 (8%)	2 (1%)	9	34
6	F	99/101 (98%)	89 (90%)	9 (9%)	1 (1%)	12	41

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
7	G	153/156 (98%)	137 (90%)	12 (8%)	4 (3%)	4	23
8	H	136/138 (99%)	129 (95%)	7 (5%)	0	100	100
9	I	125/128 (98%)	102 (82%)	11 (9%)	12 (10%)	0	3
10	J	96/105 (91%)	68 (71%)	15 (16%)	13 (14%)	0	1
11	K	117/129 (91%)	104 (89%)	12 (10%)	1 (1%)	14	43
12	L	122/132 (92%)	100 (82%)	16 (13%)	6 (5%)	1	12
13	M	123/126 (98%)	98 (80%)	20 (16%)	5 (4%)	2	15
14	N	58/61 (95%)	46 (79%)	9 (16%)	3 (5%)	1	11
15	O	86/89 (97%)	76 (88%)	9 (10%)	1 (1%)	10	37
16	P	81/88 (92%)	69 (85%)	9 (11%)	3 (4%)	2	17
17	Q	102/105 (97%)	86 (84%)	11 (11%)	5 (5%)	1	12
18	R	71/88 (81%)	57 (80%)	11 (16%)	3 (4%)	2	15
19	S	78/93 (84%)	56 (72%)	15 (19%)	7 (9%)	0	4
20	T	97/106 (92%)	78 (80%)	12 (12%)	7 (7%)	1	6
21	V	22/27 (82%)	18 (82%)	4 (18%)	0	100	100
All	All	2356/2538 (93%)	1979 (84%)	270 (12%)	107 (4%)	2	13

All (107) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	91	LEU
4	D	23	GLY
6	F	72	VAL
9	I	33	PHE
10	J	34	VAL
10	J	76	ASN
10	J	90	LEU
12	L	28	LYS
12	L	41	ARG
12	L	47	LYS
12	L	48	PRO
12	L	73	GLU
13	M	12	ASN
13	M	14	ARG
14	N	10	ALA
14	N	24	CYS
16	P	82	GLN

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Mol	Chain	Res	Type
20	T	99	LEU
2	B	207	ALA
2	B	235	SER
2	B	236	TYR
3	C	12	LEU
3	C	154	SER
4	D	4	TYR
4	D	25	ARG
4	D	69	GLY
5	E	16	THR
7	G	6	ARG
7	G	147	ALA
9	I	7	THR
9	I	29	ASN
9	I	38	GLN
9	I	101	PHE
9	I	127	LYS
10	J	72	VAL
12	L	27	LEU
14	N	25	VAL
17	Q	97	SER
18	R	21	LYS
18	R	87	ARG
19	S	9	VAL
19	S	43	GLU
20	T	45	GLN
20	T	102	GLY
2	B	9	GLU
2	B	95	GLN
2	B	221	LEU
3	C	130	VAL
3	C	146	ALA
3	C	168	ALA
3	C	206	GLU
9	I	55	ALA
9	I	121	ARG
10	J	24	VAL
10	J	39	PRO
10	J	55	LYS
10	J	85	LEU
10	J	86	MET
11	K	117	ASN

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Mol	Chain	Res	Type
13	M	106	ASN
17	Q	49	GLU
17	Q	74	LEU
17	Q	79	SER
17	Q	81	ARG
19	S	47	HIS
20	T	73	HIS
20	T	96	GLY
20	T	98	PRO
2	B	8	LYS
2	B	11	LEU
2	B	17	PHE
3	C	4	LYS
3	C	60	ALA
3	C	62	ASP
3	C	205	GLY
4	D	156	GLU
9	I	41	VAL
9	I	56	LEU
10	J	40	LEU
10	J	73	ASP
10	J	89	ASP
13	M	122	LYS
15	O	85	LEU
16	P	79	VAL
2	B	15	VAL
2	B	20	GLU
5	E	22	GLY
7	G	4	ARG
9	I	54	ASP
10	J	61	GLU
13	M	86	CYS
19	S	25	LYS
19	S	66	MET
2	B	29	ALA
4	D	5	ILE
18	R	25	THR
19	S	30	LEU
19	S	42	PRO
16	P	10	GLY
2	B	131	PRO
7	G	58	PRO

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Mol	Chain	Res	Type
9	I	100	GLY
3	C	51	GLY
3	C	66	VAL
4	D	36	ARG
2	B	227	GLY
20	T	97	ALA

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%)	173 (86%)	29 (14%)	3	14
3	C	160/188 (85%)	130 (81%)	30 (19%)	1	7
4	D	180/181 (99%)	157 (87%)	23 (13%)	4	18
5	E	115/123 (94%)	94 (82%)	21 (18%)	2	7
6	F	90/90 (100%)	76 (84%)	14 (16%)	2	12
7	G	126/127 (99%)	107 (85%)	19 (15%)	3	13
8	H	119/119 (100%)	103 (87%)	16 (13%)	4	16
9	I	98/99 (99%)	90 (92%)	8 (8%)	10	36
10	J	87/92 (95%)	70 (80%)	17 (20%)	1	6
11	K	90/99 (91%)	80 (89%)	10 (11%)	6	23
12	L	104/109 (95%)	88 (85%)	16 (15%)	2	12
13	M	100/101 (99%)	85 (85%)	15 (15%)	3	13
14	N	49/50 (98%)	42 (86%)	7 (14%)	3	14
15	O	79/80 (99%)	61 (77%)	18 (23%)	1	4
16	P	72/74 (97%)	62 (86%)	10 (14%)	3	15
17	Q	96/97 (99%)	83 (86%)	13 (14%)	4	16
18	R	64/77 (83%)	58 (91%)	6 (9%)	8	31
19	S	71/80 (89%)	55 (78%)	16 (22%)	1	4
20	T	75/82 (92%)	62 (83%)	13 (17%)	2	9

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
21	V	19/22 (86%)	15 (79%)	4 (21%)	1	5
All	All	1996/2110 (95%)	1691 (85%)	305 (15%)	3	12

All (305) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	9	GLU
2	B	15	VAL
2	B	21	ARG
2	B	22	LYS
2	B	23	ARG
2	B	39	ILE
2	B	48	MET
2	B	53	ARG
2	B	61	LEU
2	B	69	LEU
2	B	81	VAL
2	B	111	ARG
2	B	117	GLU
2	B	127	ILE
2	B	129	GLU
2	B	134	GLU
2	B	139	LYS
2	B	143	GLU
2	B	147	LYS
2	B	170	GLU
2	B	172	ILE
2	B	178	ARG
2	B	190	THR
2	B	196	LEU
2	B	197	VAL
2	B	200	ILE
2	B	215	LEU
2	B	230	VAL
2	B	231	GLU
3	C	3	ASN
3	C	4	LYS
3	C	5	ILE
3	C	30	ARG
3	C	32	LEU
3	C	34	LEU
3	C	43	LEU

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Mol	Chain	Res	Type
3	C	62	ASP
3	C	67	THR
3	C	72	LYS
3	C	76	VAL
3	C	85	ARG
3	C	89	GLU
3	C	101	LEU
3	C	102	ASN
3	C	107	GLN
3	C	128	PHE
3	C	130	VAL
3	C	138	VAL
3	C	162	GLN
3	C	165	THR
3	C	166	GLU
3	C	167	TRP
3	C	173	VAL
3	C	179	ARG
3	C	188	LEU
3	C	191	THR
3	C	192	THR
3	C	196	LEU
3	C	204	LEU
4	D	3	ARG
4	D	8	VAL
4	D	9	CYS
4	D	15	GLU
4	D	45	GLN
4	D	57	ARG
4	D	64	LEU
4	D	70	ILE
4	D	78	LEU
4	D	85	LYS
4	D	89	THR
4	D	99	SER
4	D	119	GLN
4	D	122	ARG
4	D	127	THR
4	D	135	LEU
4	D	140	VAL
4	D	150	GLU
4	D	162	LEU

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Mol	Chain	Res	Type
4	D	170	VAL
4	D	175	SER
4	D	176	LEU
4	D	188	LEU
5	E	9	LYS
5	E	10	MET
5	E	12	LEU
5	E	15	ARG
5	E	16	THR
5	E	24	ARG
5	E	34	VAL
5	E	41	VAL
5	E	43	LEU
5	E	47	LYS
5	E	56	GLN
5	E	69	VAL
5	E	80	ILE
5	E	89	ILE
5	E	92	LYS
5	E	129	ILE
5	E	135	THR
5	E	137	GLU
5	E	144	THR
5	E	150	ARG
5	E	151	LEU
6	F	10	LEU
6	F	21	LEU
6	F	25	ILE
6	F	28	ARG
6	F	40	VAL
6	F	47	ARG
6	F	69	GLU
6	F	70	ASP
6	F	72	VAL
6	F	73	ASN
6	F	78	GLU
6	F	81	ILE
6	F	91	VAL
6	F	92	LYS
7	G	8	GLU
7	G	12	LEU
7	G	24	THR

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Mol	Chain	Res	Type
7	G	32	ARG
7	G	49	ILE
7	G	50	ILE
7	G	52	GLU
7	G	69	VAL
7	G	80	VAL
7	G	90	GLU
7	G	98	SER
7	G	113	GLU
7	G	115	ARG
7	G	124	LEU
7	G	138	LYS
7	G	140	ASP
7	G	141	VAL
7	G	149	ARG
7	G	155	ARG
8	H	12	ARG
8	H	24	THR
8	H	25	ASP
8	H	26	VAL
8	H	35	ILE
8	H	39	LEU
8	H	50	ARG
8	H	63	LEU
8	H	83	ILE
8	H	85	ARG
8	H	88	LYS
8	H	91	ARG
8	H	97	VAL
8	H	109	ILE
8	H	119	LEU
8	H	121	ASP
9	I	3	GLN
9	I	12	GLU
9	I	28	VAL
9	I	53	VAL
9	I	79	LEU
9	I	102	LEU
9	I	108	VAL
9	I	127	LYS
10	J	3	LYS
10	J	4	ILE

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Mol	Chain	Res	Type
10	J	6	ILE
10	J	19	SER
10	J	30	SER
10	J	35	SER
10	J	66	ARG
10	J	71	LEU
10	J	74	ILE
10	J	75	ILE
10	J	80	LYS
10	J	84	GLN
10	J	86	MET
10	J	87	THR
10	J	90	LEU
10	J	92	THR
10	J	98	ILE
11	K	29	ILE
11	K	47	VAL
11	K	48	ILE
11	K	54	ARG
11	K	79	SER
11	K	91	ARG
11	K	99	GLN
11	K	106	LYS
11	K	114	VAL
11	K	127	LYS
12	L	7	ILE
12	L	13	LYS
12	L	28	LYS
12	L	33	ARG
12	L	37	CYS
12	L	39	VAL
12	L	40	VAL
12	L	43	VAL
12	L	54	LYS
12	L	55	VAL
12	L	60	LEU
12	L	78	GLN
12	L	79	GLU
12	L	81	SER
12	L	83	VAL
12	L	93	LEU
13	M	4	ILE

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Mol	Chain	Res	Type
13	M	9	ILE
13	M	12	ASN
13	M	13	LYS
13	M	16	ASP
13	M	34	LEU
13	M	37	THR
13	M	46	LYS
13	M	47	ASP
13	M	88	ARG
13	M	105	THR
13	M	106	ASN
13	M	110	ARG
13	M	115	LYS
13	M	125	ARG
14	N	9	LYS
14	N	12	ARG
14	N	22	THR
14	N	32	SER
14	N	33	VAL
14	N	41	ARG
14	N	58	LYS
15	O	4	THR
15	O	5	LYS
15	O	6	GLU
15	O	7	GLU
15	O	10	LYS
15	O	13	GLN
15	O	22	THR
15	O	31	LEU
15	O	34	LEU
15	O	39	LEU
15	O	41	GLU
15	O	57	LEU
15	O	65	ARG
15	O	70	LEU
15	O	71	GLN
15	O	81	LEU
15	O	82	ILE
15	O	88	ARG
16	P	1	MET
16	P	2	VAL
16	P	20	VAL

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Mol	Chain	Res	Type
16	P	31	LYS
16	P	45	THR
16	P	62	VAL
16	P	76	GLN
16	P	80	PHE
16	P	81	ARG
16	P	82	GLN
17	Q	10	VAL
17	Q	12	SER
17	Q	34	LYS
17	Q	35	VAL
17	Q	38	ARG
17	Q	52	LYS
17	Q	55	ASP
17	Q	59	ILE
17	Q	60	ILE
17	Q	74	LEU
17	Q	78	GLU
17	Q	100	LYS
17	Q	101	ARG
18	R	21	LYS
18	R	47	THR
18	R	54	ARG
18	R	69	THR
18	R	86	VAL
18	R	88	LYS
19	S	6	LYS
19	S	7	LYS
19	S	9	VAL
19	S	12	ASP
19	S	18	LYS
19	S	20	LEU
19	S	25	LYS
19	S	27	GLU
19	S	39	THR
19	S	43	GLU
19	S	62	ILE
19	S	63	THR
19	S	65	ASN
19	S	77	THR
19	S	79	THR
19	S	81	ARG

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Mol	Chain	Res	Type
20	T	19	SER
20	T	21	LYS
20	T	27	LYS
20	T	30	LYS
20	T	38	LYS
20	T	53	LEU
20	T	56	MET
20	T	64	ASP
20	T	68	LYS
20	T	75	ASN
20	T	80	ARG
20	T	84	LEU
20	T	86	ARG
21	V	6	ARG
21	V	8	THR
21	V	12	LYS
21	V	15	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	16	HIS
3	C	6	HIS
3	C	118	GLN
4	D	42	GLN
4	D	74	GLN
4	D	119	GLN
4	D	161	ASN
6	F	13	ASN
6	F	27	GLN
7	G	84	ASN
7	G	153	HIS
11	K	13	GLN
13	M	106	ASN
16	P	82	GLN
19	S	47	HIS
19	S	65	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	1506/1507 (99%)	300 (19%)	46 (3%)
22	X	8/17 (47%)	1 (12%)	0
23	W	3/6 (50%)	0	0
All	All	1517/1530 (99%)	301 (19%)	46 (3%)

All (301) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	7	G
1	A	9	G
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	54	C
1	A	59	A
1	A	60	A
1	A	61	G
1	A	88	A
1	A	101	A
1	A	115	G
1	A	116	A
1	A	121	C
1	A	129(A)	G
1	A	130	A
1	A	131	C
1	A	163	C
1	A	181	G
1	A	182	U
1	A	188	C
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(L)	U
1	A	195	A
1	A	197	A
1	A	202	U

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Mol	Chain	Res	Type
1	A	203	U
1	A	204	U
1	A	216	G
1	A	220	G
1	A	243	A
1	A	244	U
1	A	247	G
1	A	251	G
1	A	262	A
1	A	266	G
1	A	267	C
1	A	279	A
1	A	289	G
1	A	299	G
1	A	301	G
1	A	319	G
1	A	321	A
1	A	328	C
1	A	329	A
1	A	332	G
1	A	345	C
1	A	347	G
1	A	350	G
1	A	352	C
1	A	353	A
1	A	354	G
1	A	367	U
1	A	373	A
1	A	384	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

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Mol	Chain	Res	Type
1	A	417	C
1	A	420	U
1	A	421	U
1	A	422	C
1	A	423	G
1	A	424	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	436	C
1	A	439	A
1	A	440	A
1	A	448	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
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	518	C
1	A	521	G
1	A	527	G
1	A	531	U
1	A	532	A
1	A	533	A
1	A	547	A
1	A	559	A
1	A	560	U
1	A	561	U
1	A	562	C

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Mol	Chain	Res	Type
1	A	564	C
1	A	572	A
1	A	573	A
1	A	576	G
1	A	577	G
1	A	579	G
1	A	582	U
1	A	588	G
1	A	596	C
1	A	607	A
1	A	634	C
1	A	653	A
1	A	665	A
1	A	671	G
1	A	687	A
1	A	694	A
1	A	695	A
1	A	702	A
1	A	703	G
1	A	721	G
1	A	723	U
1	A	731	G
1	A	734	G
1	A	748	C
1	A	749	C
1	A	755	G
1	A	759	A
1	A	774	G
1	A	777	A
1	A	785	G
1	A	787	A
1	A	793	U
1	A	794	A
1	A	795	C
1	A	812	C
1	A	813	U
1	A	817	C
1	A	828	A
1	A	839	U
1	A	840	C
1	A	841	U
1	A	848	C

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Mol	Chain	Res	Type
1	A	859	A
1	A	873	A
1	A	874	G
1	A	876	G
1	A	902	G
1	A	914	A
1	A	916	G
1	A	919	A
1	A	926	G
1	A	927	G
1	A	928	G
1	A	932	C
1	A	934	C
1	A	935	A
1	A	942	G
1	A	960	U
1	A	961	U
1	A	965	A
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	984	C
1	A	991	U
1	A	992	U
1	A	993	G
1	A	1002	G
1	A	1005	A
1	A	1007	C
1	A	1011	G
1	A	1024	G
1	A	1025	U
1	A	1026	G
1	A	1029	C
1	A	1030	C
1	A	1032	G
1	A	1033	G
1	A	1049	U

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Mol	Chain	Res	Type
1	A	1050	G
1	A	1053	G
1	A	1054	C
1	A	1055	A
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	1138	G
1	A	1139	G
1	A	1146	A
1	A	1152	A
1	A	1154	G
1	A	1157	A
1	A	1158	C
1	A	1159	U
1	A	1171	G
1	A	1183	A
1	A	1184	G
1	A	1190	G
1	A	1191	A
1	A	1196	U
1	A	1197	G
1	A	1198	G
1	A	1200	C
1	A	1201	A
1	A	1202	G
1	A	1211	U
1	A	1212	U

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Mol	Chain	Res	Type
1	A	1213	A
1	A	1224	G
1	A	1225	A
1	A	1226	C
1	A	1227	A
1	A	1238	A
1	A	1257	U
1	A	1258	G
1	A	1266	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	1290	G
1	A	1301	U
1	A	1302	U
1	A	1305	G
1	A	1306	A
1	A	1312	G
1	A	1320	C
1	A	1322	C
1	A	1338	G
1	A	1340	A
1	A	1346	A
1	A	1347	G
1	A	1353	G
1	A	1360	A
1	A	1361	G
1	A	1362	C
1	A	1363	A
1	A	1368	G
1	A	1370	G
1	A	1398	A
1	A	1446	A
1	A	1447	G
1	A	1448	C
1	A	1454	G
1	A	1487	G
1	A	1492	A
1	A	1498	U

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Mol	Chain	Res	Type
1	A	1499	A
1	A	1502	A
1	A	1503	A
1	A	1504	G
1	A	1505	G
1	A	1506	U
1	A	1517	G
1	A	1520	G
1	A	1529	G
1	A	1530	G
1	A	1533	C
1	A	1534	A
1	A	1539	C
22	X	32	RSP

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	250	A
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	532	A
1	A	559	A
1	A	701	C
1	A	812	C
1	A	840	C
1	A	960	U
1	A	965	A
1	A	975	A
1	A	1049	U

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Mol	Chain	Res	Type
1	A	1054	C
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	1190	G
1	A	1201	A
1	A	1212	U
1	A	1224	G
1	A	1281	U
1	A	1285	A
1	A	1300	G
1	A	1305	G
1	A	1346	A
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 ⓘ

1 non-standard protein/DNA/RNA residue is modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
22	RSP	X	32	22	17,21,22	0.77	0	21,30,33	0.65	1 (4%)

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
22	RSP	X	32	22	-	2/7/25/26	0/2/2/2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
22	X	32	RSP	S2-C2-N3	-2.01	118.01	121.27

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
22	X	32	RSP	O4'-C4'-C5'-O5'
22	X	32	RSP	C3'-C4'-C5'-O5'

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 98 ligands modelled in this entry, 97 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
25	PAR	A	1693	-	44,45,45	1.38	9 (20%)	63,67,67	1.65	12 (19%)

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
25	PAR	A	1693	-	-	3/18/94/94	0/4/4/4

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
25	A	1693	PAR	C52-C42	3.42	1.59	1.52
25	A	1693	PAR	C13-C23	2.74	1.56	1.52
25	A	1693	PAR	O54-C14	2.65	1.48	1.41
25	A	1693	PAR	O43-C13	2.33	1.45	1.41
25	A	1693	PAR	C22-C12	-2.30	1.48	1.53
25	A	1693	PAR	O54-C54	2.23	1.49	1.44
25	A	1693	PAR	C44-C54	2.18	1.57	1.53
25	A	1693	PAR	O33-C14	2.17	1.47	1.41
25	A	1693	PAR	C62-C52	2.02	1.57	1.52

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
25	A	1693	PAR	O33-C14-C24	6.61	118.90	108.08
25	A	1693	PAR	O54-C54-C44	3.13	115.33	109.70
25	A	1693	PAR	C34-C24-N24	-2.98	104.95	111.05
25	A	1693	PAR	C14-O33-C33	-2.94	111.01	117.98
25	A	1693	PAR	O11-C11-O51	2.89	118.29	110.69
25	A	1693	PAR	O34-C34-C44	-2.82	103.74	110.38
25	A	1693	PAR	O52-C13-O43	-2.58	108.73	111.37
25	A	1693	PAR	O51-C51-C61	2.36	112.29	106.44
25	A	1693	PAR	C11-O51-C51	2.26	118.14	113.72
25	A	1693	PAR	C13-O52-C52	-2.24	112.68	117.98
25	A	1693	PAR	O43-C13-C23	-2.21	102.17	104.98
25	A	1693	PAR	O11-C11-C21	-2.04	104.74	108.08

There are no chirality outliers.

All (3) torsion outliers are listed below:

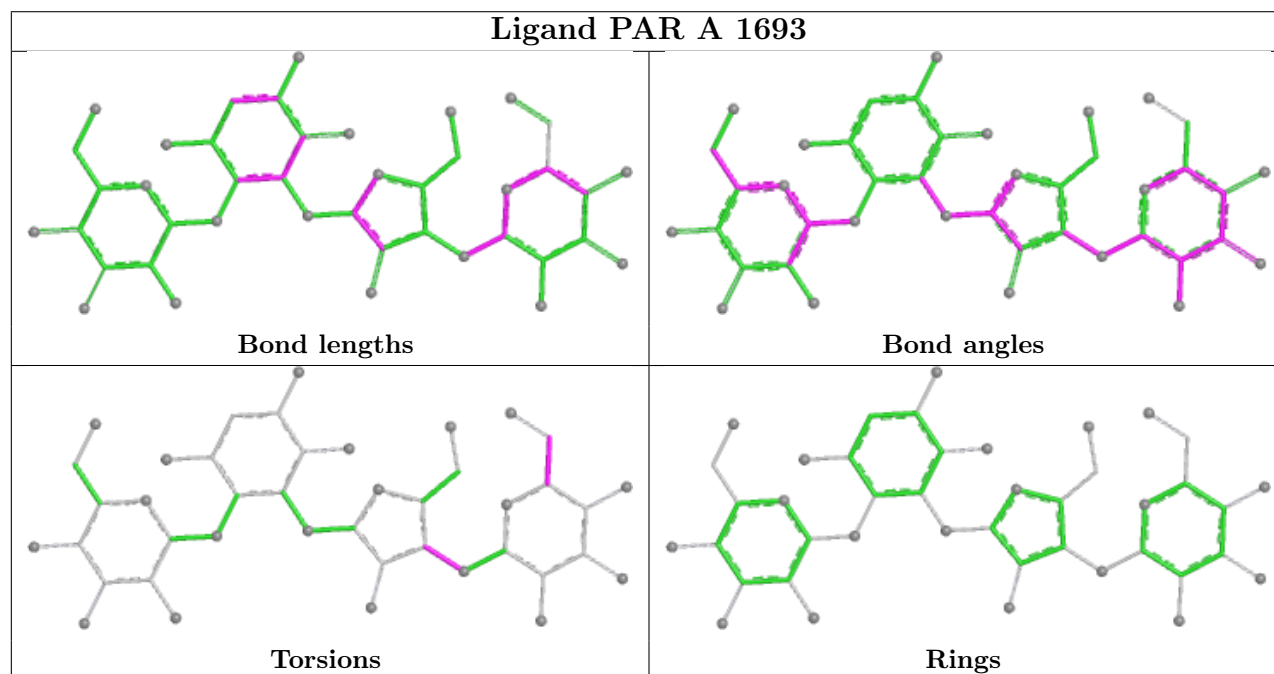
Mol	Chain	Res	Type	Atoms
25	A	1693	PAR	C44-C54-C64-N64
25	A	1693	PAR	O54-C54-C64-N64
25	A	1693	PAR	C23-C33-O33-C14

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
25	A	1693	PAR	1	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%)	-0.31	31 (2%) 63 45	69, 107, 204, 419	0
2	B	234/256 (91%)	0.03	4 (1%) 69 51	91, 147, 211, 278	0
3	C	206/239 (86%)	0.24	8 (3%) 43 29	87, 138, 201, 253	0
4	D	208/209 (99%)	0.15	9 (4%) 40 26	78, 120, 175, 210	0
5	E	150/162 (92%)	-0.09	2 (1%) 75 57	71, 102, 147, 189	0
6	F	101/101 (100%)	-0.03	2 (1%) 65 47	98, 137, 178, 233	0
7	G	155/156 (99%)	0.06	4 (2%) 57 39	89, 132, 192, 229	0
8	H	138/138 (100%)	-0.26	2 (1%) 73 56	55, 96, 139, 190	0
9	I	127/128 (99%)	0.30	6 (4%) 36 25	78, 142, 183, 208	0
10	J	98/105 (93%)	0.81	14 (14%) 6 6	89, 170, 259, 294	0
11	K	119/129 (92%)	0.03	6 (5%) 34 23	69, 110, 169, 283	0
12	L	124/132 (93%)	0.39	7 (5%) 30 21	61, 110, 166, 216	0
13	M	125/126 (99%)	0.27	7 (5%) 30 21	98, 139, 217, 295	0
14	N	60/61 (98%)	0.38	5 (8%) 17 13	94, 125, 176, 244	0
15	O	88/89 (98%)	0.23	5 (5%) 29 20	72, 115, 177, 236	0
16	P	83/88 (94%)	0.17	3 (3%) 46 31	71, 99, 121, 215	0
17	Q	104/105 (99%)	0.36	4 (3%) 44 30	71, 107, 185, 241	0
18	R	73/88 (82%)	-0.06	3 (4%) 41 27	89, 119, 198, 283	0
19	S	80/93 (86%)	0.42	6 (7%) 20 15	103, 161, 199, 248	0
20	T	99/106 (93%)	0.28	5 (5%) 33 23	80, 106, 172, 200	0
21	V	24/27 (88%)	0.74	3 (12%) 8 7	98, 129, 165, 197	0
22	X	9/17 (52%)	0.33	2 (22%) 2 2	121, 149, 208, 277	0
23	W	4/6 (66%)	1.00	1 (25%) 2 2	112, 114, 131, 149	0
All	All	3916/4068 (96%)	-0.01	139 (3%) 47 31	55, 118, 201, 419	0

All (139) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
11	K	129	SER	10.9
11	K	128	ALA	8.2
13	M	126	LYS	7.5
7	G	33	ASP	7.1
19	S	3	ARG	6.4
11	K	127	LYS	5.9
3	C	161	GLU	5.8
12	L	128	ALA	4.8
20	T	9	ASN	4.8
12	L	47	LYS	4.7
20	T	68	LYS	4.6
1	A	1533	C	4.6
4	D	2	GLY	4.6
1	A	407	G	4.5
4	D	77	ASN	4.4
6	F	94	GLN	4.3
12	L	19	ARG	4.3
17	Q	105	ALA	4.3
14	N	13	THR	4.2
1	A	417	C	4.1
1	A	422	C	4.1
1	A	462	G	4.1
21	V	25	LYS	4.0
1	A	416	G	4.0
13	M	8	GLU	4.0
10	J	45	ARG	3.9
15	O	48	LYS	3.9
14	N	4	LYS	3.9
19	S	2	PRO	3.9
1	A	424	G	3.8
10	J	73	ASP	3.8
1	A	1541	U	3.8
19	S	37	ARG	3.8
23	W	4	A	3.8
1	A	1539	C	3.7
1	A	410	G	3.7
13	M	124	PRO	3.7
14	N	30	ALA	3.7
10	J	4	ILE	3.7
9	I	128	ARG	3.6
1	A	1361	G	3.6
10	J	43	ARG	3.6

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Mol	Chain	Res	Type	RSRZ
18	R	31	LEU	3.6
1	A	217	C	3.6
1	A	1031	G	3.5
3	C	178	LEU	3.5
1	A	421	U	3.5
10	J	34	VAL	3.5
12	L	18	VAL	3.4
21	V	6	ARG	3.2
7	G	8	GLU	3.2
1	A	420	U	3.1
19	S	73	GLU	3.1
1	A	408	A	3.1
13	M	123	ALA	3.1
7	G	106	GLN	3.0
8	H	1	MET	3.0
1	A	1126	U	3.0
1	A	1540	U	3.0
10	J	58	ASP	3.0
3	C	42	LEU	3.0
6	F	97	PHE	3.0
13	M	125	ARG	2.9
15	O	2	PRO	2.9
9	I	30	GLY	2.9
8	H	18	ARG	2.9
4	D	31	CYS	2.9
2	B	191	ASP	2.8
17	Q	13	ASP	2.8
21	V	2	GLY	2.8
16	P	29	ASP	2.8
20	T	64	ASP	2.8
9	I	126	SER	2.8
3	C	177	THR	2.8
1	A	413	G	2.7
18	R	32	ARG	2.7
13	M	35	GLU	2.7
22	X	30	G	2.7
10	J	72	VAL	2.7
1	A	202	U	2.7
3	C	44	GLU	2.6
13	M	6	GLY	2.6
20	T	69	GLY	2.6
1	A	1534	A	2.6

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Mol	Chain	Res	Type	RSRZ
10	J	5	ARG	2.6
19	S	81	ARG	2.6
9	I	106	ALA	2.6
4	D	3	ARG	2.6
12	L	89	ARG	2.6
5	E	27	ARG	2.5
10	J	37	PRO	2.5
11	K	126	ARG	2.5
17	Q	100	LYS	2.5
10	J	18	ALA	2.4
15	O	58	MET	2.4
5	E	79	GLU	2.4
10	J	100	THR	2.4
2	B	16	HIS	2.4
3	C	176	HIS	2.4
4	D	22	LYS	2.4
1	A	1362	C	2.4
3	C	167	TRP	2.4
22	X	33	U	2.4
4	D	81	GLU	2.3
17	Q	58	GLU	2.3
10	J	71	LEU	2.3
19	S	5	LEU	2.3
14	N	20	ALA	2.3
12	L	17	LYS	2.3
1	A	1280	A	2.3
18	R	30	ASP	2.3
11	K	51	LYS	2.2
7	G	5	ARG	2.2
1	A	251	G	2.2
1	A	423	G	2.2
1	A	425	G	2.2
4	D	72	GLU	2.2
1	A	532	A	2.2
15	O	59	MET	2.2
11	K	52	GLY	2.2
1	A	474	G	2.2
1	A	477	G	2.2
2	B	133	LYS	2.1
3	C	129	ALA	2.1
4	D	24	GLU	2.1
14	N	2	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	1051	C	2.1
20	T	73	HIS	2.1
9	I	7	THR	2.1
15	O	54	ARG	2.1
2	B	132	LYS	2.1
10	J	66	ARG	2.1
1	A	460	A	2.0
12	L	26	ALA	2.0
16	P	65	GLN	2.0
10	J	6	ILE	2.0
16	P	80	PHE	2.0
9	I	66	ARG	2.0
4	D	5	ILE	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
22	RSP	X	32	20/21	0.70	0.15	238,239,242,242	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
24	MG	A	1607	1/1	0.57	0.34	106,106,106,106	0
24	MG	A	1675	1/1	0.63	0.36	98,98,98,98	0
24	MG	A	1653	1/1	0.64	0.45	62,62,62,62	0
24	MG	A	1656	1/1	0.70	0.34	92,92,92,92	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
24	MG	A	1640	1/1	0.73	0.32	93,93,93,93	0
24	MG	A	1615	1/1	0.75	0.23	94,94,94,94	0
24	MG	A	1680	1/1	0.75	0.47	110,110,110,110	0
24	MG	A	1623	1/1	0.77	0.36	56,56,56,56	0
24	MG	A	1684	1/1	0.77	0.32	94,94,94,94	0
24	MG	N	101	1/1	0.77	0.27	94,94,94,94	0
24	MG	A	1678	1/1	0.78	0.30	77,77,77,77	0
24	MG	A	1634	1/1	0.78	0.46	96,96,96,96	0
24	MG	W	101	1/1	0.78	0.22	105,105,105,105	0
24	MG	A	1659	1/1	0.80	0.25	108,108,108,108	0
24	MG	A	1667	1/1	0.80	0.41	109,109,109,109	0
24	MG	A	1671	1/1	0.80	0.33	121,121,121,121	0
24	MG	A	1604	1/1	0.81	0.17	97,97,97,97	0
24	MG	A	1682	1/1	0.82	0.31	60,60,60,60	0
24	MG	A	1603	1/1	0.84	0.33	102,102,102,102	0
24	MG	A	1629	1/1	0.84	0.28	69,69,69,69	0
24	MG	A	1650	1/1	0.85	0.33	113,113,113,113	0
24	MG	A	1639	1/1	0.86	0.25	61,61,61,61	0
24	MG	A	1669	1/1	0.86	0.40	98,98,98,98	0
24	MG	A	1679	1/1	0.86	0.32	99,99,99,99	0
24	MG	A	1663	1/1	0.86	0.28	96,96,96,96	0
24	MG	A	1676	1/1	0.87	0.23	87,87,87,87	0
24	MG	A	1601	1/1	0.87	0.21	100,100,100,100	0
24	MG	A	1686	1/1	0.87	0.40	87,87,87,87	0
24	MG	A	1687	1/1	0.87	0.33	99,99,99,99	0
24	MG	A	1689	1/1	0.87	0.16	123,123,123,123	0
24	MG	A	1631	1/1	0.87	0.67	85,85,85,85	0
24	MG	A	1633	1/1	0.87	0.20	70,70,70,70	0
24	MG	A	1605	1/1	0.88	0.40	81,81,81,81	0
24	MG	A	1627	1/1	0.88	0.35	57,57,57,57	0
24	MG	A	1621	1/1	0.88	0.29	78,78,78,78	0
24	MG	A	1660	1/1	0.88	0.29	107,107,107,107	0
24	MG	A	1649	1/1	0.89	0.40	70,70,70,70	0
24	MG	A	1664	1/1	0.89	0.17	88,88,88,88	0
24	MG	A	1665	1/1	0.89	0.20	101,101,101,101	0
24	MG	A	1655	1/1	0.90	0.18	90,90,90,90	0
24	MG	A	1616	1/1	0.90	0.31	54,54,54,54	0
24	MG	A	1651	1/1	0.91	0.36	78,78,78,78	0
24	MG	A	1611	1/1	0.91	0.23	76,76,76,76	0
24	MG	A	1690	1/1	0.91	0.43	97,97,97,97	0
24	MG	A	1637	1/1	0.91	0.32	83,83,83,83	0
24	MG	A	1625	1/1	0.91	0.34	62,62,62,62	0

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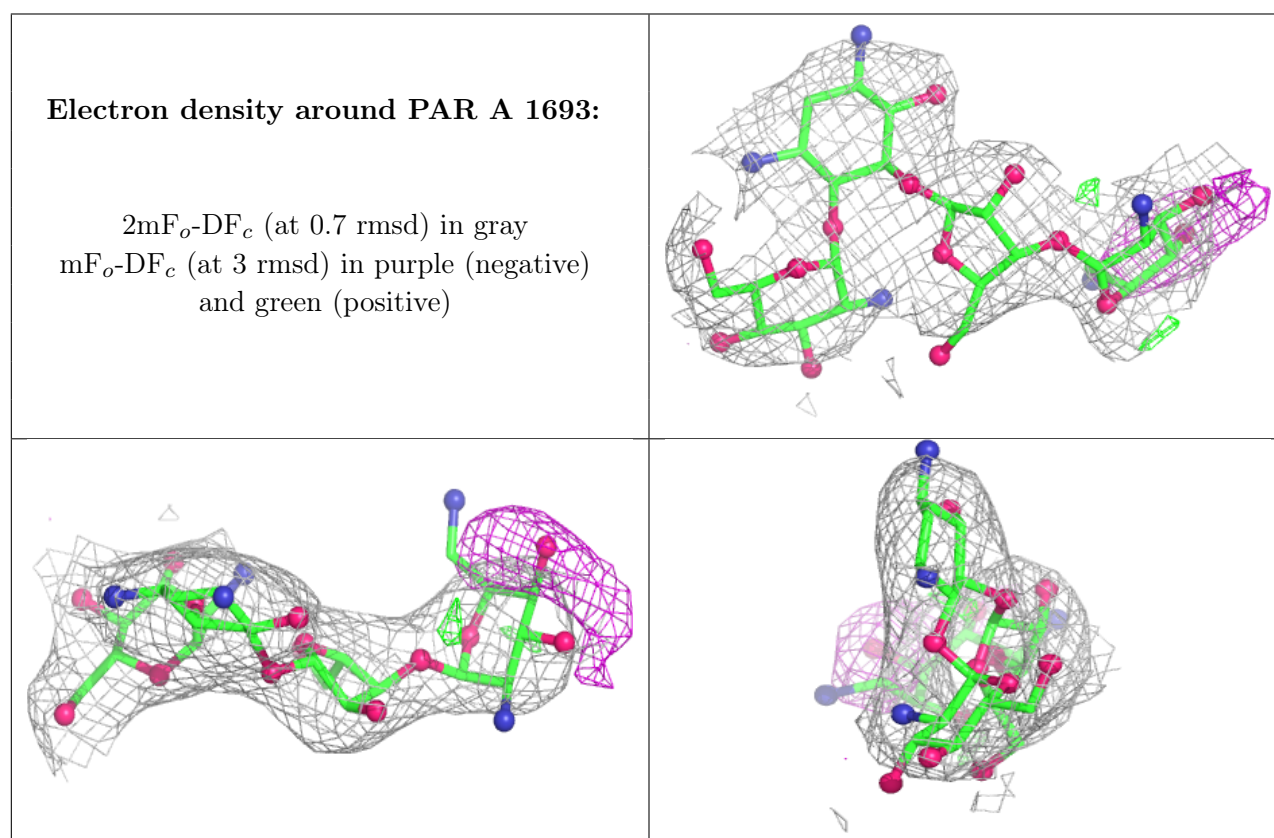
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
24	MG	A	1635	1/1	0.92	0.26	76,76,76,76	0
24	MG	A	1681	1/1	0.92	0.32	106,106,106,106	0
24	MG	A	1610	1/1	0.92	0.22	108,108,108,108	0
24	MG	A	1602	1/1	0.92	0.17	83,83,83,83	0
24	MG	A	1658	1/1	0.92	0.49	86,86,86,86	0
24	MG	A	1674	1/1	0.92	0.27	87,87,87,87	0
24	MG	A	1617	1/1	0.92	0.42	66,66,66,66	0
24	MG	A	1619	1/1	0.92	0.42	83,83,83,83	0
24	MG	A	1612	1/1	0.92	0.32	83,83,83,83	0
24	MG	A	1613	1/1	0.92	0.11	106,106,106,106	0
25	PAR	A	1693	42/42	0.92	0.14	90,95,99,100	0
24	MG	A	1606	1/1	0.93	0.44	106,106,106,106	0
24	MG	A	1609	1/1	0.93	0.55	96,96,96,96	0
24	MG	A	1677	1/1	0.93	0.22	92,92,92,92	0
24	MG	A	1694	1/1	0.93	0.21	83,83,83,83	0
24	MG	A	1672	1/1	0.93	0.26	84,84,84,84	0
24	MG	A	1685	1/1	0.93	0.16	84,84,84,84	0
24	MG	A	1638	1/1	0.93	0.30	62,62,62,62	0
24	MG	A	1657	1/1	0.94	0.15	65,65,65,65	0
24	MG	A	1626	1/1	0.94	0.08	103,103,103,103	0
24	MG	A	1641	1/1	0.94	0.23	77,77,77,77	0
24	MG	A	1644	1/1	0.94	0.13	93,93,93,93	0
24	MG	A	1662	1/1	0.94	0.27	74,74,74,74	0
24	MG	A	1688	1/1	0.94	0.33	83,83,83,83	0
24	MG	A	1668	1/1	0.95	0.32	111,111,111,111	0
24	MG	A	1630	1/1	0.95	0.22	85,85,85,85	0
24	MG	A	1683	1/1	0.95	0.32	76,76,76,76	0
24	MG	A	1628	1/1	0.95	0.39	68,68,68,68	0
24	MG	J	201	1/1	0.95	0.10	82,82,82,82	0
24	MG	A	1666	1/1	0.95	0.48	78,78,78,78	0
24	MG	A	1673	1/1	0.95	0.13	57,57,57,57	0
24	MG	A	1652	1/1	0.95	0.16	65,65,65,65	0
24	MG	A	1648	1/1	0.96	0.09	66,66,66,66	0
24	MG	A	1614	1/1	0.96	0.35	80,80,80,80	0
24	MG	A	1670	1/1	0.96	0.48	99,99,99,99	0
24	MG	A	1620	1/1	0.96	0.24	58,58,58,58	0
24	MG	A	1642	1/1	0.96	0.18	91,91,91,91	0
24	MG	A	1636	1/1	0.96	0.19	76,76,76,76	0
24	MG	A	1646	1/1	0.96	0.43	72,72,72,72	0
24	MG	A	1691	1/1	0.97	0.45	98,98,98,98	0
24	MG	A	1618	1/1	0.97	0.30	68,68,68,68	0
24	MG	A	1632	1/1	0.97	0.30	66,66,66,66	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
24	MG	A	1624	1/1	0.97	0.37	87,87,87,87	0
24	MG	A	1643	1/1	0.97	0.12	80,80,80,80	0
24	MG	A	1654	1/1	0.97	0.41	66,66,66,66	0
26	ZN	D	301	1/1	0.97	0.17	118,118,118,118	0
24	MG	A	1622	1/1	0.98	0.16	69,69,69,69	0
24	MG	A	1692	1/1	0.98	0.09	98,98,98,98	0
24	MG	A	1608	1/1	0.98	0.05	85,85,85,85	0
24	MG	A	1647	1/1	0.98	0.20	67,67,67,67	0
24	MG	A	1661	1/1	0.99	0.03	58,58,58,58	0
24	MG	A	1645	1/1	0.99	0.39	71,71,71,71	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.