



Full wwPDB EM Validation Report ⓘ

Mar 19, 2026 – 07:57 PM UTC

PDB ID : 9IOB / pdb_00009iob
EMDB ID : EMD-60725
Title : Cryo-EM structure of the hexameric DRT9-ncRNA complex
Authors : Zhang, J.T.; Song, X.Y.; Xia, Y.S.; Liu, Y.J.; Jia, N.
Deposited on : 2024-07-08
Resolution : 2.62 Å(reported)

This is a Full wwPDB EM 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/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : **NOT EXECUTED**
MapQ : **FAILED**
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

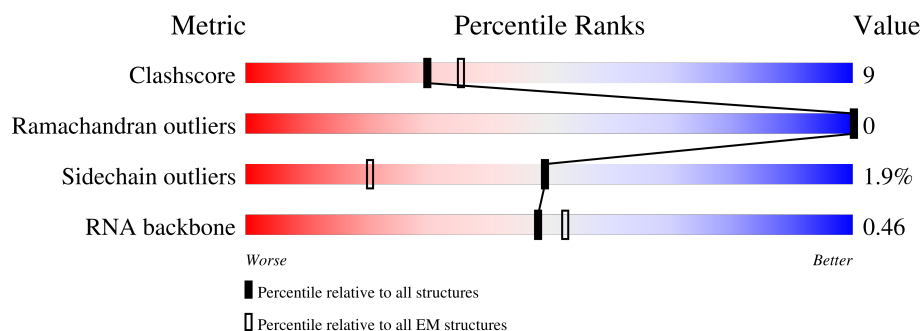
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.62 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)	EM structures (#Entries)
Clashscore	229148	23984
Ramachandran outliers	224038	23583
Sidechain outliers	223484	23102
RNA backbone	8273	3508

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	499	
1	C	499	
1	E	499	
1	G	499	
1	I	499	
1	K	499	
2	B	177	

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Mol	Chain	Length	Quality of chain
2	D	177	 51% 34% 15%
2	F	177	 51% 37% 12%
2	H	177	 51% 34% 14%
2	J	177	 53% 33% 14%
2	L	177	 50% 36% 14%

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 82602 atoms, of which 35946 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called RNA-dependent DNA polymerase.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	488	Total	C	H	N	O	S	0	0
			8124	2590	4103	702	720	9		
1	C	488	Total	C	H	N	O	S	0	0
			8124	2590	4103	702	720	9		
1	E	488	Total	C	H	N	O	S	0	0
			8124	2590	4103	702	720	9		
1	G	488	Total	C	H	N	O	S	0	0
			8124	2590	4103	702	720	9		
1	I	488	Total	C	H	N	O	S	0	0
			8124	2590	4103	702	720	9		
1	K	488	Total	C	H	N	O	S	0	0
			8124	2590	4103	702	720	9		

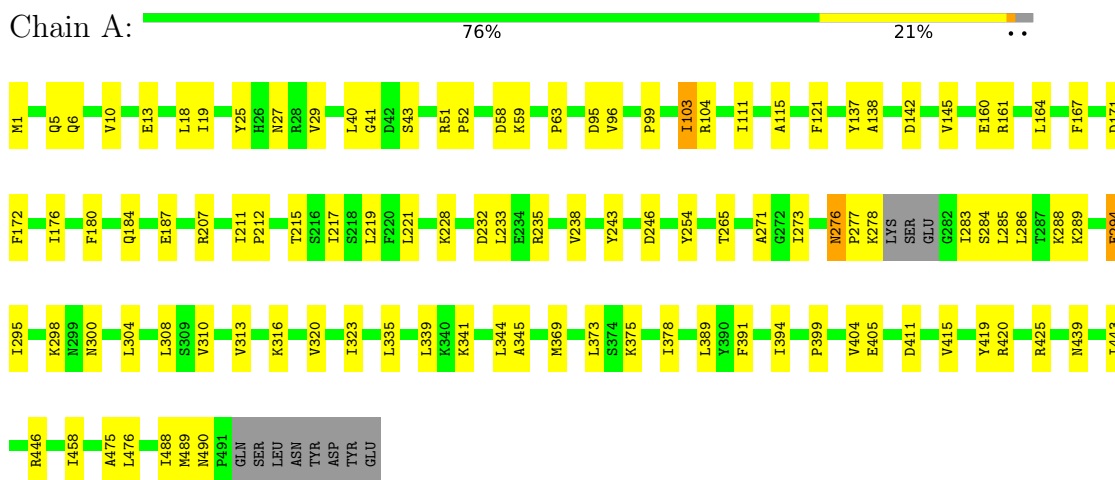
- Molecule 2 is a RNA chain called RNA (177-MER).

Mol	Chain	Residues	Atoms						AltConf	Trace
2	B	177	Total	C	H	N	O	P	0	0
			5643	1678	1888	641	1259	177		
2	D	177	Total	C	H	N	O	P	0	0
			5643	1678	1888	641	1259	177		
2	F	177	Total	C	H	N	O	P	0	0
			5643	1678	1888	641	1259	177		
2	H	177	Total	C	H	N	O	P	0	0
			5643	1678	1888	641	1259	177		
2	J	177	Total	C	H	N	O	P	0	0
			5643	1678	1888	641	1259	177		
2	L	177	Total	C	H	N	O	P	0	0
			5643	1678	1888	641	1259	177		

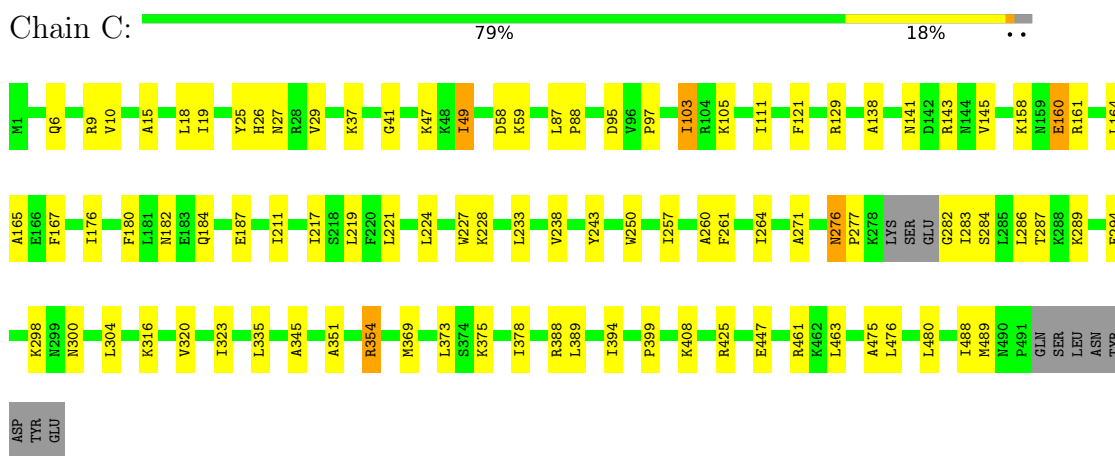
3 Residue-property plots

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. 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. 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.

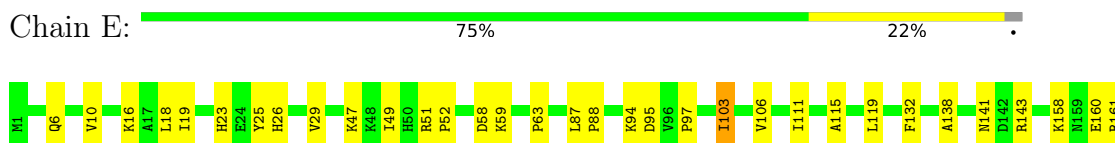
- Molecule 1: RNA-dependent DNA polymerase

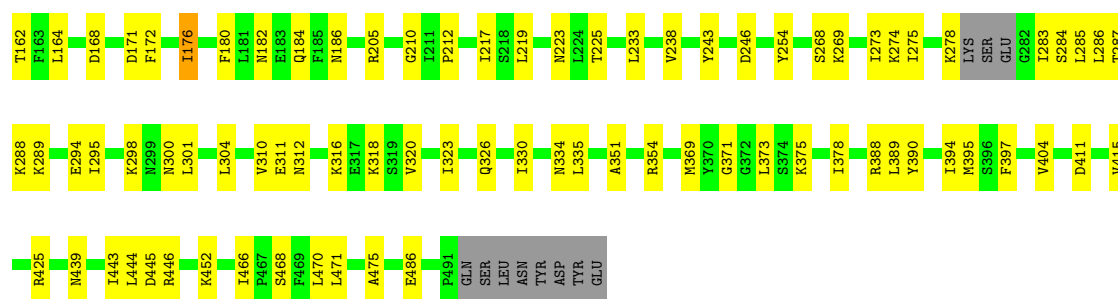


- Molecule 1: RNA-dependent DNA polymerase

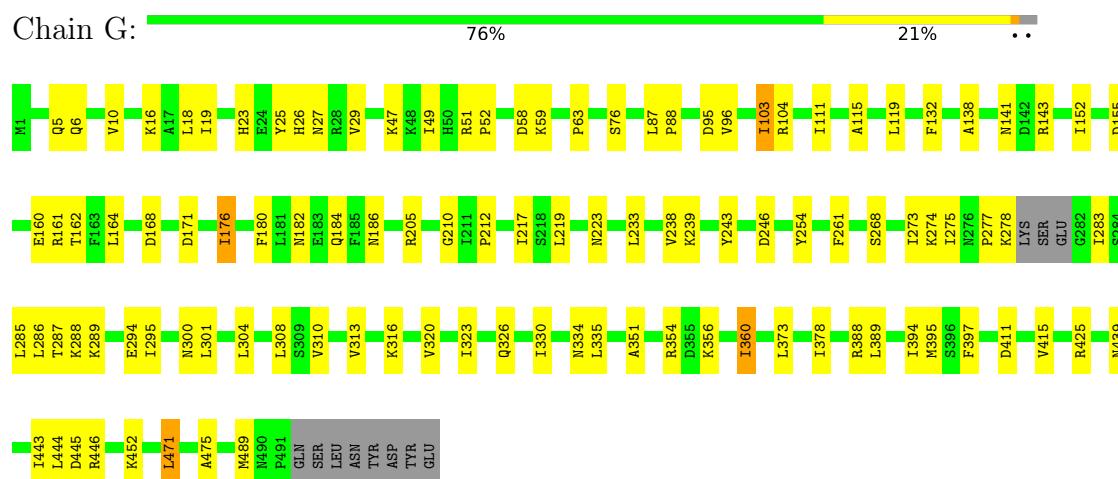


- Molecule 1: RNA-dependent DNA polymerase

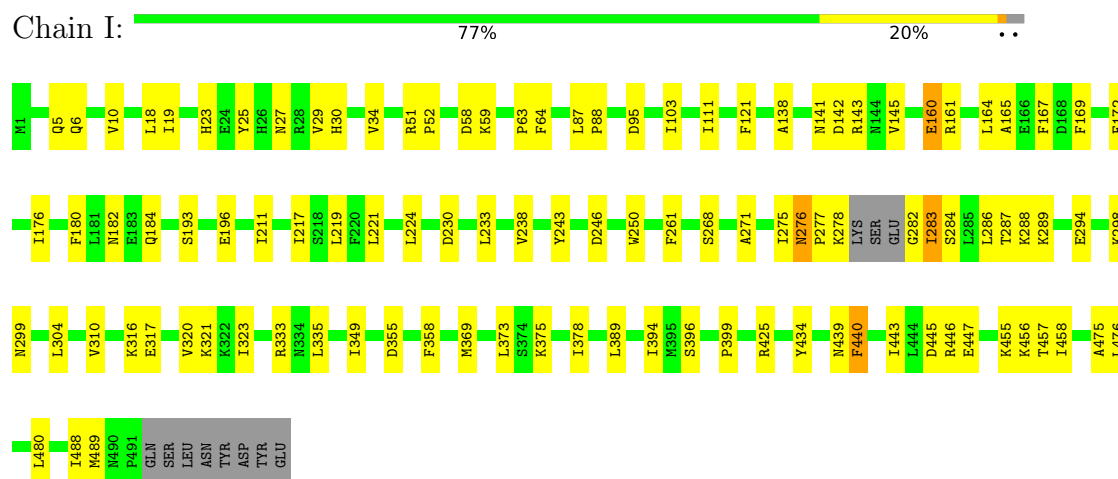




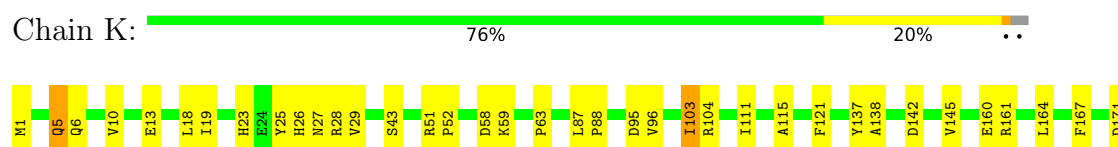
• Molecule 1: RNA-dependent DNA polymerase

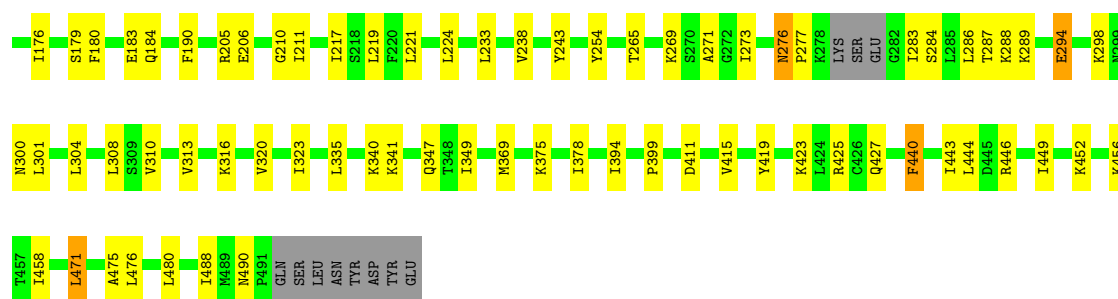


• Molecule 1: RNA-dependent DNA polymerase

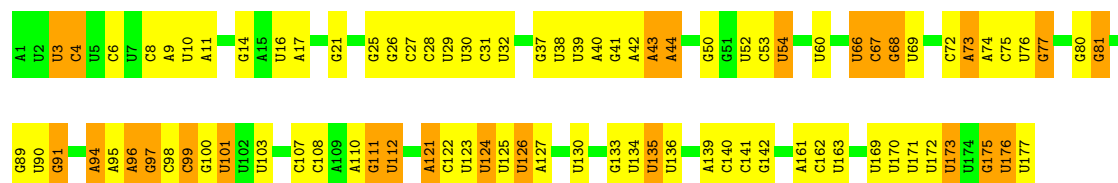


• Molecule 1: RNA-dependent DNA polymerase

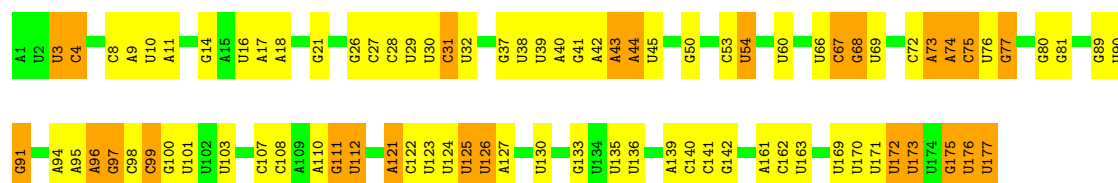




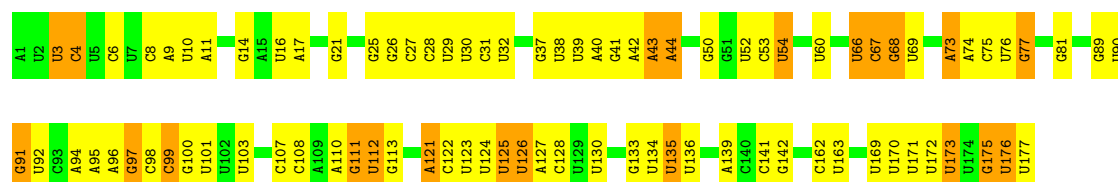
• Molecule 2: RNA (177-MER)



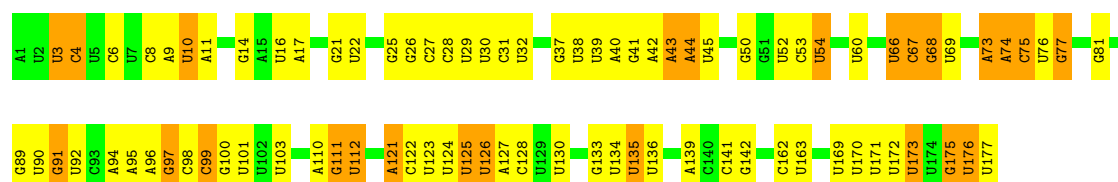
• Molecule 2: RNA (177-MER)



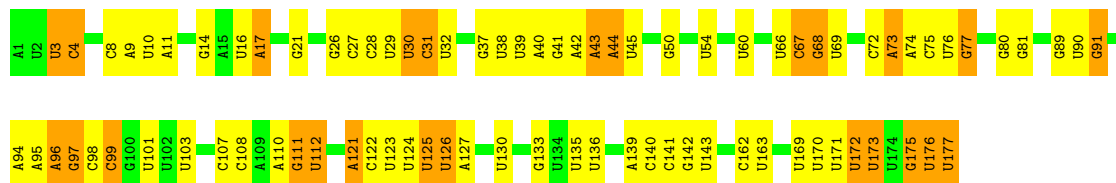
• Molecule 2: RNA (177-MER)



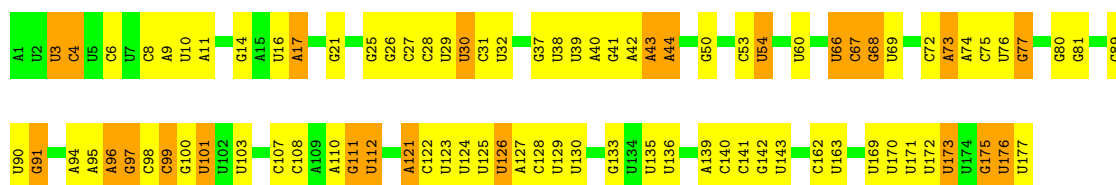
• Molecule 2: RNA (177-MER)



● Molecule 2: RNA (177-MER)

Chain J:  53% 33% 14%

● Molecule 2: RNA (177-MER)

Chain L:  50% 36% 14%

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	220612	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.22	0/4109	0.41	0/5521
1	C	0.22	0/4109	0.41	0/5521
1	E	0.22	0/4109	0.42	0/5521
1	G	0.22	0/4109	0.41	0/5521
1	I	0.22	0/4109	0.41	0/5521
1	K	0.21	0/4109	0.41	0/5521
2	B	0.17	0/4193	0.30	0/6529
2	D	0.16	0/4193	0.30	0/6529
2	F	0.17	0/4193	0.30	0/6529
2	H	0.17	0/4193	0.30	0/6529
2	J	0.16	0/4193	0.29	0/6529
2	L	0.17	0/4193	0.30	0/6529
All	All	0.19	0/49812	0.36	0/72300

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	4021	4103	4101	82	0
1	C	4021	4103	4101	79	0
1	E	4021	4103	4101	79	0
1	G	4021	4103	4101	75	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	I	4021	4103	4101	83	0
1	K	4021	4103	4101	85	0
2	B	3755	1888	1888	57	0
2	D	3755	1888	1888	56	0
2	F	3755	1888	1888	52	0
2	H	3755	1888	1888	53	0
2	J	3755	1888	1888	52	0
2	L	3755	1888	1888	55	0
All	All	46656	35946	35934	736	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 9.

All (736) 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:D:18:A:N7	2:D:175:G:N2	2.20	0.90
1:A:294:GLU:OE1	1:E:254:TYR:OH	1.97	0.82
1:G:254:TYR:OH	1:K:294:GLU:OE1	1.96	0.82
1:C:121:PHE:CD1	1:C:217:ILE:HG22	2.18	0.79
1:I:121:PHE:CD1	1:I:217:ILE:HG22	2.19	0.78
2:D:67:C:O2'	2:J:67:C:O2'	2.00	0.78
1:I:455:LYS:HE2	1:I:455:LYS:HA	1.67	0.77
1:E:404:VAL:HG13	1:E:470:LEU:HD21	1.67	0.75
2:H:90:U:O2'	2:H:91:G:O5'	2.04	0.74
2:F:90:U:O2'	2:F:91:G:O5'	2.05	0.74
2:B:90:U:O2'	2:B:91:G:O5'	2.06	0.74
2:D:67:C:HO2'	2:J:67:C:HO2'	1.36	0.73
1:I:58:ASP:OD2	1:I:59:LYS:N	2.22	0.73
2:L:90:U:O2'	2:L:91:G:O5'	2.07	0.72
1:A:187:GLU:OE2	1:K:269:LYS:NZ	2.22	0.71
1:G:25:TYR:O	1:G:29:VAL:HG23	1.90	0.71
1:A:25:TYR:O	1:A:29:VAL:HG23	1.91	0.70
1:C:25:TYR:O	1:C:29:VAL:HG23	1.91	0.70
1:G:171:ASP:OD2	1:G:274:LYS:NZ	2.20	0.70
1:K:25:TYR:O	1:K:29:VAL:HG23	1.90	0.70
1:I:25:TYR:O	1:I:29:VAL:HG23	1.91	0.69
2:D:90:U:O2'	2:D:91:G:O5'	2.08	0.69
2:B:90:U:O2	2:B:97:G:O6	2.11	0.69
1:K:121:PHE:CD1	1:K:217:ILE:HG22	2.28	0.69
1:A:1:MET:HE1	1:A:5:GLN:HB3	1.74	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:67:C:O2'	2:L:67:C:O2'	2.12	0.68
1:C:447:GLU:N	1:C:447:GLU:OE1	2.26	0.68
2:J:90:U:O2'	2:J:91:G:O5'	2.08	0.68
2:L:90:U:O2	2:L:97:G:O6	2.11	0.68
1:K:1:MET:HE1	1:K:5:GLN:HB3	1.75	0.68
2:J:90:U:O2	2:J:97:G:O6	2.12	0.67
2:D:90:U:O2	2:D:97:G:O6	2.13	0.67
1:E:25:TYR:O	1:E:29:VAL:HG23	1.94	0.67
2:L:111:G:O2'	2:L:112:U:O2	2.12	0.67
1:A:121:PHE:CD1	1:A:217:ILE:HG22	2.31	0.66
1:I:321:LYS:NZ	2:J:108:C:O3'	2.29	0.66
2:B:111:G:O2'	2:B:112:U:O2	2.11	0.66
2:B:67:C:O2'	2:H:67:C:O2'	2.12	0.66
1:I:447:GLU:OE1	1:I:447:GLU:N	2.29	0.65
2:H:90:U:O2	2:H:97:G:O6	2.14	0.65
2:F:90:U:O2	2:F:97:G:O6	2.15	0.64
2:F:111:G:O2'	2:F:112:U:O2	2.16	0.63
2:D:111:G:O2'	2:D:112:U:O2	2.17	0.63
2:H:111:G:O2'	2:H:112:U:O2	2.16	0.62
1:C:158:LYS:HB3	1:C:158:LYS:HZ2	1.65	0.61
2:H:90:U:HO2'	2:H:91:G:C5'	2.14	0.61
1:G:47:LYS:HD3	1:G:49:ILE:HD11	1.83	0.61
1:G:288:LYS:HG2	1:G:310:VAL:HG11	1.83	0.61
1:C:294:GLU:OE1	1:K:254:TYR:OH	2.19	0.60
2:F:141:C:H3'	2:F:142:G:H21	1.66	0.60
1:A:254:TYR:OH	1:I:294:GLU:OE1	2.17	0.60
2:J:111:G:O2'	2:J:112:U:O2	2.19	0.60
1:I:457:THR:C	1:I:458:ILE:HD12	2.27	0.60
2:D:67:C:HO2'	2:J:67:C:C2'	2.14	0.60
2:H:141:C:H3'	2:H:142:G:H21	1.67	0.60
1:I:355:ASP:OD1	1:I:358:PHE:N	2.34	0.60
2:D:98:C:HO2'	2:D:99:C:H6	1.50	0.60
1:G:26:HIS:HB2	1:G:49:ILE:HD12	1.85	0.59
1:A:164:LEU:HD11	1:A:286:LEU:HD22	1.83	0.59
1:A:265:THR:OG1	1:I:160:GLU:OE2	2.13	0.59
2:B:90:U:HO2'	2:B:91:G:C5'	2.16	0.59
2:D:90:U:HO2'	2:D:91:G:C5'	2.13	0.59
1:G:171:ASP:OD2	1:G:273:ILE:HD11	2.03	0.59
2:D:141:C:H3'	2:D:142:G:H21	1.68	0.59
1:C:187:GLU:OE2	1:E:269:LYS:NZ	2.29	0.59
1:K:179:SER:O	1:K:183:GLU:HG3	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:90:U:HO2'	2:J:91:G:C5'	2.14	0.58
2:J:141:C:H3'	2:J:142:G:H21	1.67	0.58
1:C:121:PHE:CG	1:C:217:ILE:HG22	2.38	0.58
2:D:171:U:H2'	2:D:172:U:H6	1.69	0.58
1:E:288:LYS:HG2	1:E:310:VAL:HG11	1.85	0.58
2:B:67:C:O2'	2:H:67:C:O3'	2.20	0.58
1:E:143:ARG:NH1	1:E:486:GLU:OE2	2.36	0.58
1:G:143:ARG:HD2	1:G:489:MET:HB3	1.85	0.58
2:B:171:U:H2'	2:B:172:U:H6	1.69	0.58
1:E:87:LEU:HD23	1:E:205:ARG:NH1	2.18	0.58
1:G:261:PHE:CD1	1:K:161:ARG:HD3	2.39	0.58
2:F:171:U:H2'	2:F:172:U:H6	1.69	0.57
2:B:53:C:HO2'	2:B:54:U:H6	1.51	0.57
2:F:90:U:HO2'	2:F:91:G:C5'	2.17	0.57
1:G:87:LEU:HD23	1:G:205:ARG:NH1	2.18	0.57
2:H:171:U:H2'	2:H:172:U:H6	1.68	0.57
1:E:164:LEU:HD22	1:E:286:LEU:HD13	1.86	0.57
2:J:171:U:H2'	2:J:172:U:H6	1.69	0.57
1:E:294:GLU:HG3	1:I:283:ILE:HG23	1.87	0.57
2:L:98:C:HO2'	2:L:99:C:H6	1.50	0.57
2:B:170:U:H2'	2:B:171:U:H6	1.70	0.57
1:E:171:ASP:OD2	1:E:273:ILE:HD11	2.04	0.57
2:L:171:U:H2'	2:L:172:U:H6	1.69	0.57
2:J:170:U:H2'	2:J:171:U:H6	1.70	0.57
1:C:164:LEU:CD1	1:C:286:LEU:HD13	2.34	0.56
2:F:67:C:O3'	2:L:67:C:O2'	2.22	0.56
1:I:121:PHE:CG	1:I:217:ILE:HG22	2.40	0.56
2:F:170:U:H2'	2:F:171:U:H6	1.71	0.56
1:I:375:LYS:NZ	2:J:80:G:O6	2.39	0.56
2:H:171:U:C2	2:H:172:U:C5	2.93	0.56
2:D:170:U:H2'	2:D:171:U:H6	1.70	0.56
1:C:161:ARG:HG2	1:C:287:THR:HG22	1.86	0.55
2:D:171:U:C2	2:D:172:U:C5	2.94	0.55
1:K:164:LEU:HD11	1:K:286:LEU:HD22	1.88	0.55
2:L:90:U:HO2'	2:L:91:G:C5'	2.17	0.55
1:I:439:ASN:O	1:I:443:ILE:HG22	2.06	0.55
2:H:170:U:H2'	2:H:171:U:H6	1.70	0.55
1:A:1:MET:HE1	1:A:5:GLN:CB	2.36	0.55
1:C:283:ILE:HG23	1:G:294:GLU:HG3	1.88	0.55
2:F:171:U:C2	2:F:172:U:C5	2.95	0.55
2:L:170:U:H2'	2:L:171:U:H6	1.71	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:141:C:H3'	2:L:142:G:H21	1.72	0.55
2:B:141:C:H3'	2:B:142:G:H21	1.72	0.55
1:I:161:ARG:NH2	1:I:294:GLU:OE2	2.40	0.55
2:F:68:G:OP2	2:L:68:G:OP2	2.25	0.54
2:B:170:U:H2'	2:B:171:U:C6	2.43	0.54
2:F:67:C:HO2'	2:L:67:C:C2'	2.20	0.54
2:B:171:U:C2	2:B:172:U:C5	2.96	0.54
1:C:95:ASP:HB3	1:C:103:ILE:HG12	1.90	0.54
1:E:10:VAL:HG12	1:E:115:ALA:HB1	1.90	0.54
2:D:170:U:H2'	2:D:171:U:C6	2.43	0.54
1:G:439:ASN:O	1:G:443:ILE:HG22	2.07	0.54
2:J:171:U:C2	2:J:172:U:C5	2.96	0.54
1:A:294:GLU:HB3	1:E:295:ILE:HB	1.90	0.54
1:C:257:ILE:CG2	1:C:283:ILE:HD12	2.38	0.54
2:F:122:C:H2'	2:F:123:U:O4'	2.08	0.54
2:J:170:U:H2'	2:J:171:U:C6	2.43	0.54
1:K:288:LYS:HG3	1:K:310:VAL:HG11	1.89	0.53
2:B:68:G:OP2	2:H:68:G:OP2	2.25	0.53
1:E:439:ASN:O	1:E:443:ILE:HG22	2.08	0.53
1:G:10:VAL:HG12	1:G:115:ALA:HB1	1.89	0.53
2:H:122:C:H2'	2:H:123:U:O4'	2.09	0.53
1:C:461:ARG:HE	1:C:463:LEU:HD21	1.73	0.53
2:D:123:U:O2'	2:D:126:U:OP2	2.25	0.53
1:K:335:LEU:HB3	1:K:425:ARG:HD2	1.91	0.53
1:A:288:LYS:HG3	1:A:310:VAL:HG11	1.89	0.53
2:F:170:U:H2'	2:F:171:U:C6	2.44	0.53
2:L:171:U:C2	2:L:172:U:C5	2.97	0.53
2:H:170:U:H2'	2:H:171:U:C6	2.44	0.53
1:K:490:ASN:OD1	1:K:490:ASN:O	2.27	0.53
1:E:171:ASP:OD2	1:E:274:LYS:NZ	2.19	0.53
1:A:490:ASN:O	1:A:490:ASN:OD1	2.27	0.53
1:G:378:ILE:CG2	1:G:471:LEU:HD22	2.39	0.53
2:B:67:C:C2'	2:H:67:C:HO2'	2.21	0.52
2:J:122:C:H2'	2:J:123:U:O4'	2.09	0.52
2:L:170:U:H2'	2:L:171:U:C6	2.43	0.52
1:A:294:GLU:HG3	1:E:283:ILE:HG23	1.91	0.52
2:B:3:U:O2'	2:B:4:C:OP1	2.25	0.52
1:G:168:ASP:OD1	1:G:278:LYS:NZ	2.31	0.52
1:I:142:ASP:HB3	1:I:145:VAL:HG23	1.92	0.52
2:B:122:C:H2'	2:B:123:U:O4'	2.10	0.52
1:K:142:ASP:HB3	1:K:145:VAL:HG23	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:375:LYS:NZ	2:L:80:G:O6	2.42	0.52
2:D:122:C:H2'	2:D:123:U:O4'	2.09	0.52
1:E:58:ASP:OD2	1:E:59:LYS:N	2.43	0.52
2:B:40:A:N7	2:B:41:G:N2	2.58	0.52
1:E:330:ILE:O	1:E:334:ASN:ND2	2.41	0.52
1:C:47:LYS:HD3	1:C:49:ILE:HD13	1.92	0.52
1:G:295:ILE:HB	1:K:294:GLU:HB3	1.91	0.52
2:H:123:U:O2'	2:H:126:U:OP2	2.27	0.52
2:B:68:G:H8	1:C:182:ASN:ND2	2.07	0.51
1:C:15:ALA:O	1:C:19:ILE:HG13	2.10	0.51
1:C:129:ARG:HB3	1:C:129:ARG:NH1	2.24	0.51
2:L:122:C:H2'	2:L:123:U:O4'	2.10	0.51
1:A:142:ASP:HB3	1:A:145:VAL:HG23	1.90	0.51
1:K:87:LEU:HD22	1:K:88:PRO:HD2	1.92	0.51
1:G:283:ILE:HG23	1:K:294:GLU:HG3	1.91	0.51
2:H:27:C:C2	2:H:28:C:C6	2.99	0.51
1:E:373:LEU:HD22	1:E:389:LEU:HD11	1.91	0.51
1:E:180:PHE:O	1:E:184:GLN:HG2	2.11	0.51
2:J:40:A:N7	2:J:41:G:N2	2.59	0.51
1:A:276:ASN:OD1	1:A:276:ASN:N	2.44	0.51
1:C:143:ARG:HD2	1:C:489:MET:HB3	1.93	0.51
1:C:180:PHE:O	1:C:184:GLN:HG2	2.10	0.51
1:K:13:GLU:HA	1:K:13:GLU:OE1	2.10	0.51
1:K:145:VAL:HG11	1:K:304:LEU:HD21	1.92	0.51
2:H:53:C:HO2'	2:H:54:U:H6	1.58	0.51
1:A:13:GLU:HA	1:A:13:GLU:OE1	2.11	0.51
2:D:67:C:C2'	2:J:67:C:HO2'	2.23	0.51
1:G:373:LEU:HD22	1:G:389:LEU:HD11	1.92	0.51
2:J:3:U:O2'	2:J:4:C:OP1	2.29	0.51
1:K:1:MET:HE1	1:K:5:GLN:CB	2.41	0.51
1:K:206:GLU:OE1	1:K:206:GLU:HA	2.11	0.51
1:I:164:LEU:CD1	1:I:286:LEU:HD13	2.41	0.50
1:I:182:ASN:ND2	2:L:68:G:H8	2.09	0.50
1:I:246:ASP:OD1	1:I:278:LYS:NZ	2.43	0.50
2:D:162:C:H2'	2:D:163:U:H6	1.75	0.50
2:F:27:C:C2	2:F:28:C:C6	2.99	0.50
1:A:172:PHE:CD2	1:A:273:ILE:HG23	2.46	0.50
1:C:294:GLU:CD	1:K:254:TYR:HH	2.18	0.50
1:I:268:SER:OG	1:I:275:ILE:HG13	2.12	0.50
2:F:3:U:O2'	2:F:4:C:OP1	2.28	0.50
1:C:300:ASN:C	1:C:300:ASN:OD1	2.54	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:378:ILE:HG23	1:C:475:ALA:HB2	1.94	0.50
1:G:87:LEU:HD22	1:G:88:PRO:HD2	1.94	0.50
2:J:162:C:H2'	2:J:163:U:H6	1.76	0.50
1:A:375:LYS:NZ	2:B:80:G:O6	2.44	0.50
2:H:162:C:H2'	2:H:163:U:H6	1.77	0.50
1:E:26:HIS:HB2	1:E:49:ILE:HD12	1.93	0.50
2:B:98:C:HO2'	2:B:99:C:H6	1.56	0.50
1:C:164:LEU:HD22	1:C:250:TRP:HB3	1.94	0.50
1:E:94:LYS:O	1:E:106:VAL:HG12	2.12	0.50
1:G:96:VAL:HG22	1:G:104:ARG:O	2.12	0.50
1:K:378:ILE:HG23	1:K:475:ALA:HB2	1.94	0.50
2:L:27:C:C2	2:L:28:C:C6	3.00	0.50
2:L:162:C:H2'	2:L:163:U:H6	1.77	0.50
2:F:162:C:H2'	2:F:163:U:H6	1.77	0.50
2:J:27:C:C2	2:J:28:C:C6	2.99	0.50
2:B:123:U:O2'	2:B:126:U:OP2	2.29	0.49
1:I:30:HIS:O	1:I:34:VAL:HG23	2.11	0.49
1:K:276:ASN:N	1:K:276:ASN:OD1	2.44	0.49
2:B:27:C:C2	2:B:28:C:C6	3.00	0.49
2:D:27:C:C2	2:D:28:C:C6	3.00	0.49
1:G:180:PHE:O	1:G:184:GLN:HG2	2.11	0.49
2:L:3:U:O2'	2:L:4:C:OP1	2.25	0.49
2:D:40:A:N7	2:D:41:G:N2	2.60	0.49
1:I:276:ASN:N	1:I:276:ASN:OD1	2.45	0.49
1:K:378:ILE:CG2	1:K:471:LEU:HD22	2.42	0.49
2:B:162:C:H2'	2:B:163:U:H6	1.77	0.49
2:B:171:U:H2'	2:B:172:U:C6	2.48	0.49
1:E:161:ARG:NH2	1:E:294:GLU:OE2	2.44	0.49
2:F:25:G:H2'	2:F:26:G:H8	1.76	0.49
1:G:58:ASP:OD2	1:G:59:LYS:N	2.46	0.49
1:K:171:ASP:OD2	1:K:273:ILE:HD11	2.12	0.49
2:L:171:U:H2'	2:L:172:U:C6	2.48	0.49
1:A:439:ASN:O	1:A:443:ILE:HG22	2.12	0.49
2:J:27:C:H42	2:J:133:G:H1	1.60	0.49
1:A:96:VAL:HG22	1:A:104:ARG:O	2.12	0.49
1:E:378:ILE:HG23	1:E:475:ALA:HB2	1.94	0.49
1:E:87:LEU:HD22	1:E:88:PRO:HD2	1.95	0.49
1:E:411:ASP:OD1	1:E:468:SER:OG	2.26	0.49
2:F:123:U:O2'	2:F:126:U:OP2	2.28	0.49
2:H:3:U:HO2'	2:H:4:C:P	2.36	0.49
1:K:1:MET:HE3	1:K:1:MET:CA	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:40:A:N7	2:L:41:G:N2	2.61	0.49
2:D:3:U:O2'	2:D:4:C:OP1	2.29	0.49
1:E:168:ASP:OD1	1:E:278:LYS:NZ	2.28	0.49
1:K:96:VAL:HG22	1:K:104:ARG:O	2.13	0.49
1:A:145:VAL:HG11	1:A:304:LEU:HD21	1.93	0.48
1:C:161:ARG:NH2	1:C:294:GLU:OE2	2.44	0.48
1:I:378:ILE:HG23	1:I:475:ALA:HB2	1.95	0.48
2:D:27:C:H42	2:D:133:G:H1	1.60	0.48
2:D:172:U:C2	2:D:173:U:C5	3.01	0.48
1:G:330:ILE:O	1:G:334:ASN:ND2	2.40	0.48
1:C:463:LEU:HD23	1:C:463:LEU:N	2.29	0.48
1:E:47:LYS:HE2	1:E:49:ILE:HD11	1.94	0.48
1:I:164:LEU:HD22	1:I:250:TRP:HB3	1.96	0.48
1:K:19:ILE:HG13	1:K:63:PRO:HB2	1.96	0.48
2:F:3:U:HO2'	2:F:4:C:P	2.37	0.48
1:K:316:LYS:O	1:K:320:VAL:HG23	2.14	0.48
1:C:316:LYS:O	1:C:320:VAL:HG23	2.14	0.48
1:E:212:PRO:CG	1:E:217:ILE:HD11	2.44	0.48
1:E:316:LYS:O	1:E:320:VAL:HG23	2.14	0.48
1:I:288:LYS:HG2	1:I:310:VAL:HG11	1.94	0.48
1:A:378:ILE:HG23	1:A:475:ALA:HB2	1.95	0.48
1:K:161:ARG:NH2	1:K:294:GLU:OE2	2.45	0.48
1:C:276:ASN:OD1	1:C:276:ASN:N	2.46	0.48
1:E:6:GLN:O	1:E:10:VAL:HG23	2.13	0.48
1:G:6:GLN:O	1:G:10:VAL:HG23	2.14	0.48
1:G:161:ARG:NH2	1:G:294:GLU:OE2	2.44	0.48
2:J:98:C:HO2'	2:J:99:C:H6	1.55	0.48
2:L:3:U:HO2'	2:L:4:C:P	2.37	0.48
2:F:98:C:HO2'	2:F:99:C:H6	1.59	0.48
2:H:25:G:H2'	2:H:26:G:H8	1.77	0.48
1:A:171:ASP:OD2	1:A:273:ILE:HD11	2.13	0.48
2:L:3:U:O2'	2:L:4:C:P	2.72	0.47
1:A:246:ASP:OD1	1:A:278:LYS:NZ	2.46	0.47
1:A:335:LEU:HB3	1:A:425:ARG:HD2	1.95	0.47
1:E:335:LEU:HB3	1:E:425:ARG:HD2	1.96	0.47
1:G:326:GLN:O	1:G:330:ILE:HD12	2.14	0.47
1:I:172:PHE:CD1	1:I:275:ILE:HG12	2.49	0.47
2:J:172:U:C2	2:J:173:U:C5	3.02	0.47
1:A:283:ILE:HG21	1:I:294:GLU:OE2	2.14	0.47
1:K:190:PHE:CE2	1:K:224:LEU:HD12	2.48	0.47
1:C:335:LEU:HB3	1:C:425:ARG:HD2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:161:ARG:HD3	1:I:261:PHE:CE1	2.50	0.47
1:G:316:LYS:O	1:G:320:VAL:HG23	2.14	0.47
2:H:91:G:O2'	2:H:92:U:OP2	2.31	0.47
1:I:180:PHE:O	1:I:184:GLN:HG2	2.13	0.47
1:A:145:VAL:HG11	1:A:304:LEU:CD2	2.44	0.47
1:G:212:PRO:CG	1:G:217:ILE:HD11	2.44	0.47
2:J:123:U:O2'	2:J:126:U:OP2	2.28	0.47
2:B:3:U:O2'	2:B:4:C:P	2.73	0.47
1:E:162:THR:HG22	1:E:164:LEU:CD1	2.45	0.47
1:I:335:LEU:HB3	1:I:425:ARG:HD2	1.97	0.47
2:B:3:U:HO2'	2:B:4:C:P	2.37	0.47
1:C:58:ASP:OD2	1:C:59:LYS:N	2.47	0.47
1:C:211:ILE:HG21	1:C:221:LEU:HB2	1.97	0.47
2:D:172:U:H2'	2:D:173:U:C6	2.50	0.47
2:F:171:U:H2'	2:F:172:U:C6	2.49	0.47
2:H:171:U:H2'	2:H:172:U:C6	2.49	0.47
1:I:316:LYS:O	1:I:320:VAL:HG23	2.14	0.47
1:I:440:PHE:CE1	1:I:456:LYS:HD2	2.49	0.47
2:J:28:C:H2'	2:J:29:U:H5'	1.97	0.47
1:K:145:VAL:HG11	1:K:304:LEU:CD2	2.44	0.47
2:L:172:U:C2	2:L:173:U:C5	3.03	0.47
2:B:60:U:H3	2:B:73:A:H61	1.62	0.47
1:G:268:SER:OG	1:G:275:ILE:HG13	2.14	0.47
1:G:378:ILE:HG23	1:G:475:ALA:HB2	1.96	0.47
1:A:19:ILE:HG13	1:A:63:PRO:HB2	1.96	0.47
1:K:180:PHE:O	1:K:184:GLN:HG2	2.15	0.47
1:A:211:ILE:HG21	1:A:221:LEU:HB2	1.97	0.47
1:A:316:LYS:O	1:A:320:VAL:HG23	2.15	0.47
1:E:268:SER:OG	1:E:275:ILE:HG13	2.15	0.47
1:G:23:HIS:HA	1:G:26:HIS:NE2	2.30	0.47
2:H:40:A:N7	2:H:41:G:N2	2.63	0.47
2:H:43:A:H4'	2:H:44:A:O5'	2.15	0.47
1:K:18:LEU:HD13	1:K:111:ILE:HD11	1.97	0.47
1:K:27:ASN:ND2	2:L:140:C:H1'	2.29	0.47
2:L:27:C:H42	2:L:133:G:H1	1.63	0.47
2:L:60:U:H3	2:L:73:A:H61	1.62	0.47
2:B:43:A:H4'	2:B:44:A:O5'	2.15	0.46
1:C:27:ASN:ND2	2:D:140:C:H1'	2.30	0.46
1:G:186:ASN:OD1	1:G:186:ASN:C	2.58	0.46
1:I:95:ASP:HB3	1:I:103:ILE:CG1	2.45	0.46
1:K:211:ILE:HG21	1:K:221:LEU:HB2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:PHE:O	1:A:184:GLN:HG2	2.16	0.46
2:B:27:C:H42	2:B:133:G:H1	1.62	0.46
2:B:172:U:C2	2:B:173:U:C5	3.04	0.46
2:F:43:A:H4'	2:F:44:A:O5'	2.15	0.46
1:G:246:ASP:OD1	1:G:278:LYS:CE	2.63	0.46
2:H:3:U:O2'	2:H:4:C:OP1	2.28	0.46
1:A:95:ASP:HB3	1:A:103:ILE:HG12	1.98	0.46
1:C:257:ILE:HG22	1:C:283:ILE:HD12	1.97	0.46
1:C:261:PHE:CE1	1:G:161:ARG:HD3	2.49	0.46
2:J:3:U:O2'	2:J:4:C:P	2.73	0.46
2:J:43:A:H4'	2:J:44:A:O5'	2.15	0.46
2:B:90:U:H2'	2:B:91:G:C8	2.51	0.46
2:D:3:U:O2'	2:D:4:C:P	2.74	0.46
2:D:43:A:H4'	2:D:44:A:O5'	2.15	0.46
1:E:326:GLN:O	1:E:330:ILE:HD12	2.16	0.46
2:L:43:A:H4'	2:L:44:A:O5'	2.15	0.46
1:E:95:ASP:HB3	1:E:103:ILE:HG12	1.97	0.46
2:F:172:U:C2	2:F:173:U:C5	3.04	0.46
1:I:261:PHE:HZ	1:I:277:PRO:HG3	1.79	0.46
1:K:28:ARG:HH11	1:K:28:ARG:HG2	1.81	0.46
1:K:58:ASP:OD2	1:K:59:LYS:N	2.48	0.46
1:K:164:LEU:HD13	1:K:301:LEU:HD22	1.98	0.46
1:A:300:ASN:OD1	1:A:300:ASN:C	2.59	0.46
2:H:172:U:C2	2:H:173:U:C5	3.04	0.46
1:I:19:ILE:HG13	1:I:63:PRO:HB2	1.97	0.46
2:J:172:U:H2'	2:J:173:U:C6	2.51	0.46
2:L:28:C:H2'	2:L:29:U:H5'	1.97	0.46
1:C:224:LEU:HD12	1:C:224:LEU:O	2.16	0.46
1:E:182:ASN:ND2	2:H:68:G:H8	2.13	0.46
1:E:286:LEU:HG	1:E:298:LYS:O	2.15	0.46
2:H:3:U:O2'	2:H:4:C:P	2.73	0.46
1:K:95:ASP:HB3	1:K:103:ILE:HG12	1.98	0.46
1:A:304:LEU:O	1:A:316:LYS:HE3	2.15	0.46
2:B:172:U:H2'	2:B:173:U:C6	2.51	0.46
1:C:160:GLU:OE2	1:K:265:THR:OG1	2.25	0.46
1:C:233:LEU:O	1:C:238:VAL:HG22	2.16	0.46
1:G:335:LEU:HB3	1:G:425:ARG:HD2	1.98	0.46
1:I:161:ARG:HG2	1:I:287:THR:HG22	1.98	0.46
2:B:28:C:H2'	2:B:29:U:H5'	1.98	0.46
1:C:261:PHE:HZ	1:C:277:PRO:HG3	1.81	0.46
1:E:23:HIS:HA	1:E:26:HIS:NE2	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:3:U:O2'	2:F:4:C:P	2.73	0.46
2:F:68:G:H8	1:G:182:ASN:ND2	2.13	0.46
1:G:95:ASP:HB3	1:G:103:ILE:HG12	1.96	0.46
1:G:162:THR:HG22	1:G:164:LEU:CD1	2.46	0.46
1:G:164:LEU:HD22	1:G:286:LEU:HD13	1.97	0.46
1:G:411:ASP:O	1:G:415:VAL:HG23	2.15	0.46
1:I:246:ASP:OD1	1:I:278:LYS:CE	2.64	0.46
2:L:90:U:H2'	2:L:91:G:C8	2.51	0.46
1:A:1:MET:C	1:A:1:MET:HE3	2.42	0.45
1:A:27:ASN:ND2	2:B:140:C:H1'	2.31	0.45
1:A:308:LEU:HD22	1:A:313:VAL:HG22	1.98	0.45
2:B:94:A:C5	2:F:113:G:C6	3.04	0.45
1:E:233:LEU:HD22	1:E:238:VAL:HG21	1.98	0.45
1:I:243:TYR:CD2	1:I:304:LEU:HD11	2.51	0.45
1:I:480:LEU:HG	1:I:488:ILE:HD13	1.97	0.45
2:L:172:U:H2'	2:L:173:U:C6	2.51	0.45
2:D:60:U:H3	2:D:73:A:H61	1.64	0.45
1:E:411:ASP:O	1:E:415:VAL:HG23	2.16	0.45
2:L:123:U:O2'	2:L:126:U:OP2	2.29	0.45
1:C:95:ASP:HB3	1:C:103:ILE:CG1	2.46	0.45
1:C:145:VAL:HG11	1:C:304:LEU:HD21	1.98	0.45
2:D:28:C:H2'	2:D:29:U:H5'	1.98	0.45
2:F:40:A:N7	2:F:41:G:N2	2.64	0.45
2:B:175:G:C2	2:B:176:U:C4	3.05	0.45
1:I:176:ILE:CD1	1:I:271:ALA:HB1	2.46	0.45
2:J:90:U:H2'	2:J:91:G:C8	2.51	0.45
1:K:87:LEU:HD23	1:K:205:ARG:NH1	2.31	0.45
1:A:1:MET:HE2	1:A:6:GLN:CG	2.47	0.45
1:A:18:LEU:HD13	1:A:111:ILE:HD11	1.97	0.45
1:A:254:TYR:HH	1:I:294:GLU:CD	2.22	0.45
1:A:490:ASN:OD1	1:A:490:ASN:C	2.59	0.45
1:G:445:ASP:OD1	1:G:446:ARG:N	2.50	0.45
2:J:60:U:H3	2:J:73:A:H61	1.64	0.45
1:K:347:GLN:O	1:K:349:ILE:HD13	2.16	0.45
2:L:175:G:C2	2:L:176:U:C4	3.05	0.45
1:A:43:SER:OG	1:A:341:LYS:HD2	2.17	0.45
1:G:233:LEU:HD22	1:G:238:VAL:HG21	1.97	0.45
1:I:224:LEU:O	1:I:224:LEU:HD12	2.17	0.45
1:E:186:ASN:OD1	1:E:186:ASN:C	2.58	0.45
1:G:351:ALA:O	1:G:354:ARG:HG2	2.17	0.45
1:K:340:LYS:O	1:K:340:LYS:HG2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:172:U:H2'	2:H:173:U:C6	2.52	0.45
2:J:29:U:C2	2:J:30:U:C5	3.05	0.45
1:A:369:MET:HA	1:A:394:ILE:HG22	1.99	0.45
2:D:68:G:OP2	2:J:68:G:P	2.75	0.45
1:E:445:ASP:OD1	1:E:446:ARG:N	2.50	0.45
1:I:169:PHE:CE2	1:I:275:ILE:HD13	2.52	0.45
1:K:1:MET:HE2	1:K:6:GLN:HG3	1.98	0.45
2:L:76:U:HO2'	2:L:77:G:P	2.39	0.45
2:D:175:G:C2	2:D:176:U:C4	3.05	0.45
1:G:19:ILE:HG13	1:G:63:PRO:HB2	1.99	0.45
2:H:28:C:H2'	2:H:29:U:H5'	1.99	0.45
1:K:399:PRO:HB2	1:K:488:ILE:HD11	1.98	0.45
2:L:107:C:H2'	2:L:108:C:C6	2.52	0.45
1:A:161:ARG:NH2	1:A:294:GLU:OE2	2.47	0.44
1:A:243:TYR:CD2	1:A:304:LEU:HD11	2.52	0.44
1:G:444:LEU:HD21	1:G:452:LYS:HB3	1.99	0.44
1:K:1:MET:HE2	1:K:6:GLN:CG	2.47	0.44
1:K:284:SER:OG	1:K:298:LYS:HE3	2.16	0.44
1:A:235:ARG:CZ	1:A:235:ARG:HB3	2.47	0.44
2:D:68:G:P	2:J:68:G:OP2	2.75	0.44
1:E:246:ASP:OD1	1:E:278:LYS:CE	2.65	0.44
2:F:60:U:H3	2:F:73:A:H61	1.65	0.44
1:G:261:PHE:HZ	1:G:277:PRO:HG3	1.82	0.44
1:K:43:SER:OG	1:K:341:LYS:HD2	2.17	0.44
1:K:304:LEU:O	1:K:316:LYS:HE3	2.17	0.44
1:A:1:MET:HE3	1:A:1:MET:CA	2.47	0.44
2:F:28:C:H2'	2:F:29:U:H5'	1.98	0.44
2:F:172:U:H2'	2:F:173:U:C6	2.52	0.44
2:H:27:C:H42	2:H:133:G:H1	1.65	0.44
1:K:308:LEU:HD22	1:K:313:VAL:HG22	1.98	0.44
1:K:490:ASN:OD1	1:K:490:ASN:C	2.60	0.44
1:A:283:ILE:HG23	1:I:294:GLU:HG3	1.98	0.44
1:A:391:PHE:CD1	1:A:476:LEU:HD12	2.52	0.44
1:E:161:ARG:HG2	1:E:287:THR:HG22	1.99	0.44
2:F:175:G:C2	2:F:176:U:C4	3.06	0.44
2:L:53:C:HO2'	2:L:54:U:H6	1.65	0.44
1:E:19:ILE:HG13	1:E:63:PRO:HB2	1.98	0.44
1:G:51:ARG:O	1:G:52:PRO:C	2.61	0.44
1:C:461:ARG:HB2	1:C:463:LEU:HD21	2.00	0.44
2:D:76:U:HO2'	2:D:77:G:P	2.40	0.44
2:F:169:U:H2'	2:F:170:U:C6	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:76:U:HO2'	2:H:77:G:P	2.40	0.44
1:I:27:ASN:ND2	2:J:140:C:H1'	2.32	0.44
1:K:440:PHE:CE2	1:K:456:LYS:HD2	2.53	0.44
1:K:480:LEU:HG	1:K:488:ILE:HD13	1.99	0.44
1:C:351:ALA:O	1:C:354:ARG:HG2	2.17	0.44
1:E:371:GLY:O	1:E:390:TYR:OH	2.35	0.44
2:H:76:U:O2'	2:H:77:G:H8	2.01	0.44
2:J:76:U:O2'	2:J:77:G:H8	2.01	0.44
2:J:171:U:H2'	2:J:172:U:C6	2.50	0.44
1:A:58:ASP:OD2	1:A:59:LYS:N	2.51	0.44
1:C:167:PHE:HA	1:C:277:PRO:HA	1.99	0.44
2:D:68:G:OP2	2:J:68:G:OP2	2.34	0.44
2:F:91:G:O2'	2:F:92:U:OP2	2.31	0.44
1:I:6:GLN:O	1:I:10:VAL:HG22	2.17	0.44
1:I:233:LEU:O	1:I:238:VAL:HG22	2.17	0.44
1:I:317:GLU:OE1	1:I:321:LYS:HE2	2.17	0.44
1:K:300:ASN:OD1	1:K:300:ASN:O	2.36	0.44
2:L:100:G:O2'	2:L:101:U:O5'	2.35	0.44
1:A:284:SER:OG	1:A:298:LYS:HE3	2.18	0.44
1:C:261:PHE:CD1	1:G:161:ARG:HD3	2.53	0.44
1:C:283:ILE:HG21	1:G:294:GLU:OE2	2.18	0.44
2:D:72:C:H2'	2:D:73:A:O4'	2.17	0.44
2:J:72:C:H2'	2:J:73:A:O4'	2.17	0.44
2:J:175:G:C2	2:J:176:U:C4	3.06	0.44
2:L:99:C:O2'	2:L:100:G:H5'	2.18	0.44
1:A:176:ILE:CD1	1:A:271:ALA:HB1	2.48	0.43
1:A:233:LEU:O	1:A:238:VAL:HG22	2.17	0.43
1:C:165:ALA:HA	1:C:282:GLY:O	2.18	0.43
1:G:300:ASN:OD1	1:G:300:ASN:O	2.36	0.43
1:G:378:ILE:HG22	1:G:471:LEU:HD22	1.99	0.43
2:H:60:U:H3	2:H:73:A:H61	1.66	0.43
1:I:323:ILE:HG12	1:I:394:ILE:HG12	2.00	0.43
1:K:164:LEU:CD1	1:K:286:LEU:HD13	2.48	0.43
1:C:408:LYS:HB2	1:C:408:LYS:HE3	1.82	0.43
2:D:171:U:H2'	2:D:172:U:C6	2.49	0.43
1:E:304:LEU:O	1:E:316:LYS:HE3	2.18	0.43
1:A:10:VAL:HG12	1:A:115:ALA:HB1	2.00	0.43
1:A:404:VAL:HG23	1:A:405:GLU:OE2	2.19	0.43
1:C:6:GLN:O	1:C:10:VAL:HG22	2.18	0.43
1:E:176:ILE:HG22	1:E:210:GLY:HA2	2.00	0.43
1:G:243:TYR:CD2	1:G:304:LEU:HD11	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:304:LEU:O	1:G:316:LYS:HE3	2.17	0.43
2:H:66:U:OP2	2:H:67:C:N4	2.52	0.43
1:I:145:VAL:HG11	1:I:304:LEU:HD21	2.00	0.43
2:J:169:U:H2'	2:J:170:U:C6	2.53	0.43
2:L:76:U:O2'	2:L:77:G:H8	2.00	0.43
1:A:339:LEU:HD22	1:A:344:LEU:HD21	1.99	0.43
1:C:304:LEU:O	1:C:316:LYS:HE3	2.18	0.43
2:D:90:U:H2'	2:D:91:G:C8	2.53	0.43
2:D:169:U:H2'	2:D:170:U:C6	2.53	0.43
1:E:141:ASN:HB3	2:F:125:U:H1'	2.01	0.43
1:E:243:TYR:CD2	1:E:304:LEU:HD11	2.53	0.43
1:E:375:LYS:NZ	1:E:466:ILE:O	2.46	0.43
2:F:27:C:H42	2:F:133:G:H1	1.66	0.43
2:H:169:U:H2'	2:H:170:U:C6	2.53	0.43
1:K:369:MET:HA	1:K:394:ILE:HG22	1.99	0.43
2:B:76:U:O2'	2:B:77:G:H8	2.01	0.43
1:C:243:TYR:CD2	1:C:304:LEU:HD11	2.53	0.43
1:C:294:GLU:OE2	1:K:283:ILE:HG21	2.18	0.43
1:C:323:ILE:HG12	1:C:394:ILE:HG12	1.99	0.43
1:G:95:ASP:HB3	1:G:103:ILE:CG1	2.48	0.43
1:I:399:PRO:HA	1:I:476:LEU:HD21	2.00	0.43
1:I:445:ASP:OD1	1:I:446:ARG:N	2.52	0.43
2:J:26:G:H2'	2:J:27:C:H6	1.84	0.43
2:B:169:U:H2'	2:B:170:U:C6	2.53	0.43
1:C:164:LEU:HD12	1:C:286:LEU:HD13	1.99	0.43
1:C:233:LEU:HD22	1:C:238:VAL:HG21	1.99	0.43
2:D:161:A:H2'	2:D:162:C:H6	1.83	0.43
1:E:51:ARG:O	1:E:52:PRO:C	2.62	0.43
2:F:90:U:H2'	2:F:91:G:C8	2.53	0.43
1:I:434:TYR:CD1	1:I:434:TYR:N	2.87	0.43
1:A:137:TYR:CG	1:A:145:VAL:HG22	2.54	0.43
2:B:96:A:C2	2:B:97:G:C8	3.07	0.43
2:B:99:C:O2'	2:B:100:G:H5'	2.18	0.43
1:C:87:LEU:HD12	1:C:88:PRO:HD2	2.00	0.43
1:E:97:PRO:HA	1:E:103:ILE:HD12	2.01	0.43
1:E:300:ASN:OD1	1:E:300:ASN:O	2.37	0.43
1:I:18:LEU:HD13	1:I:111:ILE:HD11	2.01	0.43
1:K:1:MET:HE3	1:K:1:MET:C	2.44	0.43
1:K:176:ILE:CD1	1:K:271:ALA:HB1	2.49	0.43
1:C:26:HIS:HD1	2:D:140:C:HO2'	1.65	0.43
2:H:175:G:C2	2:H:176:U:C4	3.06	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:143:ARG:HD3	1:I:489:MET:HB3	1.99	0.43
2:L:25:G:H2'	2:L:26:G:H8	1.82	0.43
1:A:399:PRO:CB	1:A:488:ILE:HD11	2.49	0.43
2:B:161:A:H2'	2:B:162:C:H6	1.84	0.43
1:C:97:PRO:HA	1:C:103:ILE:HD12	2.00	0.43
2:B:26:G:H2'	2:B:27:C:H6	1.84	0.43
1:C:37:LYS:HB3	1:C:37:LYS:NZ	2.34	0.43
1:C:176:ILE:CD1	1:C:271:ALA:HB1	2.48	0.43
1:C:227:TRP:CH2	1:C:228:LYS:HD3	2.54	0.43
2:F:76:U:O2'	2:F:77:G:H8	2.01	0.43
1:G:141:ASN:HB3	2:H:125:U:H1'	2.00	0.43
1:K:233:LEU:O	1:K:238:VAL:HG22	2.19	0.43
1:A:212:PRO:HG3	1:A:217:ILE:HD11	2.01	0.42
2:B:100:G:O2'	2:B:101:U:O5'	2.36	0.42
1:C:138:ALA:HB2	1:C:219:LEU:HB2	2.01	0.42
1:E:444:LEU:HD21	1:E:452:LYS:HB3	2.00	0.42
1:I:304:LEU:O	1:I:316:LYS:HE3	2.18	0.42
1:A:373:LEU:HD22	1:A:389:LEU:HD11	2.01	0.42
1:A:419:TYR:CD2	1:A:446:ARG:HB2	2.54	0.42
2:B:25:G:H2'	2:B:26:G:H8	1.83	0.42
2:B:107:C:H2'	2:B:108:C:C6	2.54	0.42
1:E:294:GLU:OE2	1:I:283:ILE:HG21	2.19	0.42
1:G:176:ILE:HG22	1:G:210:GLY:HA2	2.00	0.42
1:G:233:LEU:O	1:G:238:VAL:HG22	2.20	0.42
1:I:211:ILE:HG21	1:I:221:LEU:HB2	2.02	0.42
1:K:161:ARG:HG2	1:K:287:THR:HG22	2.01	0.42
1:K:243:TYR:CD2	1:K:304:LEU:HD11	2.54	0.42
1:K:399:PRO:HA	1:K:476:LEU:HD21	2.00	0.42
2:L:169:U:H2'	2:L:170:U:C6	2.54	0.42
1:A:212:PRO:CG	1:A:217:ILE:HD11	2.49	0.42
1:C:164:LEU:HD22	1:C:250:TRP:CB	2.49	0.42
1:E:95:ASP:HB3	1:E:103:ILE:CG1	2.49	0.42
2:F:66:U:OP2	2:F:67:C:N4	2.52	0.42
1:G:10:VAL:HG21	1:G:119:LEU:HB2	2.00	0.42
1:I:488:ILE:HG13	1:I:489:MET:N	2.34	0.42
2:L:66:U:OP2	2:L:67:C:N4	2.52	0.42
2:B:66:U:OP2	2:B:67:C:N4	2.52	0.42
2:B:121:A:C6	2:B:122:C:N4	2.88	0.42
1:C:158:LYS:HB3	1:C:158:LYS:NZ	2.34	0.42
1:E:351:ALA:O	1:E:354:ARG:HG2	2.19	0.42
1:E:395:MET:HE2	1:E:395:MET:HA	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:134:U:O2'	2:F:135:U:C6	2.70	0.42
1:A:138:ALA:HB2	1:A:219:LEU:HB2	2.00	0.42
1:C:284:SER:OG	1:C:298:LYS:HE3	2.18	0.42
1:E:138:ALA:HB2	1:E:219:LEU:HB2	2.02	0.42
1:G:138:ALA:HB2	1:G:219:LEU:HB2	2.02	0.42
1:I:284:SER:OG	1:I:298:LYS:HE3	2.19	0.42
1:I:458:ILE:HD12	1:I:458:ILE:N	2.34	0.42
1:K:87:LEU:HD23	1:K:205:ARG:CZ	2.49	0.42
1:A:1:MET:HE2	1:A:6:GLN:HG3	2.01	0.42
1:A:411:ASP:O	1:A:415:VAL:HG23	2.19	0.42
2:H:90:U:H2'	2:H:91:G:C8	2.54	0.42
2:H:169:U:H2'	2:H:170:U:H6	1.84	0.42
1:K:419:TYR:CD2	1:K:446:ARG:HB2	2.54	0.42
1:A:212:PRO:HB2	1:A:215:THR:HG21	2.01	0.42
1:C:145:VAL:HG11	1:C:304:LEU:CD2	2.49	0.42
1:C:375:LYS:NZ	2:D:80:G:O6	2.42	0.42
1:E:10:VAL:HG21	1:E:119:LEU:HB2	2.01	0.42
1:E:300:ASN:OD1	1:E:300:ASN:C	2.63	0.42
1:G:161:ARG:HG2	1:G:287:THR:HG22	2.01	0.42
1:G:308:LEU:HD22	1:G:313:VAL:HG22	2.01	0.42
1:G:378:ILE:HG21	1:G:471:LEU:HD22	2.00	0.42
1:K:6:GLN:O	1:K:10:VAL:HG23	2.19	0.42
1:K:10:VAL:HG12	1:K:115:ALA:HB1	2.01	0.42
1:A:167:PHE:HA	1:A:277:PRO:HA	2.02	0.42
1:A:300:ASN:OD1	1:A:300:ASN:O	2.37	0.42
2:D:26:G:H2'	2:D:27:C:H6	1.84	0.42
2:F:26:G:H2'	2:F:27:C:H6	1.85	0.42
2:F:107:C:H2'	2:F:108:C:C6	2.55	0.42
2:L:121:A:C6	2:L:122:C:N4	2.87	0.42
2:D:76:U:O2'	2:D:77:G:H8	2.02	0.42
2:D:107:C:H2'	2:D:108:C:C6	2.55	0.42
1:E:233:LEU:O	1:E:238:VAL:HG22	2.20	0.42
1:I:19:ILE:HG23	1:I:64:PHE:CE1	2.55	0.42
1:K:138:ALA:HB2	1:K:219:LEU:HB2	2.02	0.42
2:L:26:G:H2'	2:L:27:C:H6	1.84	0.42
1:A:40:LEU:O	1:A:43:SER:OG	2.29	0.42
1:A:323:ILE:HG12	1:A:394:ILE:HG12	2.02	0.42
2:J:121:A:C6	2:J:122:C:N4	2.88	0.42
1:K:51:ARG:O	1:K:52:PRO:C	2.62	0.42
1:K:137:TYR:CG	1:K:145:VAL:HG22	2.55	0.42
1:K:323:ILE:HG12	1:K:394:ILE:HG12	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:GLN:O	1:A:10:VAL:HG23	2.19	0.41
1:A:488:ILE:HG13	1:A:489:MET:N	2.35	0.41
1:E:132:PHE:CD1	1:E:223:ASN:HB3	2.55	0.41
1:E:161:ARG:HD3	1:I:261:PHE:CD1	2.54	0.41
1:G:76:SER:OG	2:H:10:U:O2	2.37	0.41
1:I:141:ASN:HB3	2:J:125:U:H1'	2.02	0.41
1:I:349:ILE:HG21	1:I:434:TYR:CD2	2.55	0.41
1:I:396:SER:O	1:I:489:MET:HE3	2.20	0.41
1:K:176:ILE:HG22	1:K:210:GLY:HA2	2.02	0.41
1:K:378:ILE:HG22	1:K:471:LEU:HD22	2.02	0.41
2:B:68:G:H8	1:C:182:ASN:HD22	1.68	0.41
1:C:9:ARG:HG3	1:C:9:ARG:HH11	1.85	0.41
2:D:31:C:O2	2:D:31:C:O4'	2.39	0.41
2:D:121:A:C6	2:D:122:C:N4	2.88	0.41
1:E:158:LYS:HB3	1:E:158:LYS:HE2	1.89	0.41
2:F:67:C:O2'	2:L:67:C:O3'	2.32	0.41
1:K:443:ILE:HD12	1:K:443:ILE:O	2.20	0.41
1:C:480:LEU:HD21	1:C:488:ILE:HD11	2.02	0.41
2:D:44:A:H2'	2:D:45:U:O4'	2.21	0.41
1:E:285:LEU:HB2	1:E:295:ILE:HD11	2.02	0.41
1:E:369:MET:HA	1:E:394:ILE:HG22	2.02	0.41
1:I:299:ASN:N	1:I:299:ASN:OD1	2.52	0.41
2:J:107:C:H2'	2:J:108:C:C6	2.55	0.41
2:B:169:U:H2'	2:B:170:U:H6	1.85	0.41
1:C:369:MET:HA	1:C:394:ILE:HG22	2.02	0.41
2:D:67:C:O2'	2:J:67:C:O3'	2.39	0.41
2:D:99:C:O2'	2:D:100:G:H5'	2.21	0.41
1:E:95:ASP:OD1	1:E:95:ASP:N	2.52	0.41
2:F:67:C:HO2'	2:L:67:C:C3'	2.33	0.41
2:H:134:U:O2'	2:H:135:U:C6	2.70	0.41
1:I:19:ILE:HG22	1:I:23:HIS:CD2	2.55	0.41
2:J:17:A:N1	2:J:143:U:O2'	2.50	0.41
2:L:96:A:C2	2:L:97:G:C8	3.08	0.41
2:B:26:G:H2'	2:B:27:C:C6	2.55	0.41
1:C:41:GLY:HA3	1:C:345:ALA:O	2.21	0.41
2:F:99:C:O2'	2:F:100:G:H5'	2.20	0.41
1:G:18:LEU:HD13	1:G:111:ILE:HD11	2.02	0.41
1:G:395:MET:HA	1:G:395:MET:HE2	2.03	0.41
1:I:138:ALA:HB2	1:I:219:LEU:HB2	2.02	0.41
1:I:230:ASP:OD1	1:I:230:ASP:N	2.54	0.41
2:J:3:U:HO2'	2:J:4:C:P	2.42	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:340:LYS:O	1:K:340:LYS:CG	2.69	0.41
2:B:134:U:O2'	2:B:135:U:C6	2.73	0.41
1:C:399:PRO:HA	1:C:476:LEU:HD21	2.02	0.41
1:E:323:ILE:HG12	1:E:394:ILE:HG12	2.02	0.41
2:F:53:C:HO2'	2:F:54:U:H6	1.68	0.41
2:F:121:A:C6	2:F:122:C:N4	2.89	0.41
2:B:161:A:H2'	2:B:162:C:C6	2.56	0.41
1:C:95:ASP:OD1	1:C:95:ASP:N	2.53	0.41
2:D:161:A:H2'	2:D:162:C:C6	2.56	0.41
1:E:172:PHE:HE1	1:E:225:THR:HG21	1.86	0.41
1:G:300:ASN:OD1	1:G:300:ASN:C	2.64	0.41
2:H:52:U:H2'	2:H:53:C:C6	2.56	0.41
2:L:17:A:N1	2:L:143:U:O2'	2.49	0.41
1:C:18:LEU:HD13	1:C:111:ILE:HD11	2.02	0.41
2:D:96:A:C2	2:D:97:G:C8	3.09	0.41
1:E:378:ILE:HG21	1:E:471:LEU:HG	2.01	0.41
2:H:99:C:O2'	2:H:100:G:H5'	2.20	0.41
1:I:167:PHE:HA	1:I:277:PRO:HA	2.02	0.41
1:I:184:GLN:HB2	1:I:224:LEU:HD11	2.03	0.41
1:I:233:LEU:HD22	1:I:238:VAL:HG21	2.01	0.41
2:J:96:A:C2	2:J:97:G:C8	3.09	0.41
1:K:23:HIS:HA	1:K:26:HIS:NE2	2.36	0.41
1:K:95:ASP:OD1	1:K:95:ASP:N	2.53	0.41
1:K:167:PHE:HA	1:K:277:PRO:HA	2.02	0.41
1:K:423:LYS:O	1:K:427:GLN:HG3	2.19	0.41
1:A:51:ARG:O	1:A:52:PRO:C	2.63	0.41
1:A:95:ASP:OD1	1:A:95:ASP:N	2.53	0.41
1:A:228:LYS:O	1:A:232:ASP:OD1	2.39	0.41
1:A:285:LEU:HB2	1:A:295:ILE:HD11	2.03	0.41
1:A:295:ILE:HB	1:I:294:GLU:HB3	2.03	0.41
1:A:420:ARG:NH2	2:B:81:G:O6	2.53	0.41
2:B:72:C:H2'	2:B:73:A:O4'	2.21	0.41
2:D:176:U:H5''	2:D:177:U:OP2	2.21	0.41
2:F:122:C:O2'	2:F:123:U:H5'	2.20	0.41
1:G:285:LEU:HB2	1:G:295:ILE:HD11	2.02	0.41
2:H:121:A:C6	2:H:122:C:N4	2.88	0.41
1:I:145:VAL:HG11	1:I:304:LEU:CD2	2.51	0.41
1:I:164:LEU:HD22	1:I:250:TRP:CB	2.50	0.41
1:I:193:SER:OG	1:I:196:GLU:HG3	2.21	0.41
2:J:31:C:O2	2:J:31:C:O4'	2.38	0.41
2:J:44:A:H2'	2:J:45:U:O4'	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:411:ASP:O	1:K:415:VAL:HG23	2.20	0.41
2:L:29:U:O2'	2:L:30:U:OP2	2.37	0.41
2:L:128:C:H2'	2:L:129:U:C6	2.56	0.41
2:B:52:U:H2'	2:B:53:C:C6	2.56	0.41
1:E:404:VAL:CG1	1:E:470:LEU:HD21	2.42	0.41
2:F:52:U:H2'	2:F:53:C:C6	2.56	0.41
2:F:169:U:H2'	2:F:170:U:H6	1.85	0.41
1:C:184:GLN:HB2	1:C:224:LEU:HD11	2.02	0.40
1:E:18:LEU:HD13	1:E:111:ILE:HD11	2.03	0.40
1:E:397:PHE:HE2	2:F:128:C:C4	2.39	0.40
1:G:95:ASP:OD1	1:G:95:ASP:N	2.54	0.40
1:G:354:ARG:HB3	1:G:354:ARG:CZ	2.52	0.40
1:G:356:LYS:O	1:G:360:ILE:HD12	2.21	0.40
2:H:74:A:HO2'	2:H:75:C:H6	1.67	0.40
1:I:165:ALA:HA	1:I:282:GLY:O	2.20	0.40
1:I:373:LEU:HD22	1:I:389:LEU:HD11	2.03	0.40
2:L:26:G:H2'	2:L:27:C:C6	2.56	0.40
2:L:72:C:H2'	2:L:73:A:O4'	2.21	0.40
2:L:123:U:C6	2:L:126:U:H5'	2.56	0.40
1:A:95:ASP:HB3	1:A:103:ILE:CG1	2.50	0.40
1:C:373:LEU:HD22	1:C:389:LEU:HD11	2.04	0.40
2:D:169:U:H2'	2:D:170:U:H6	1.85	0.40
1:E:10:VAL:HG11	1:E:115:ALA:O	2.22	0.40
1:G:397:PHE:HE2	2:H:128:C:C4	2.40	0.40
1:I:369:MET:HA	1:I:394:ILE:HG22	2.02	0.40
1:K:444:LEU:HD22	1:K:449:ILE:HA	2.02	0.40
1:C:224:LEU:O	1:C:227:TRP:HD1	2.05	0.40
1:C:300:ASN:OD1	1:C:300:ASN:O	2.40	0.40
2:D:53:C:HO2'	2:D:54:U:H6	1.69	0.40
2:H:44:A:H2'	2:H:45:U:O4'	2.21	0.40
2:H:122:C:O2'	2:H:123:U:H5'	2.21	0.40
1:I:87:LEU:HD12	1:I:88:PRO:HD2	2.04	0.40
1:I:164:LEU:HD12	1:I:286:LEU:HD13	2.03	0.40
1:K:95:ASP:HB3	1:K:103:ILE:CG1	2.51	0.40
1:A:41:GLY:HA3	1:A:345:ALA:O	2.21	0.40
1:A:207:ARG:HD3	1:A:212:PRO:HD3	2.03	0.40
2:B:67:C:C3'	2:H:67:C:HO2'	2.34	0.40
2:D:74:A:HO2'	2:D:75:C:H6	1.67	0.40
1:E:316:LYS:NZ	1:E:318:LYS:HB2	2.37	0.40
2:F:26:G:H2'	2:F:27:C:C6	2.56	0.40
1:G:27:ASN:OD1	2:H:22:U:H1'	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:132:PHE:CD1	1:G:223:ASN:HB3	2.56	0.40
1:I:51:ARG:O	1:I:52:PRO:C	2.64	0.40
2:J:176:U:H5''	2:J:177:U:OP2	2.21	0.40
1:K:1:MET:HE3	1:K:1:MET:HA	2.03	0.40
2:L:99:C:C2'	2:L:100:G:H5'	2.51	0.40
1:A:99:PRO:HG3	2:B:124:U:P	2.61	0.40
1:C:141:ASN:HB3	2:D:125:U:H1'	2.04	0.40
1:C:260:ALA:O	1:C:264:ILE:HD13	2.22	0.40
1:E:284:SER:OG	1:E:298:LYS:HE3	2.21	0.40
1:G:155:ASP:OD1	1:G:239:LYS:CE	2.69	0.40
1:G:323:ILE:HG12	1:G:394:ILE:HG12	2.02	0.40
2:H:26:G:H2'	2:H:27:C:C6	2.57	0.40
2:H:98:C:HO2'	2:H:99:C:H6	1.59	0.40
2:J:169:U:H2'	2:J:170:U:H6	1.86	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 PDB entries followed by that with respect to all EM entries.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	484/499 (97%)	477 (99%)	7 (1%)	0	100	100
1	C	484/499 (97%)	478 (99%)	6 (1%)	0	100	100
1	E	484/499 (97%)	478 (99%)	6 (1%)	0	100	100
1	G	484/499 (97%)	478 (99%)	6 (1%)	0	100	100
1	I	484/499 (97%)	477 (99%)	7 (1%)	0	100	100
1	K	484/499 (97%)	478 (99%)	6 (1%)	0	100	100
All	All	2904/2994 (97%)	2866 (99%)	38 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	445/456 (98%)	439 (99%)	6 (1%)	61	81
1	C	445/456 (98%)	437 (98%)	8 (2%)	51	75
1	E	445/456 (98%)	436 (98%)	9 (2%)	48	73
1	G	445/456 (98%)	434 (98%)	11 (2%)	42	68
1	I	445/456 (98%)	438 (98%)	7 (2%)	55	78
1	K	445/456 (98%)	435 (98%)	10 (2%)	45	70
All	All	2670/2736 (98%)	2619 (98%)	51 (2%)	49	74

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	103	ILE
1	A	160	GLU
1	A	276	ASN
1	A	289	LYS
1	A	294	GLU
1	A	458	ILE
1	C	49	ILE
1	C	103	ILE
1	C	105	LYS
1	C	160	GLU
1	C	276	ASN
1	C	289	LYS
1	C	354	ARG
1	C	388	ARG
1	E	16	LYS
1	E	103	ILE
1	E	160	GLU
1	E	176	ILE
1	E	289	LYS
1	E	301	LEU
1	E	311	GLU
1	E	312	ASN

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Mol	Chain	Res	Type
1	E	388	ARG
1	G	5	GLN
1	G	16	LYS
1	G	103	ILE
1	G	152	ILE
1	G	160	GLU
1	G	176	ILE
1	G	289	LYS
1	G	301	LEU
1	G	360	ILE
1	G	388	ARG
1	G	471	LEU
1	I	5	GLN
1	I	160	GLU
1	I	276	ASN
1	I	283	ILE
1	I	289	LYS
1	I	333	ARG
1	I	440	PHE
1	K	5	GLN
1	K	103	ILE
1	K	160	GLU
1	K	276	ASN
1	K	289	LYS
1	K	294	GLU
1	K	440	PHE
1	K	452	LYS
1	K	458	ILE
1	K	471	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	A	53	ASN
1	A	146	HIS
1	A	231	GLN
1	A	334	ASN
1	A	377	GLN
1	A	438	HIS
1	A	477	GLN
1	C	5	GLN

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Mol	Chain	Res	Type
1	C	27	ASN
1	C	110	GLN
1	C	146	HIS
1	C	182	ASN
1	C	186	ASN
1	C	213	GLN
1	C	334	ASN
1	C	347	GLN
1	C	377	GLN
1	C	481	GLN
1	E	146	HIS
1	E	252	GLN
1	G	5	GLN
1	G	146	HIS
1	G	299	ASN
1	G	312	ASN
1	G	477	GLN
1	I	27	ASN
1	I	146	HIS
1	I	186	ASN
1	I	334	ASN
1	I	481	GLN
1	K	6	GLN
1	K	27	ASN
1	K	141	ASN
1	K	146	HIS
1	K	334	ASN
1	K	377	GLN
1	K	438	HIS
1	K	477	GLN
1	K	481	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	176/177 (99%)	56 (31%)	2 (1%)
2	D	176/177 (99%)	56 (31%)	2 (1%)
2	F	176/177 (99%)	56 (31%)	2 (1%)
2	H	176/177 (99%)	56 (31%)	2 (1%)
2	J	176/177 (99%)	56 (31%)	2 (1%)
2	L	176/177 (99%)	56 (31%)	2 (1%)

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Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
All	All	1056/1062 (99%)	336 (31%)	12 (1%)

All (336) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	3	U
2	B	4	C
2	B	6	C
2	B	8	C
2	B	9	A
2	B	10	U
2	B	11	A
2	B	14	G
2	B	16	U
2	B	17	A
2	B	21	G
2	B	30	U
2	B	31	C
2	B	32	U
2	B	37	G
2	B	38	U
2	B	39	U
2	B	42	A
2	B	43	A
2	B	44	A
2	B	50	G
2	B	54	U
2	B	66	U
2	B	67	C
2	B	68	G
2	B	69	U
2	B	73	A
2	B	74	A
2	B	75	C
2	B	77	G
2	B	81	G
2	B	89	G
2	B	91	G
2	B	94	A
2	B	95	A
2	B	96	A
2	B	97	G

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Mol	Chain	Res	Type
2	B	99	C
2	B	101	U
2	B	103	U
2	B	110	A
2	B	111	G
2	B	112	U
2	B	121	A
2	B	124	U
2	B	125	U
2	B	126	U
2	B	127	A
2	B	130	U
2	B	135	U
2	B	136	U
2	B	139	A
2	B	173	U
2	B	175	G
2	B	176	U
2	B	177	U
2	D	3	U
2	D	4	C
2	D	8	C
2	D	9	A
2	D	10	U
2	D	11	A
2	D	14	G
2	D	16	U
2	D	17	A
2	D	21	G
2	D	30	U
2	D	31	C
2	D	32	U
2	D	37	G
2	D	38	U
2	D	39	U
2	D	42	A
2	D	43	A
2	D	44	A
2	D	50	G
2	D	54	U
2	D	66	U
2	D	67	C

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Mol	Chain	Res	Type
2	D	68	G
2	D	69	U
2	D	73	A
2	D	74	A
2	D	75	C
2	D	77	G
2	D	81	G
2	D	89	G
2	D	91	G
2	D	94	A
2	D	95	A
2	D	96	A
2	D	97	G
2	D	99	C
2	D	101	U
2	D	103	U
2	D	110	A
2	D	111	G
2	D	112	U
2	D	121	A
2	D	124	U
2	D	125	U
2	D	126	U
2	D	127	A
2	D	130	U
2	D	135	U
2	D	136	U
2	D	139	A
2	D	172	U
2	D	173	U
2	D	175	G
2	D	176	U
2	D	177	U
2	F	3	U
2	F	4	C
2	F	6	C
2	F	8	C
2	F	9	A
2	F	10	U
2	F	11	A
2	F	14	G
2	F	16	U

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Mol	Chain	Res	Type
2	F	17	A
2	F	21	G
2	F	30	U
2	F	31	C
2	F	32	U
2	F	37	G
2	F	38	U
2	F	39	U
2	F	42	A
2	F	43	A
2	F	44	A
2	F	50	G
2	F	54	U
2	F	66	U
2	F	67	C
2	F	68	G
2	F	69	U
2	F	73	A
2	F	74	A
2	F	75	C
2	F	77	G
2	F	81	G
2	F	89	G
2	F	91	G
2	F	94	A
2	F	95	A
2	F	96	A
2	F	97	G
2	F	99	C
2	F	101	U
2	F	103	U
2	F	110	A
2	F	111	G
2	F	112	U
2	F	121	A
2	F	124	U
2	F	125	U
2	F	126	U
2	F	127	A
2	F	130	U
2	F	135	U
2	F	136	U

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Mol	Chain	Res	Type
2	F	139	A
2	F	173	U
2	F	175	G
2	F	176	U
2	F	177	U
2	H	3	U
2	H	4	C
2	H	6	C
2	H	8	C
2	H	9	A
2	H	10	U
2	H	11	A
2	H	14	G
2	H	16	U
2	H	17	A
2	H	21	G
2	H	30	U
2	H	31	C
2	H	32	U
2	H	37	G
2	H	38	U
2	H	39	U
2	H	42	A
2	H	43	A
2	H	44	A
2	H	50	G
2	H	54	U
2	H	66	U
2	H	67	C
2	H	68	G
2	H	69	U
2	H	73	A
2	H	74	A
2	H	75	C
2	H	77	G
2	H	81	G
2	H	89	G
2	H	91	G
2	H	94	A
2	H	95	A
2	H	96	A
2	H	97	G

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Mol	Chain	Res	Type
2	H	99	C
2	H	101	U
2	H	103	U
2	H	110	A
2	H	111	G
2	H	112	U
2	H	121	A
2	H	124	U
2	H	125	U
2	H	126	U
2	H	127	A
2	H	130	U
2	H	135	U
2	H	136	U
2	H	139	A
2	H	173	U
2	H	175	G
2	H	176	U
2	H	177	U
2	J	3	U
2	J	4	C
2	J	8	C
2	J	9	A
2	J	10	U
2	J	11	A
2	J	14	G
2	J	16	U
2	J	17	A
2	J	21	G
2	J	30	U
2	J	31	C
2	J	32	U
2	J	37	G
2	J	38	U
2	J	39	U
2	J	42	A
2	J	43	A
2	J	44	A
2	J	50	G
2	J	54	U
2	J	66	U
2	J	67	C

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Mol	Chain	Res	Type
2	J	68	G
2	J	69	U
2	J	73	A
2	J	74	A
2	J	75	C
2	J	77	G
2	J	81	G
2	J	89	G
2	J	91	G
2	J	94	A
2	J	95	A
2	J	96	A
2	J	97	G
2	J	99	C
2	J	101	U
2	J	103	U
2	J	110	A
2	J	111	G
2	J	112	U
2	J	121	A
2	J	124	U
2	J	125	U
2	J	126	U
2	J	127	A
2	J	130	U
2	J	135	U
2	J	136	U
2	J	139	A
2	J	172	U
2	J	173	U
2	J	175	G
2	J	176	U
2	J	177	U
2	L	3	U
2	L	4	C
2	L	6	C
2	L	8	C
2	L	9	A
2	L	10	U
2	L	11	A
2	L	14	G
2	L	16	U

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Mol	Chain	Res	Type
2	L	17	A
2	L	21	G
2	L	30	U
2	L	31	C
2	L	32	U
2	L	37	G
2	L	38	U
2	L	39	U
2	L	42	A
2	L	43	A
2	L	44	A
2	L	50	G
2	L	54	U
2	L	66	U
2	L	67	C
2	L	68	G
2	L	69	U
2	L	73	A
2	L	74	A
2	L	75	C
2	L	77	G
2	L	81	G
2	L	89	G
2	L	91	G
2	L	94	A
2	L	95	A
2	L	96	A
2	L	97	G
2	L	99	C
2	L	101	U
2	L	103	U
2	L	110	A
2	L	111	G
2	L	112	U
2	L	121	A
2	L	124	U
2	L	125	U
2	L	126	U
2	L	127	A
2	L	130	U
2	L	135	U
2	L	136	U

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Mol	Chain	Res	Type
2	L	139	A
2	L	173	U
2	L	175	G
2	L	176	U
2	L	177	U

All (12) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	3	U
2	B	43	A
2	D	3	U
2	D	43	A
2	F	3	U
2	F	43	A
2	H	3	U
2	H	43	A
2	J	3	U
2	J	43	A
2	L	3	U
2	L	43	A

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Map visualisation

This section contains visualisations of the EMDB entry EMD-60725. These allow visual inspection of the internal detail of the map and identification of artifacts.

Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

This section was not generated.

6.2 Central slices

This section was not generated.

6.3 Largest variance slices

This section was not generated.

6.4 Orthogonal standard-deviation projections (False-color)

This section was not generated.

6.5 Orthogonal surface views

This section was not generated.

6.6 Mask visualisation

This section was not generated. No masks/segmentation were deposited.

7 Map analysis ⓘ

This section contains the results of statistical analysis of the map.

7.1 Map-value distribution ⓘ

This section was not generated.

7.2 Volume estimate versus contour level ⓘ

This section was not generated.

7.3 Rotationally averaged power spectrum ⓘ

This section was not generated. The rotationally averaged power spectrum had issues being displayed.

8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit ⓘ

This section was not generated.