



Full wwPDB EM Validation Report ⓘ

Sep 16, 2026 – 11:18 PM EDT

PDB ID : 8FZI / pdb_00008fzi
EMDB ID : EMD-29631
Title : Cryo-EM structure of an E. coli rotated ribosome bound with RF3-GDPCP and p/E-tRNAPhe (Composite state II-B)
Authors : Rybak, M.Y.; Li, L.; Lin, J.; Gagnon, M.G.
Deposited on : 2023-01-28
Resolution : 3.10 Å(reported)
Based on initial model : 7K00

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.dev133
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

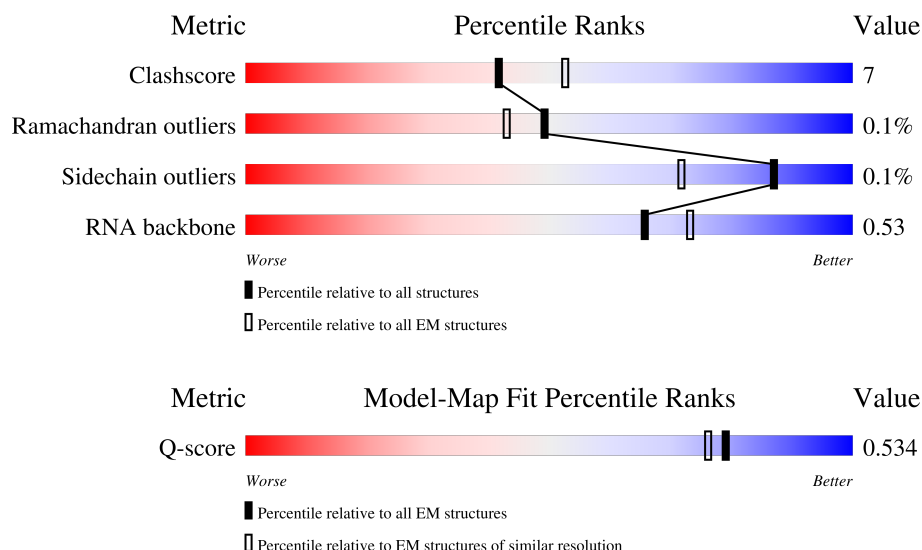
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.10 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)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	14724 (2.60 - 3.60)

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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	a	1542	
2	b	241	
3	c	233	

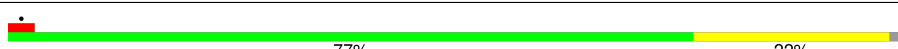
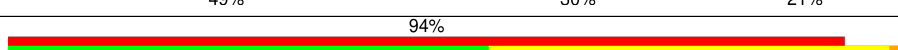
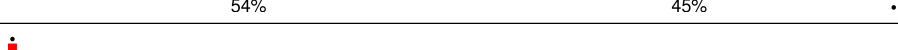
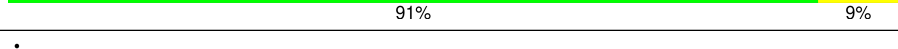
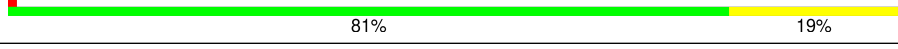
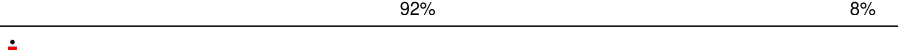
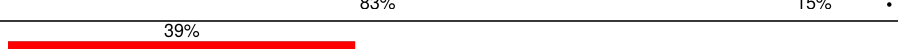
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Mol	Chain	Length	Quality of chain
4	d	206	
5	e	167	
6	f	131	
7	g	156	
8	h	130	
9	i	130	
10	j	103	
11	k	129	
12	l	124	
13	m	118	
14	n	101	
15	o	89	
16	p	82	
17	q	84	
18	r	75	
19	s	92	
20	t	87	
21	u	71	
22	x	76	
23	z	21	
24	v	529	
25	A	2904	
26	B	120	
27	C	273	
28	D	209	

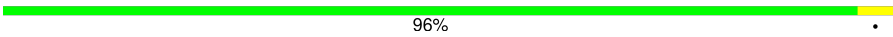
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Mol	Chain	Length	Quality of chain
29	E	201	
30	F	179	
31	G	177	
32	I	165	
33	J	142	
34	L	142	
35	M	123	
36	N	144	
37	O	136	
38	P	127	
39	Q	117	
40	R	115	
41	S	118	
42	T	103	
43	U	110	
44	V	100	
45	W	104	
46	X	94	
47	Y	85	
48	Z	78	
49	1	63	
50	2	59	
51	3	70	
52	4	57	
53	5	55	

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Mol	Chain	Length	Quality of chain
54	6	46	 96% .
55	7	65	 89% 9% .
56	8	38	 63% 37%

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
58	GCP	v	601	-	-	X	-

2 Entry composition

There are 60 unique types of molecules in this entry. The entry contains 149604 atoms, of which 0 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 RNA chain called 16S Ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	a	1528	Total	C	N	O	P	0	0
			32803	14637	6019	10619	1528		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
2	b	224	Total	C	N	O	S	0	0
			1754	1110	315	321	8		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
3	c	206	Total	C	N	O	S	0	0
			1624	1028	305	288	3		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
4	d	205	Total	C	N	O	S	0	0
			1643	1026	315	298	4		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
5	e	156	Total	C	N	O	S	0	0
			1152	717	217	212	6		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
6	f	103	Total	C	N	O	S	0	0
			839	530	151	151	7		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
7	g	152	Total	C	N	O	S	0	0
			1191	741	230	216	4		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
8	h	129	Total	C	N	O	S	0	0
			979	616	173	184	6		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
9	i	127	Total	C	N	O	S	0	0
			1022	634	206	179	3		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
10	j	98	Total	C	N	O	S	0	0
			786	493	150	142	1		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
11	k	117	Total	C	N	O	S	0	0
			877	540	173	161	3		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
k	119	IAS	ASN	conflict	UNP A0A0H3PWX2

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

Mol	Chain	Residues	Atoms					AltConf	Trace
12	l	121	Total	C	N	O	S	0	0
			942	582	193	162	5		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
13	m	115	Total	C	N	O	S	0	0
			891	552	179	157	3		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
14	n	100	Total	C	N	O	S	0	0
			805	499	164	139	3		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
15	o	88	Total	C	N	O	S	0	0
			714	439	144	130	1		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
16	p	81	Total	C	N	O	S	0	0
			643	403	127	112	1		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
17	q	79	Total	C	N	O	S	0	0
			641	406	120	112	3		

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

Mol	Chain	Residues	Atoms				AltConf	Trace
18	r	54	Total	C	N	O	0	0
			446	283	85	78		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
19	s	83	Total	C	N	O	S	0	0
			663	424	126	111	2		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
20	t	86	Total	C	N	O	S	0	0
			670	414	138	115	3		

- Molecule 21 is a protein called 30S ribosomal protein S21.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	u	55	Total	C	N	O	S	0	0
			460	287	95	77	1		

- Molecule 22 is a RNA chain called p/E phenylalanyl-tRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace	
22	x	74	Total	C	N	O	P	S	0	0
			1588	713	285	515	73	2		

- Molecule 23 is a RNA chain called F-UAA mRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	z	8	Total	C	N	O	P	0	0
			168	76	29	55	8		

- Molecule 24 is a protein called Peptide chain release factor RF3.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	v	525	Total	C	N	O	S	0	0
			4099	2595	709	773	22		

- Molecule 25 is a RNA chain called 23S Ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	A	2899	Total	C	N	O	P	0	0
			62252	27778	11456	20119	2899		

- Molecule 26 is a RNA chain called 5S Ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	B	120	Total	C	N	O	P	0	0
			2572	1145	470	837	120		

- Molecule 27 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	C	271	Total	C	N	O	S	0	0
			2082	1288	423	364	7		

- Molecule 28 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	D	209	Total	C	N	O	S	0	0
			1566	980	288	294	4		

- Molecule 29 is a protein called 50S ribosomal protein L4.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	E	201	Total	C	N	O	S	0	0
			1552	974	283	290	5		

- Molecule 30 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	F	177	Total	C	N	O	S	0	0
			1410	899	249	256	6		

- Molecule 31 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	G	176	Total	C	N	O	S	0	0
			1323	832	243	246	2		

- Molecule 32 is a protein called 50S ribosomal protein L10.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	I	131	Total	C	N	O	S	0	0
			988	625	175	183	5		

- Molecule 33 is a protein called 50S ribosomal protein L11.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	J	141	Total	C	N	O	S	0	0
			1032	651	179	196	6		

- Molecule 34 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	L	142	Total	C	N	O	S	0	0
			1129	714	212	199	4		

- Molecule 35 is a protein called 50S ribosomal protein L14.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	M	123	Total	C	N	O	S	0	0
			946	593	181	166	6		

- Molecule 36 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	N	144	Total	C	N	O	S	0	0
			1052	653	207	190	2		

- Molecule 37 is a protein called 50S ribosomal protein L16.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	O	136	Total	C	N	O	S	0	0
			1075	686	205	177	7		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
O	82	MS6	MET	conflict	UNP E6BI61

- Molecule 38 is a protein called 50S ribosomal protein L17.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	P	118	Total	C	N	O	S	0	0
			945	585	194	161	5		

- Molecule 39 is a protein called 50S ribosomal protein L18.

Mol	Chain	Residues	Atoms					AltConf	Trace
39	Q	116	Total	C	N	O		0	0
			892	552	178	162			

- Molecule 40 is a protein called 50S ribosomal protein L19.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	R	114	Total	C	N	O	S	0	0
			917	574	179	163	1		

- Molecule 41 is a protein called 50S ribosomal protein L20.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	S	117	Total	C	N	O		0	0
			947	604	192	151			

- Molecule 42 is a protein called Ribosomal protein L21.

Mol	Chain	Residues	Atoms					AltConf	Trace
42	T	103	Total	C	N	O	S	0	0
			816	516	153	145	2		

- Molecule 43 is a protein called 50S ribosomal protein L22.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	U	110	Total	C	N	O	S	0	0
			857	532	166	156	3		

- Molecule 44 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	V	93	Total	C	N	O	S	0	0
			738	466	139	131	2		

- Molecule 45 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms					AltConf	Trace
45	W	102	Total	C	N	O		0	0
			779	492	146	141			

- Molecule 46 is a protein called 50S ribosomal protein L25.

Mol	Chain	Residues	Atoms					AltConf	Trace
46	X	94	Total	C	N	O	S	0	0
			753	479	137	134	3		

- Molecule 47 is a protein called 50S ribosomal protein L27.

Mol	Chain	Residues	Atoms					AltConf	Trace
47	Y	76	Total	C	N	O	S	0	0
			582	360	117	104	1		

- Molecule 48 is a protein called 50S ribosomal protein L28.

Mol	Chain	Residues	Atoms					AltConf	Trace
48	Z	77	Total	C	N	O	S	0	0
			625	388	129	106	2		

- Molecule 49 is a protein called 50S ribosomal protein L29.

Mol	Chain	Residues	Atoms					AltConf	Trace
49	1	61	Total	C	N	O	S	0	0
			495	305	97	92	1		

- Molecule 50 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms					AltConf	Trace
50	2	58	Total	C	N	O	S	0	0
			449	281	87	79	2		

- Molecule 51 is a protein called 50S ribosomal protein L31.

Mol	Chain	Residues	Atoms					AltConf	Trace
51	3	48	Total	C	N	O	S	0	0
			373	232	66	69	6		

- Molecule 52 is a protein called 50S ribosomal protein L32.

Mol	Chain	Residues	Atoms					AltConf	Trace
52	4	55	Total	C	N	O	S	0	0
			434	263	92	78	1		

- Molecule 53 is a protein called 50S ribosomal protein L33.

Mol	Chain	Residues	Atoms				AltConf	Trace
53	5	51	Total	C	N	O	0	0
			417	269	76	72		

- Molecule 54 is a protein called 50S ribosomal protein L34.

Mol	Chain	Residues	Atoms					AltConf	Trace
54	6	46	Total	C	N	O	S	0	0
			377	228	90	57	2		

- Molecule 55 is a protein called 50S ribosomal protein L35.

Mol	Chain	Residues	Atoms					AltConf	Trace
55	7	64	Total	C	N	O	S	0	0
			504	323	105	74	2		

- Molecule 56 is a protein called 50S ribosomal protein L36.

Mol	Chain	Residues	Atoms					AltConf	Trace
56	8	38	Total	C	N	O	S	0	0
			302	185	65	48	4		

- Molecule 57 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

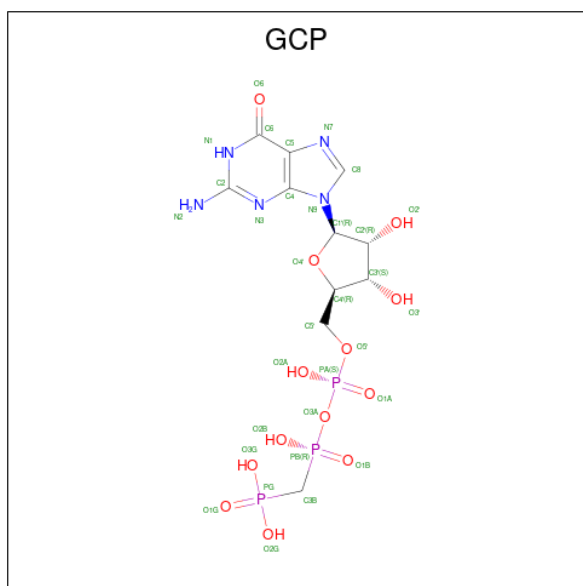
Mol	Chain	Residues	Atoms		AltConf
57	a	125	Total	Mg	0
			125	125	
57	n	1	Total	Mg	0
			1	1	
57	p	1	Total	Mg	0
			1	1	
57	x	1	Total	Mg	0
			1	1	
57	z	1	Total	Mg	0
			1	1	
57	A	572	Total	Mg	0
			572	572	
57	B	15	Total	Mg	0
			15	15	
57	C	3	Total	Mg	0
			3	3	
57	D	4	Total	Mg	0
			4	4	
57	E	1	Total	Mg	0
			1	1	
57	N	1	Total	Mg	0
			1	1	
57	P	2	Total	Mg	0
			2	2	

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Mol	Chain	Residues	Atoms		AltConf
57	S	1	Total	Mg	0
			1	1	
57	T	3	Total	Mg	0
			3	3	
57	V	2	Total	Mg	0
			2	2	
57	Y	1	Total	Mg	0
			1	1	
57	4	1	Total	Mg	0
			1	1	
57	6	1	Total	Mg	0
			1	1	
57	7	1	Total	Mg	0
			1	1	

- Molecule 58 is PHOSPHOMETHYLPHOSPHONIC ACID GUANYLATE ESTER (CCD ID: GCP) (formula: $C_{11}H_{18}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms					AltConf
58	v	1	Total	C	N	O	P	0
			32	11	5	13	3	

- Molecule 59 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
59	3	1	Total	Zn	0
			1	1	

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Mol	Chain	Residues	Atoms		AltConf
59	8	1	Total	Zn	0
			1	1	

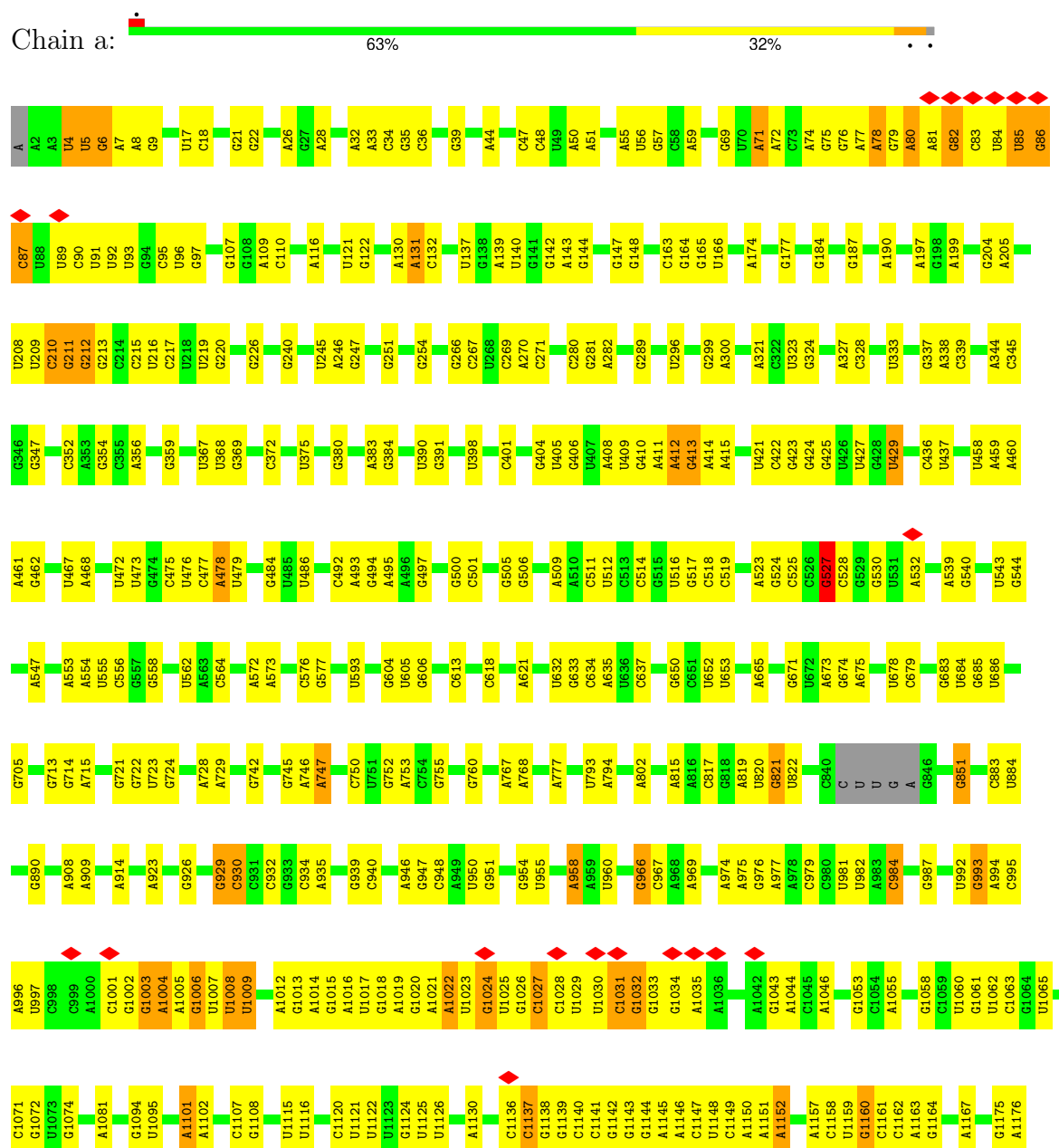
- Molecule 60 is water.

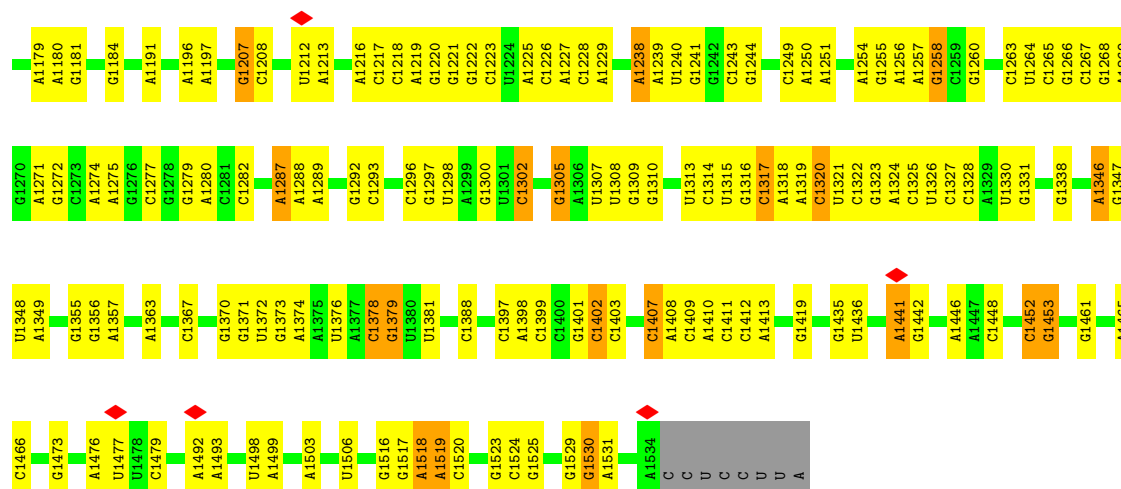
Mol	Chain	Residues	Atoms		AltConf
60	a	29	Total	O	0
			29	29	
60	n	1	Total	O	0
			1	1	
60	x	1	Total	O	0
			1	1	
60	z	1	Total	O	0
			1	1	
60	A	202	Total	O	0
			202	202	
60	B	7	Total	O	0
			7	7	
60	C	1	Total	O	0
			1	1	
60	D	3	Total	O	0
			3	3	
60	N	1	Total	O	0
			1	1	
60	P	1	Total	O	0
			1	1	
60	S	2	Total	O	0
			2	2	
60	6	1	Total	O	0
			1	1	
60	7	1	Total	O	0
			1	1	
60	8	1	Total	O	0
			1	1	

3 Residue-property plots

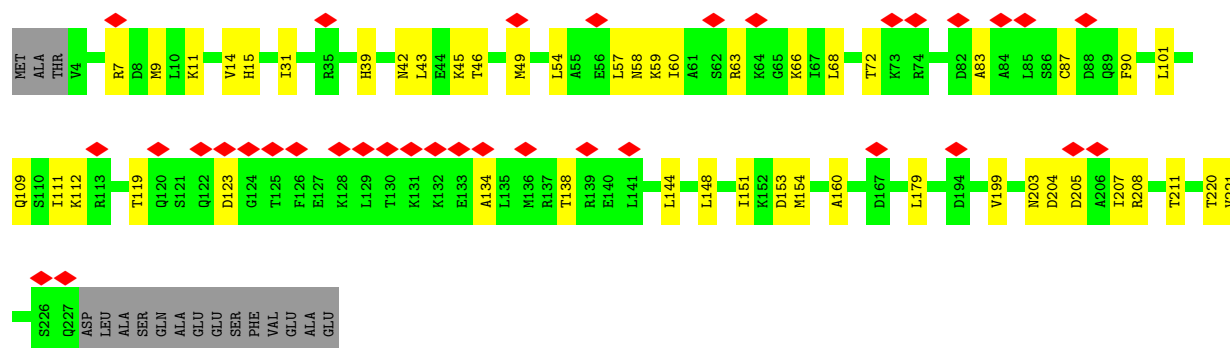
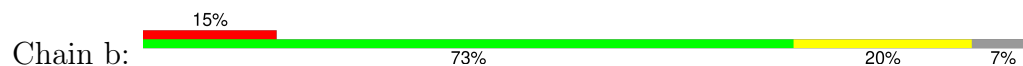
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map 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 diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). 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 Ribosomal RNA

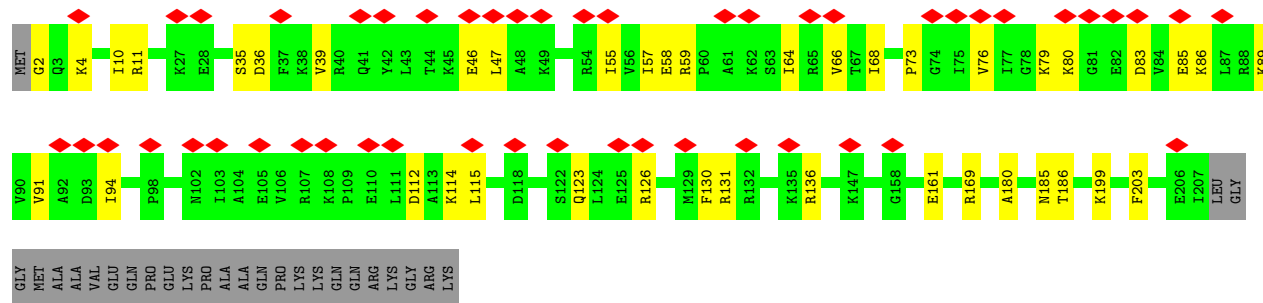




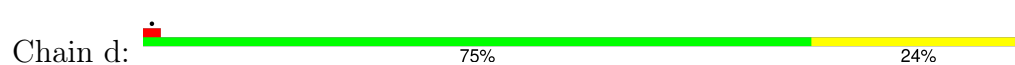
• Molecule 2: 30S ribosomal protein S2



• Molecule 3: 30S ribosomal protein S3



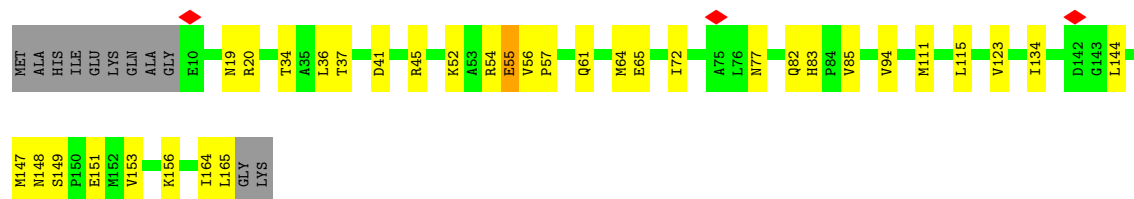
• Molecule 4: 30S ribosomal protein S4





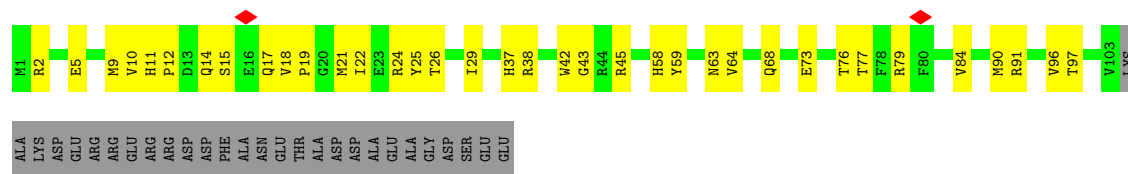
• Molecule 5: 30S ribosomal protein S5

Chain e: 73% 20% 7%



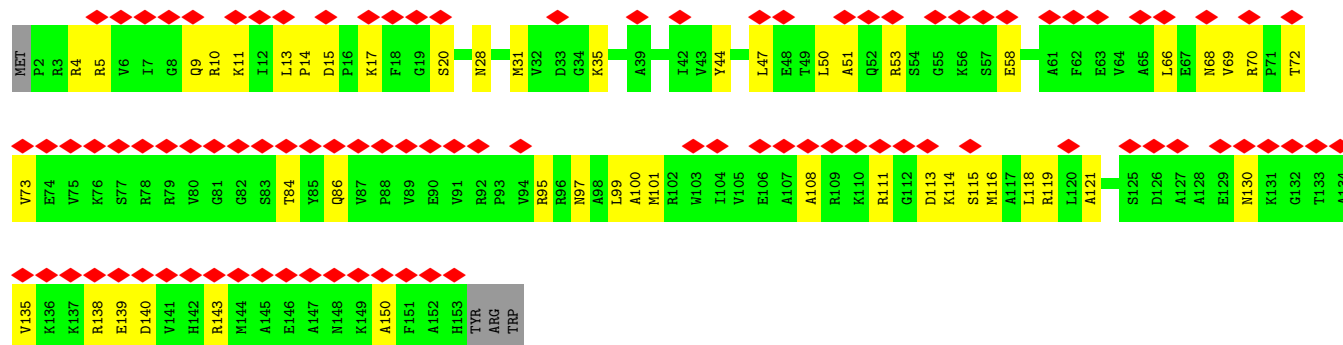
• Molecule 6: 30S ribosomal protein S6

Chain f: 51% 27% 21%



• Molecule 7: 30S ribosomal protein S7

Chain g: 60% 67% 31%



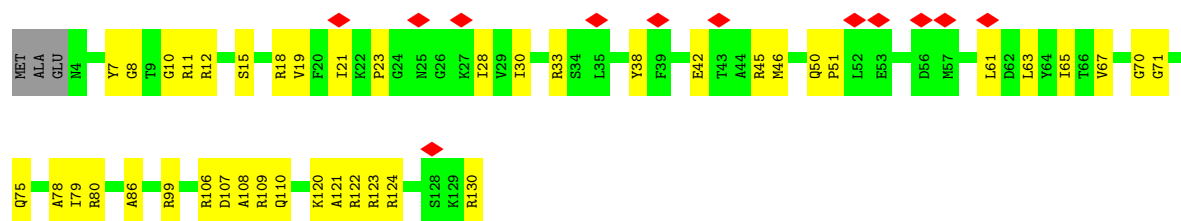
• Molecule 8: 30S ribosomal protein S8

Chain h: 82% 18%

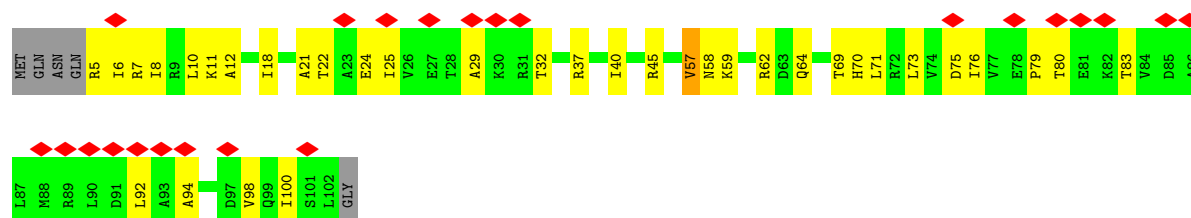


• Molecule 9: 30S ribosomal protein S9

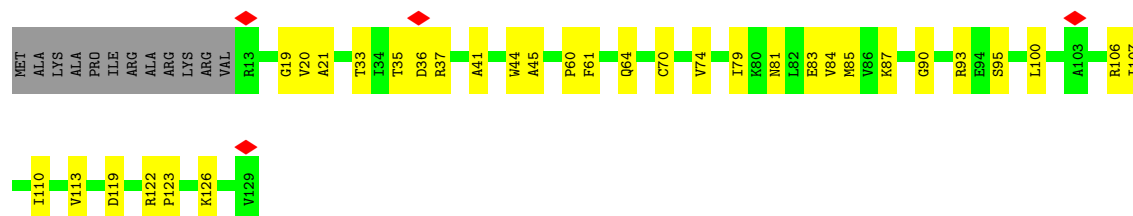
Chain i: 9% 65% 32%



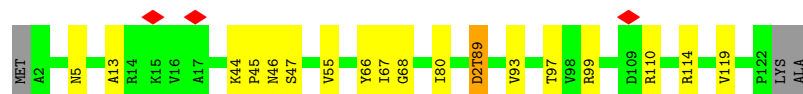
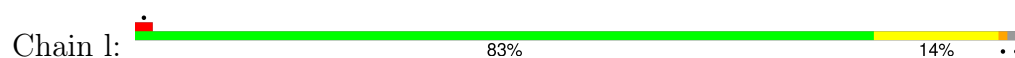
- Molecule 10: 30S ribosomal protein S10



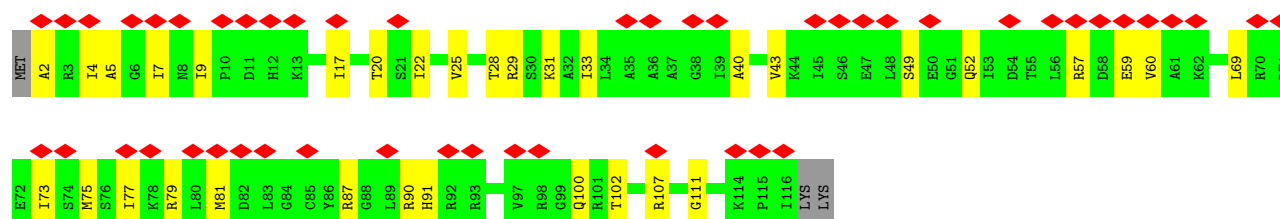
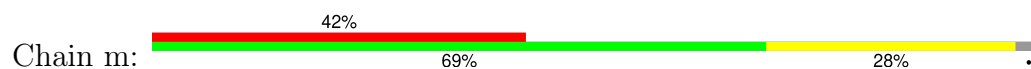
- Molecule 11: 30S ribosomal protein S11



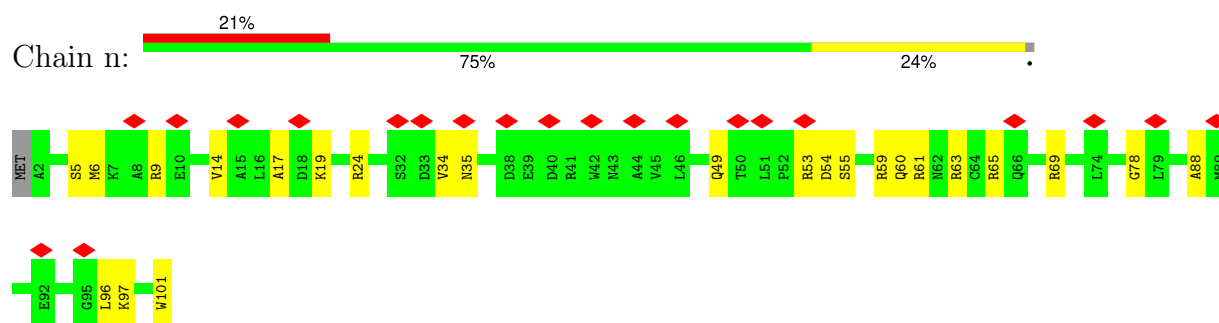
- Molecule 12: 30S ribosomal protein S12



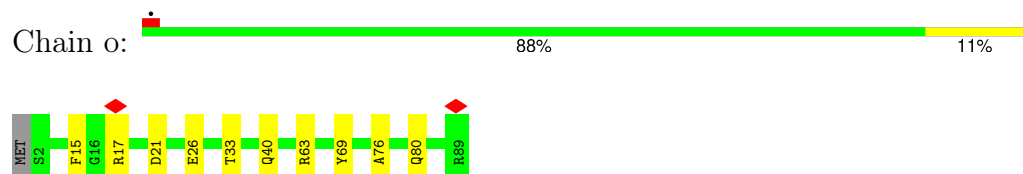
- Molecule 13: 30S ribosomal protein S13



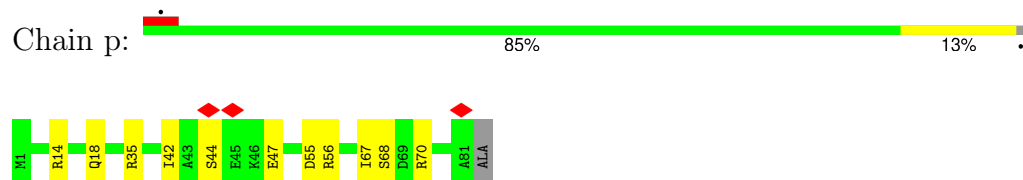
- Molecule 14: 30S ribosomal protein S14



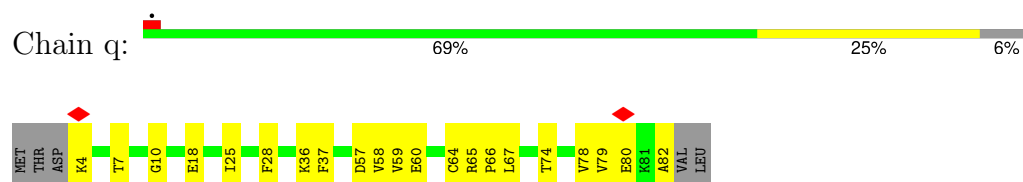
- Molecule 15: 30S ribosomal protein S15



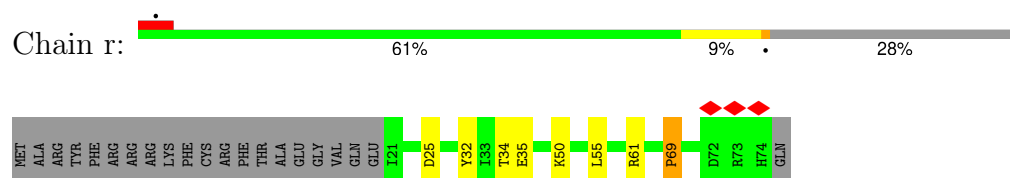
- Molecule 16: 30S ribosomal protein S16



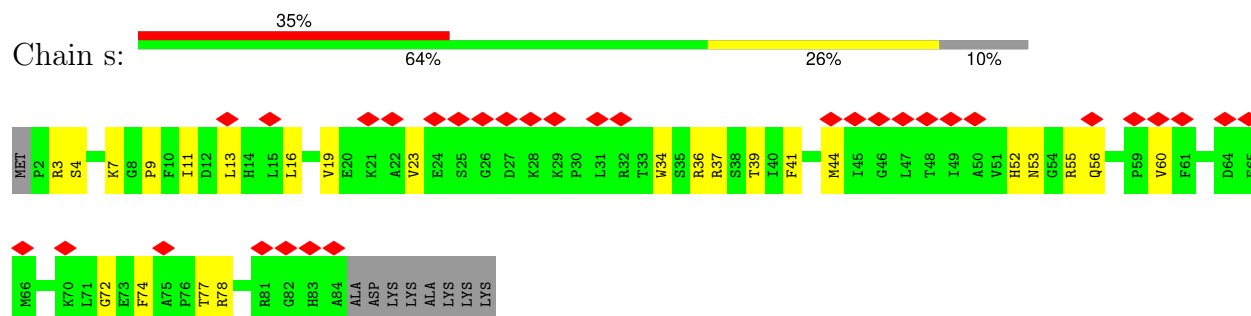
- Molecule 17: 30S ribosomal protein S17



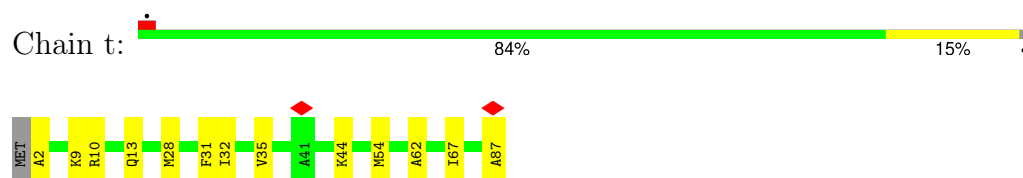
- Molecule 18: 30S ribosomal protein S18



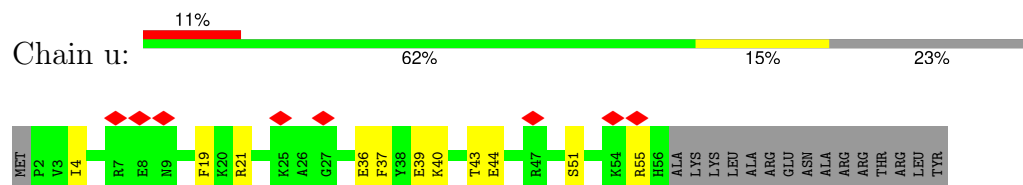
- Molecule 19: 30S ribosomal protein S19



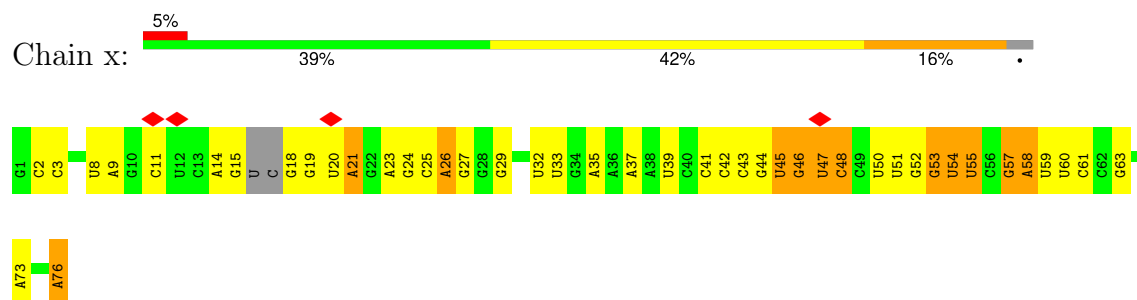
- Molecule 20: 30S ribosomal protein S20



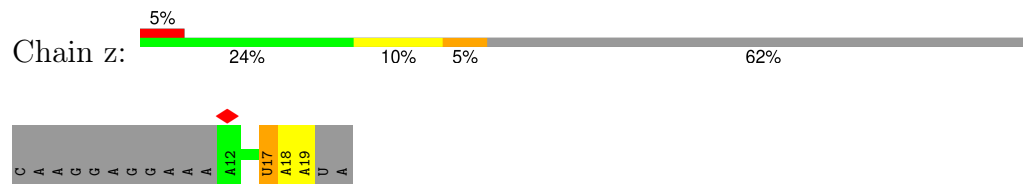
- Molecule 21: 30S ribosomal protein S21



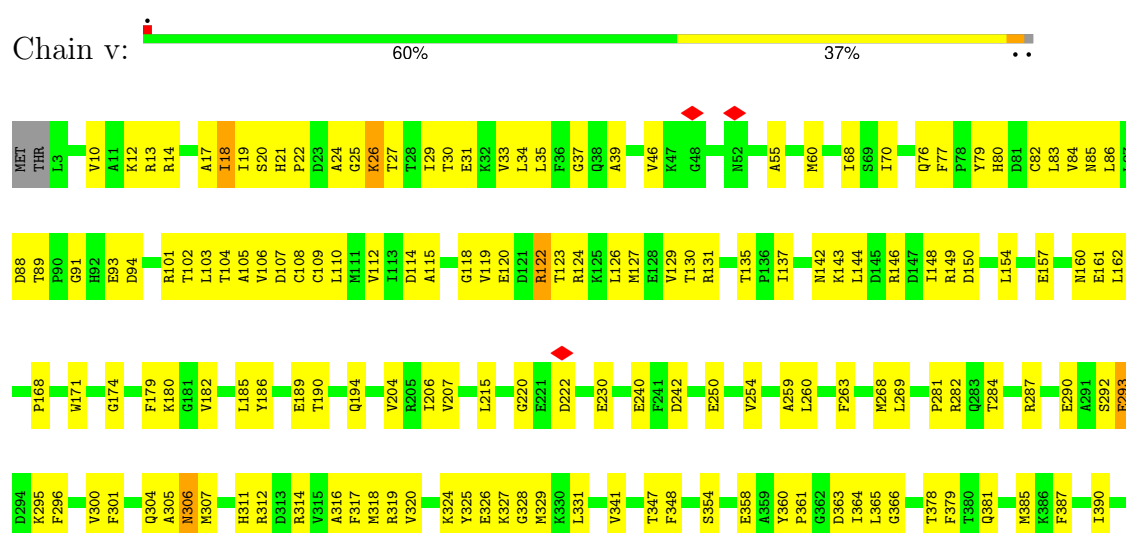
- Molecule 22: p/E phenylalanyl-tRNA

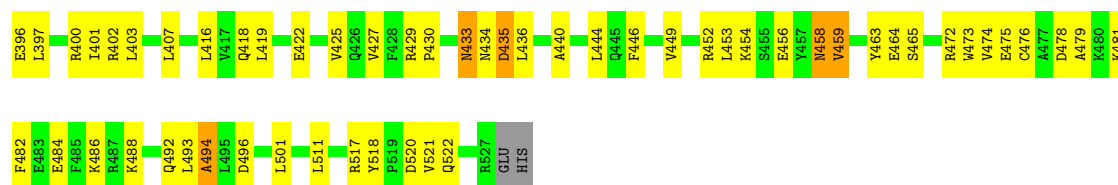


- Molecule 23: F-UAA mRNA



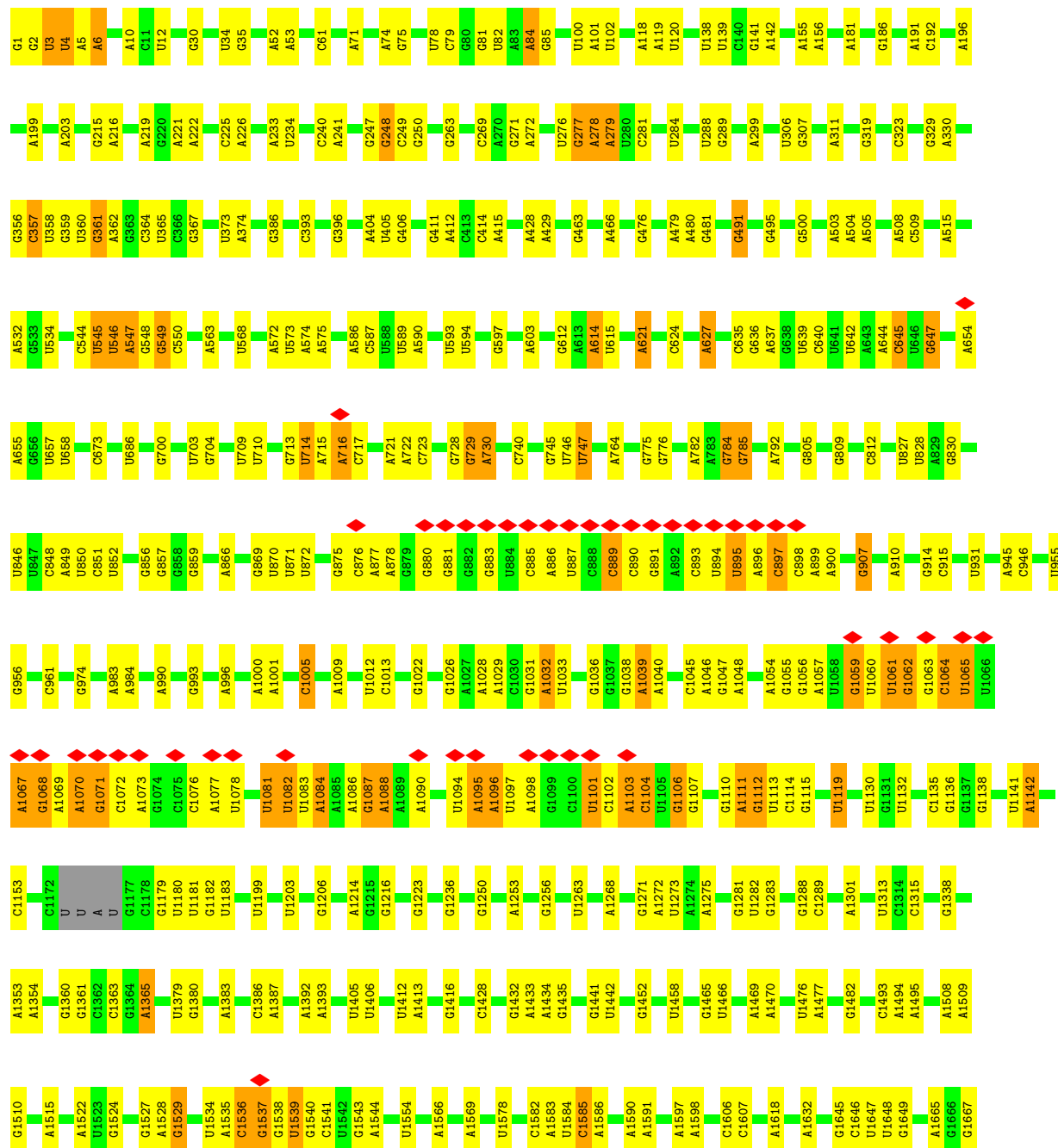
- Molecule 24: Peptide chain release factor RF3

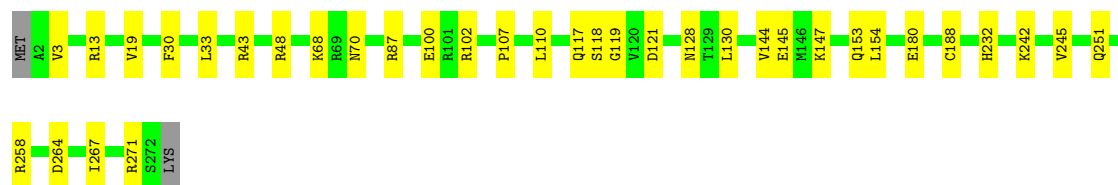




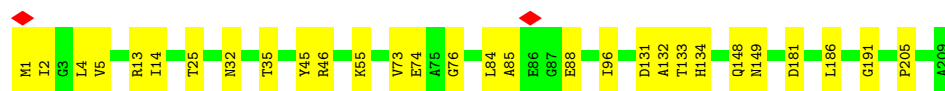
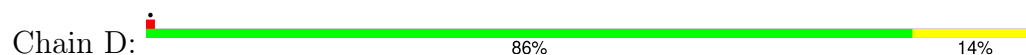
• Molecule 25: 23S Ribosomal RNA

Chain A: 72% 24%

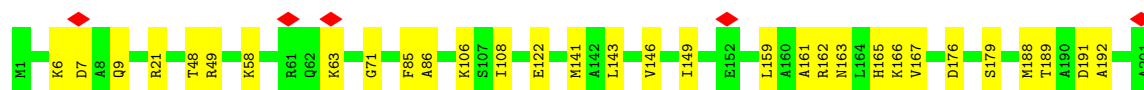
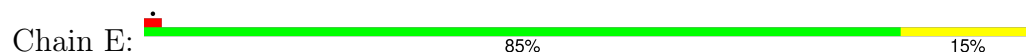




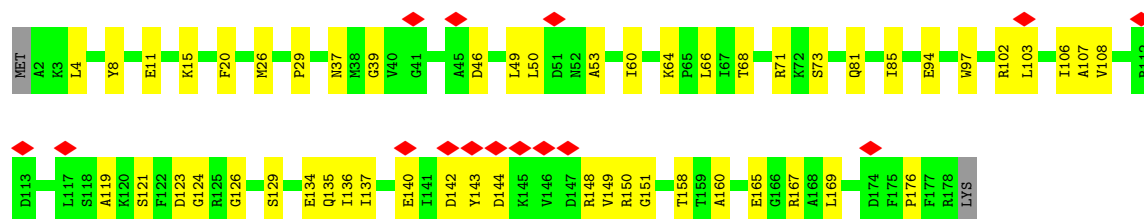
- Molecule 28: 50S ribosomal protein L3



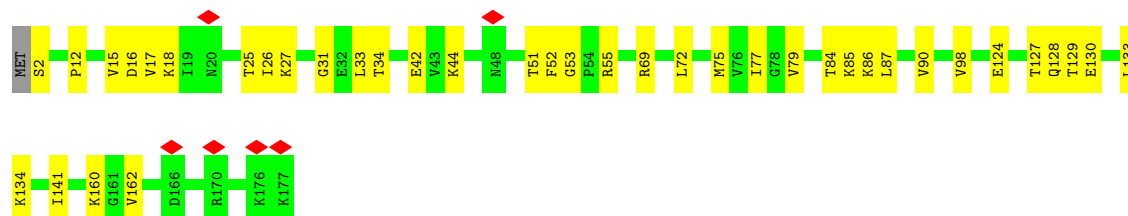
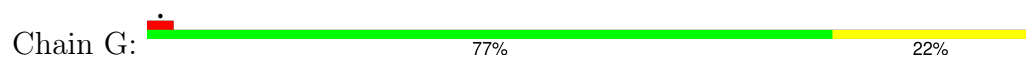
- Molecule 29: 50S ribosomal protein L4



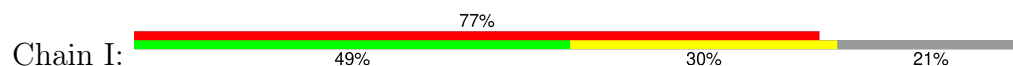
- Molecule 30: 50S ribosomal protein L5

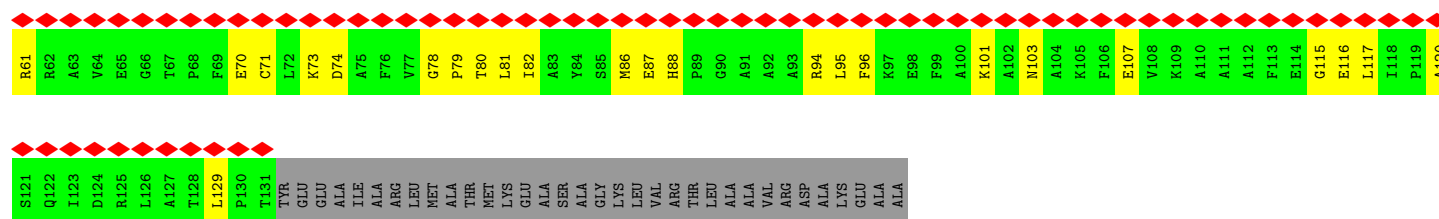


- Molecule 31: 50S ribosomal protein L6

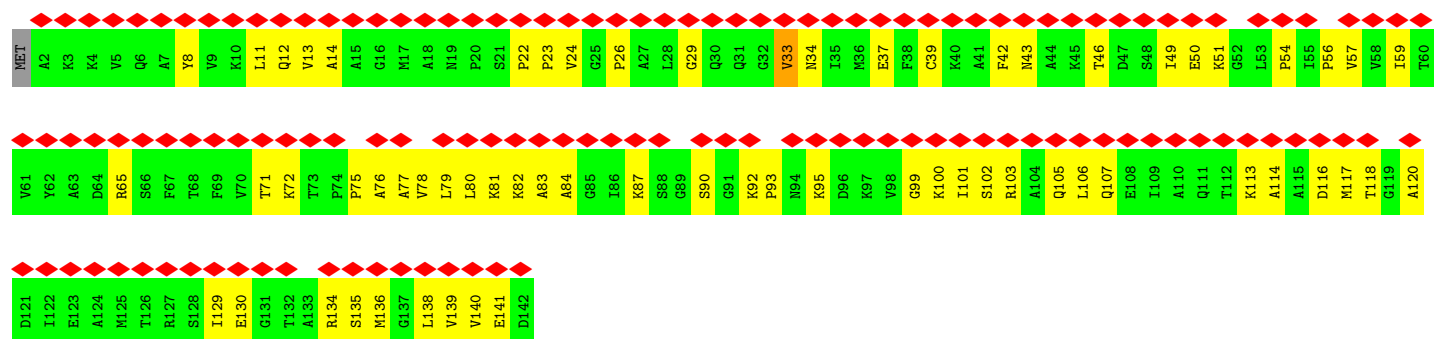


- Molecule 32: 50S ribosomal protein L10





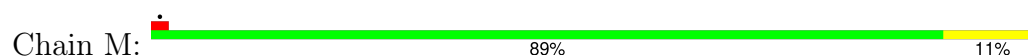
• Molecule 33: 50S ribosomal protein L11



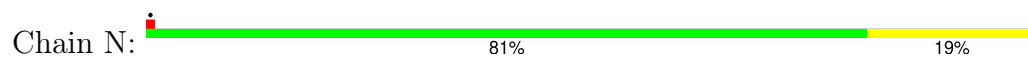
• Molecule 34: 50S ribosomal protein L13



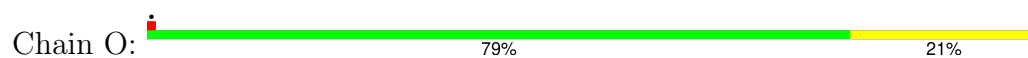
• Molecule 35: 50S ribosomal protein L14

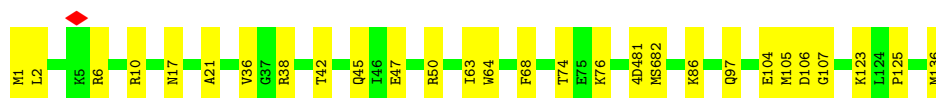


• Molecule 36: 50S ribosomal protein L15



• Molecule 37: 50S ribosomal protein L16





- Molecule 38: 50S ribosomal protein L17

Chain P: 86% 7% 7%



- Molecule 39: 50S ribosomal protein L18

Chain Q: 80% 19% .



- Molecule 40: 50S ribosomal protein L19

Chain R: 85% 14% .



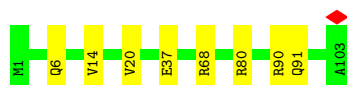
- Molecule 41: 50S ribosomal protein L20

Chain S: 91% 8% .



- Molecule 42: Ribosomal protein L21

Chain T: 92% 8% .

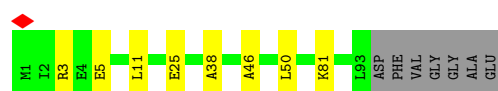
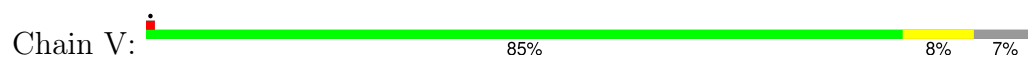


- Molecule 43: 50S ribosomal protein L22

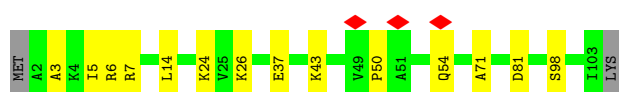
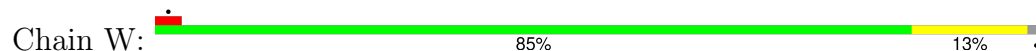
Chain U: 89% 11% .



- Molecule 44: 50S ribosomal protein L23



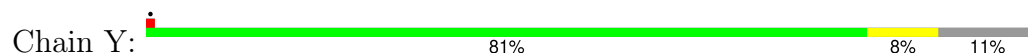
- Molecule 45: 50S ribosomal protein L24



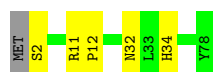
- Molecule 46: 50S ribosomal protein L25



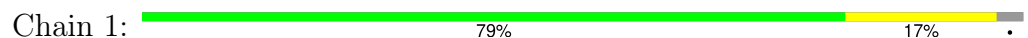
- Molecule 47: 50S ribosomal protein L27



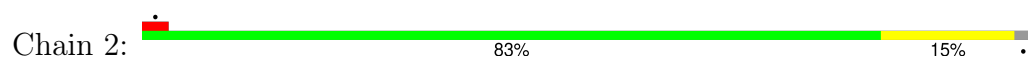
- Molecule 48: 50S ribosomal protein L28



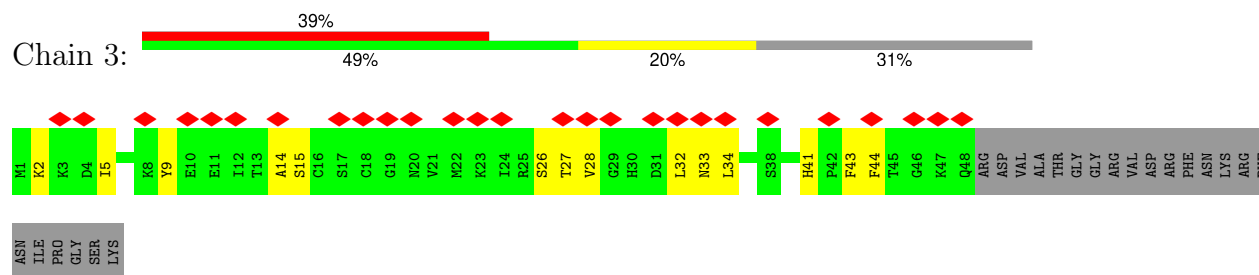
- Molecule 49: 50S ribosomal protein L29



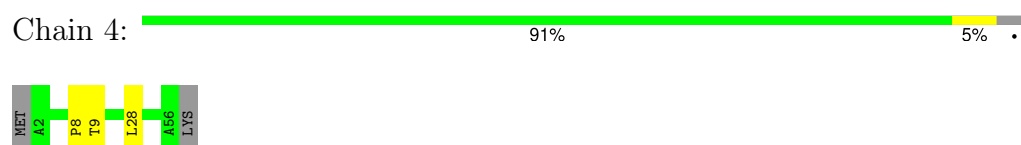
- Molecule 50: 50S ribosomal protein L30



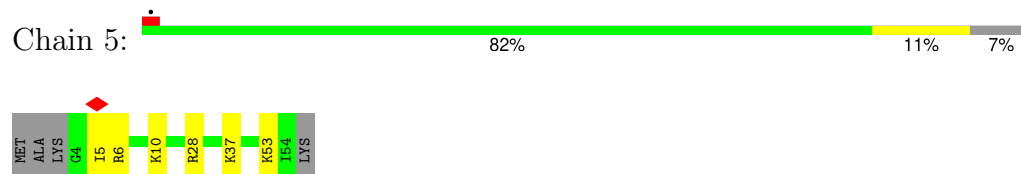
• Molecule 51: 50S ribosomal protein L31



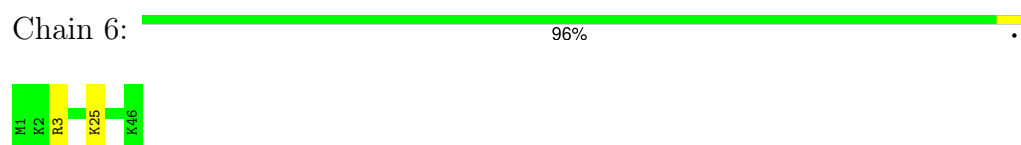
• Molecule 52: 50S ribosomal protein L32



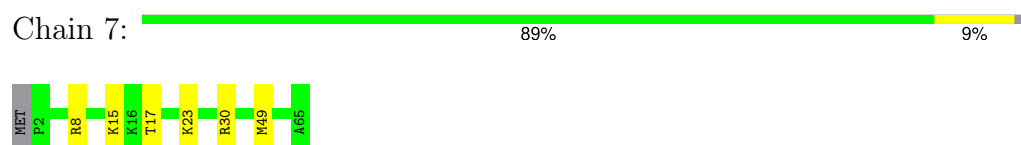
• Molecule 53: 50S ribosomal protein L33



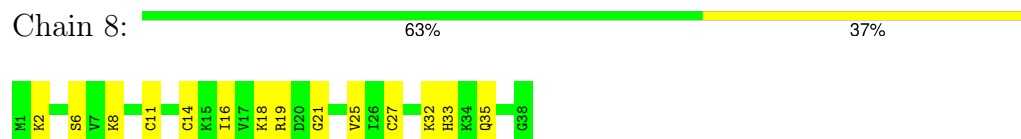
• Molecule 54: 50S ribosomal protein L34



• Molecule 55: 50S ribosomal protein L35



• Molecule 56: 50S ribosomal protein L36



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	56202	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	40.44	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	96000	Depositor
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	42.411	Depositor
Minimum map value	-25.024	Depositor
Average map value	0.001	Depositor
Map value standard deviation	1.033	Depositor
Recommended contour level	3	Depositor
Map size ($\text{\AA}$)	435.2, 435.2, 435.2	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.85, 0.85, 0.85	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MA6, 3TD, IAS, PSU, 2MG, OMC, OMU, MG, 2MA, 1MG, H2U, GCP, 6MZ, D2T, MIA, G7M, MEQ, 5MU, MS6, 4SU, 4OC, UR3, 5MC, OMG, 4D4, ZN

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	0.23	0/36450	0.30	0/56856
2	b	0.17	0/1785	0.39	0/2404
3	c	0.16	0/1651	0.35	0/2225
4	d	0.22	0/1665	0.36	0/2227
5	e	0.22	0/1165	0.37	0/1568
6	f	0.19	0/858	0.44	0/1160
7	g	0.15	0/1206	0.38	0/1617
8	h	0.23	0/989	0.35	0/1326
9	i	0.15	0/1034	0.33	0/1375
10	j	0.17	0/796	0.42	0/1077
11	k	0.21	0/884	0.36	0/1191
12	l	0.24	0/945	0.39	0/1268
13	m	0.14	0/900	0.32	0/1204
14	n	0.12	0/817	0.26	0/1088
15	o	0.25	0/722	0.39	0/964
16	p	0.22	0/653	0.38	0/877
17	q	0.23	0/650	0.41	0/871
18	r	0.36	0/453	0.67	2/609 (0.3%)
19	s	0.13	0/680	0.31	0/915
20	t	0.22	0/676	0.38	0/895
21	u	0.15	0/467	0.27	0/620
22	x	0.25	0/1602	0.31	0/2493
23	z	0.17	0/187	0.29	0/288
24	v	0.55	0/4176	0.69	2/5649 (0.0%)
25	A	0.31	0/69147	0.33	0/107869
26	B	0.23	0/2876	0.32	0/4483
27	C	0.30	0/2121	0.36	0/2852
28	D	0.30	0/1576	0.40	0/2119
29	E	0.26	0/1571	0.32	0/2113
30	F	0.17	0/1434	0.36	0/1926
31	G	0.20	0/1343	0.34	0/1816

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
32	I	0.15	0/1001	0.40	0/1350
33	J	0.13	0/1046	0.40	0/1410
34	L	0.29	0/1152	0.33	0/1551
35	M	0.29	0/955	0.36	0/1279
36	N	0.28	0/1061	0.36	0/1412
37	O	0.20	0/1073	0.31	0/1433
38	P	0.32	0/958	0.38	0/1281
39	Q	0.21	0/902	0.35	0/1209
40	R	0.28	0/929	0.34	0/1242
41	S	0.33	0/960	0.31	0/1278
42	T	0.28	0/829	0.37	0/1107
43	U	0.29	0/864	0.33	0/1156
44	V	0.26	0/744	0.34	0/994
45	W	0.25	0/787	0.46	0/1051
46	X	0.21	0/766	0.34	0/1025
47	Y	0.29	0/589	0.43	0/779
48	Z	0.29	0/635	0.34	0/848
49	1	0.23	0/496	0.32	0/660
50	2	0.28	0/453	0.31	0/605
51	3	0.10	0/380	0.28	0/508
52	4	0.31	0/440	0.33	0/588
53	5	0.23	0/424	0.34	0/565
54	6	0.32	0/380	0.30	0/498
55	7	0.28	0/513	0.40	0/676
56	8	0.26	0/303	0.39	0/397
All	All	0.28	0/160119	0.35	4/238847 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
5	e	0	1
37	O	0	1
55	7	0	1
All	All	0	3

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
18	r	69	PRO	CA-N-CD	-10.82	96.85	112.00
24	v	433	ASN	N-CA-C	7.12	119.41	108.30
18	r	69	PRO	N-CD-CG	-6.93	92.80	103.20
24	v	458	ASN	CB-CA-C	-5.02	110.81	116.63

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
55	7	30	ARG	Peptide
37	O	82	MS6	Peptide
5	e	55	GLU	Peptide

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	32803	0	16528	333	0
2	b	1754	0	1782	36	0
3	c	1624	0	1696	30	0
4	d	1643	0	1707	38	0
5	e	1152	0	1196	25	0
6	f	839	0	833	32	0
7	g	1191	0	1245	40	0
8	h	979	0	1031	15	0
9	i	1022	0	1070	34	0
10	j	786	0	828	26	0
11	k	877	0	884	21	0
12	l	942	0	999	14	0
13	m	891	0	952	24	0
14	n	805	0	844	23	0
15	o	714	0	734	7	0
16	p	643	0	661	8	0
17	q	641	0	682	14	0
18	r	446	0	472	6	0
19	s	663	0	688	21	0
20	t	670	0	719	8	0
21	u	460	0	486	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
22	x	1588	0	819	37	0
23	z	168	0	86	1	0
24	v	4099	0	4059	190	0
25	A	62252	0	31325	424	0
26	B	2572	0	1301	30	0
27	C	2082	0	2154	23	0
28	D	1566	0	1618	20	0
29	E	1552	0	1618	20	0
30	F	1410	0	1444	47	0
31	G	1323	0	1371	25	0
32	I	988	0	1025	38	0
33	J	1032	0	1085	50	0
34	L	1129	0	1162	11	0
35	M	946	0	1023	11	0
36	N	1052	0	1127	21	0
37	O	1075	0	1145	21	0
38	P	945	0	988	6	0
39	Q	892	0	923	14	0
40	R	917	0	962	11	0
41	S	947	0	1019	8	0
42	T	816	0	839	7	0
43	U	857	0	922	10	0
44	V	738	0	807	5	0
45	W	779	0	831	10	0
46	X	753	0	780	15	0
47	Y	582	0	599	5	0
48	Z	625	0	652	3	0
49	1	495	0	526	7	0
50	2	449	0	488	5	0
51	3	373	0	370	14	0
52	4	434	0	445	3	0
53	5	417	0	451	4	0
54	6	377	0	418	2	0
55	7	504	0	572	3	0
56	8	302	0	340	10	0
57	4	1	0	0	0	0
57	6	1	0	0	0	0
57	7	1	0	0	0	0
57	A	572	0	0	0	0
57	B	15	0	0	0	0
57	C	3	0	0	0	0
57	D	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
57	E	1	0	0	0	0
57	N	1	0	0	0	0
57	P	2	0	0	0	0
57	S	1	0	0	0	0
57	T	3	0	0	0	0
57	V	2	0	0	0	0
57	Y	1	0	0	0	0
57	a	125	0	0	0	0
57	n	1	0	0	0	0
57	p	1	0	0	0	0
57	x	1	0	0	0	0
57	z	1	0	0	0	0
58	v	32	0	14	9	0
59	3	1	0	0	0	0
59	8	1	0	0	0	0
60	6	1	0	0	0	0
60	7	1	0	0	0	0
60	8	1	0	0	0	0
60	A	202	0	0	2	0
60	B	7	0	0	0	0
60	C	1	0	0	0	0
60	D	3	0	0	0	0
60	N	1	0	0	0	0
60	P	1	0	0	0	0
60	S	2	0	0	0	0
60	a	29	0	0	0	0
60	n	1	0	0	0	0
60	x	1	0	0	0	0
60	z	1	0	0	0	0
All	All	149604	0	101345	1671	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 7.

All (1671) 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)
24:v:25:GLY:HA2	58:v:601:GCP:O1A	1.47	1.12
25:A:2114:A:H62	25:A:2119:A:N6	1.47	1.10
24:v:20:SER:HB3	24:v:26:LYS:HD3	1.31	1.08
24:v:21:HIS:ND1	24:v:22:PRO:HD2	1.73	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:A:2114:A:N6	25:A:2119:A:H61	1.56	1.02
25:A:284:U:H3	25:A:356:G:H1	1.09	1.00
24:v:259:ALA:HB3	58:v:601:GCP:O6	1.65	0.97
10:j:57:VAL:HG12	10:j:58:ASN:H	1.30	0.96
24:v:21:HIS:CE1	24:v:22:PRO:HD2	2.01	0.95
24:v:430:PRO:HD2	24:v:435:ASP:O	1.66	0.95
25:A:2114:A:H62	25:A:2119:A:H61	1.04	0.95
1:a:1441:A:H62	1:a:1461:G:H21	1.05	0.93
25:A:2134:A:C6	25:A:2156:G:C2	2.58	0.92
1:a:76:G:H1	1:a:93:U:H3	0.94	0.91
2:b:119:THR:O	2:b:123:ASP:HB2	1.73	0.89
46:X:83:LYS:HD2	46:X:84:PRO:HD2	1.55	0.89
24:v:127:MET:HG3	24:v:162:LEU:HD12	1.52	0.88
24:v:20:SER:HB3	24:v:26:LYS:CD	2.05	0.87
37:O:64:TRP:HB2	37:O:104:GLU:HB2	1.57	0.85
12:l:44:LYS:HD3	12:l:45:PRO:HA	1.57	0.84
13:m:7:ILE:HD11	13:m:22:ILE:HA	1.60	0.82
22:x:11:C:N4	22:x:24:G:H1	1.77	0.82
1:a:1441:A:H62	1:a:1461:G:N2	1.78	0.81
2:b:83:ALA:O	2:b:87:CYS:HB2	1.80	0.81
1:a:1221:G:OP1	19:s:36:ARG:NH2	2.15	0.80
10:j:5:ARG:HA	10:j:76:ILE:O	1.82	0.80
1:a:1441:A:N6	1:a:1461:G:H21	1.78	0.80
1:a:984:C:N3	1:a:1222:G:N2	2.29	0.79
22:x:26:A:N6	22:x:44:G:H1	1.79	0.79
25:A:544:C:O2	25:A:549:G:N2	2.14	0.79
25:A:2175:C:N3	25:A:2176:A:N6	2.31	0.79
1:a:501:C:OP1	12:l:114:ARG:NH2	2.15	0.79
32:I:25:ALA:HA	32:I:115:GLY:HA2	1.65	0.79
4:d:117:LEU:HD23	4:d:123:ILE:HD11	1.63	0.78
10:j:59:LYS:HD2	10:j:62:ARG:HH21	1.47	0.78
33:J:90:SER:HB2	33:J:93:PRO:HG3	1.66	0.78
15:o:17:ARG:NH1	15:o:26:GLU:OE2	2.18	0.77
26:B:54:G:H21	30:F:26:MET:HE1	1.47	0.77
24:v:520:ASP:OD1	24:v:521:VAL:N	2.18	0.77
7:g:28:ASN:HA	7:g:31:MET:HE3	1.66	0.76
42:T:6:GLN:NE2	42:T:37:GLU:OE2	2.16	0.76
24:v:68:ILE:HB	58:v:601:GCP:O1G	1.85	0.76
24:v:472:ARG:HD2	24:v:511:LEU:HD11	1.68	0.76
26:B:18:G:H1	26:B:65:U:H3	1.35	0.75
24:v:14:ARG:NH1	24:v:82:CYS:SG	2.59	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:I:59:LEU:HD23	32:I:61:ARG:H	1.52	0.75
37:O:136:MET:O	46:X:79:ARG:NH2	2.19	0.75
25:A:2134:A:C6	25:A:2156:G:N3	2.55	0.75
5:e:77:ASN:ND2	5:e:82:GLN:OE1	2.20	0.75
24:v:378:THR:HG21	24:v:387:PHE:H	1.51	0.75
25:A:2134:A:N6	25:A:2156:G:C4	2.54	0.75
24:v:520:ASP:OD2	24:v:522:GLN:NE2	2.20	0.74
22:x:53:G:H3'	22:x:54:5MU:H73	1.67	0.74
24:v:25:GLY:CA	58:v:601:GCP:O1A	2.33	0.74
25:A:627:A:OP1	36:N:78:ARG:NH2	2.20	0.74
25:A:1107:G:H4'	32:I:80:THR:HA	1.68	0.74
22:x:26:A:H61	22:x:44:G:H1	1.31	0.74
1:a:1008:U:H3'	1:a:1009:U:H5''	1.69	0.74
25:A:1086:A:O2'	25:A:1087:G:N7	2.21	0.74
25:A:1102:C:H2'	25:A:1103:A:H8	1.53	0.73
2:b:11:LYS:O	2:b:208:ARG:NH1	2.17	0.73
26:B:51:G:OP1	39:Q:63:LYS:NZ	2.20	0.73
24:v:402:ARG:NH1	24:v:464:GLU:OE1	2.21	0.73
24:v:21:HIS:CG	24:v:22:PRO:HD2	2.23	0.73
1:a:1009:U:O2	1:a:1020:G:N2	2.22	0.73
25:A:2134:A:C5	25:A:2156:G:N2	2.57	0.72
1:a:1379:G:N2	1:a:1381:U:O4	2.18	0.72
45:W:3:ALA:O	45:W:6:ARG:NH1	2.23	0.72
24:v:430:PRO:CD	24:v:435:ASP:O	2.37	0.72
24:v:168:PRO:HA	24:v:254:VAL:HG22	1.72	0.72
25:A:544:C:N3	25:A:549:G:N1	2.28	0.72
25:A:2134:A:C5	25:A:2156:G:C2	2.78	0.72
1:a:606:G:N2	1:a:632:U:OP1	2.21	0.71
24:v:21:HIS:CD2	24:v:120:GLU:HG2	2.24	0.71
1:a:1318:A:H1'	19:s:37:ARG:HH11	1.55	0.71
24:v:21:HIS:CG	24:v:22:PRO:CD	2.73	0.71
25:A:534:U:O2'	41:S:49:ASP:OD2	2.06	0.71
51:3:2:LYS:HD3	51:3:5:ILE:HD11	1.72	0.71
30:F:29:PRO:HB2	30:F:169:LEU:HD22	1.72	0.71
35:M:70:ARG:HG2	35:M:76:VAL:HG12	1.73	0.71
18:r:69:PRO:O	18:r:69:PRO:HD2	1.90	0.71
4:d:85:ASN:HB2	4:d:88:GLU:HG2	1.72	0.71
1:a:76:G:N2	1:a:93:U:O2	2.24	0.70
33:J:49:ILE:HG13	33:J:50:GLU:H	1.57	0.70
32:I:18:VAL:HG23	32:I:86:MET:HE2	1.73	0.70
6:f:5:GLU:OE1	6:f:63:ASN:ND2	2.25	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:v:142:ASN:OD1	24:v:143:LYS:N	2.24	0.70
9:i:65:ILE:HD12	9:i:79:ILE:HD12	1.73	0.70
19:s:11:ILE:HG12	19:s:16:LEU:HD13	1.73	0.70
24:v:260:LEU:N	58:v:601:GCP:O6	2.25	0.70
24:v:127:MET:HE1	24:v:137:ILE:HG21	1.72	0.70
9:i:106:ARG:NH1	9:i:107:ASP:O	2.25	0.70
11:k:81:ASN:OD1	11:k:106:ARG:NH2	2.25	0.70
1:a:254:G:O2'	17:q:18:GLU:O	2.09	0.69
24:v:422:GLU:N	24:v:422:GLU:OE1	2.24	0.69
1:a:1147:C:O2	9:i:18:ARG:NH1	2.24	0.69
17:q:79:VAL:HG12	17:q:80:GLU:HG3	1.73	0.69
24:v:76:GLN:OE1	24:v:85:ASN:ND2	2.25	0.69
25:A:2114:A:N6	25:A:2119:A:N6	2.23	0.69
22:x:51:U:H3	22:x:63:G:H1	0.77	0.69
24:v:171:TRP:NE1	24:v:230:GLU:OE1	2.25	0.69
10:j:25:ILE:HD11	10:j:92:LEU:HD11	1.74	0.69
24:v:20:SER:CB	24:v:26:LYS:HD3	2.17	0.69
1:a:932:C:OP1	7:g:4:ARG:NH2	2.25	0.68
31:G:16:ASP:OD2	31:G:27:LYS:NZ	2.26	0.68
6:f:11:HIS:HD2	6:f:12:PRO:HD2	1.57	0.68
25:A:61:C:OP2	49:1:47:ARG:NH2	2.26	0.68
28:D:1:MET:SD	28:D:2:ILE:N	2.65	0.68
1:a:981:U:O2'	14:n:61:ARG:NH1	2.27	0.68
25:A:2116:G:H22	25:A:2171:A:H61	1.40	0.68
28:D:46:ARG:NH1	28:D:85:ALA:O	2.27	0.68
34:L:13:ARG:NH2	34:L:49:ASP:O	2.27	0.68
25:A:2304:G:H22	25:A:2312:U:H3	1.41	0.68
9:i:120:LYS:HG2	9:i:121:ALA:H	1.59	0.67
15:o:40:GLN:HE22	25:A:716:A:H4'	1.59	0.67
24:v:304:GLN:HG2	24:v:307:MET:HE3	1.75	0.67
25:A:881:G:O6	25:A:895:U:C4	2.47	0.67
25:A:1068:G:N2	25:A:1095:A:O2'	2.27	0.67
3:c:114:LYS:HD3	3:c:185:ASN:HD21	1.59	0.67
24:v:21:HIS:HD2	24:v:122:ARG:H	1.42	0.67
25:A:2144:G:N2	25:A:2146:C:O2'	2.28	0.67
1:a:1071:C:H2'	1:a:1072:G:H8	1.59	0.67
4:d:118:VAL:HA	4:d:123:ILE:HD13	1.77	0.67
22:x:26:A:N1	22:x:44:G:N2	2.42	0.67
24:v:433:ASN:CG	24:v:434:ASN:H	2.03	0.67
33:J:81:LYS:HD2	33:J:87:LYS:HA	1.77	0.67
33:J:100:LYS:HB3	33:J:141:GLU:HB2	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:Q:24:THR:HG22	39:Q:42:PRO:HD3	1.76	0.67
1:a:1071:C:OP1	5:e:54:ARG:NH2	2.28	0.67
25:A:2123:G:N2	25:A:2172:U:OP1	2.26	0.67
1:a:1163:A:H2'	1:a:1164:G:H8	1.59	0.67
25:A:1365:A:OP2	48:Z:2:SER:OG	2.13	0.67
1:a:1307:U:OP1	13:m:100:GLN:NE2	2.26	0.67
24:v:77:PHE:HE1	24:v:86:LEU:HD23	1.60	0.67
24:v:416:LEU:HB3	24:v:427:VAL:HG21	1.76	0.67
25:A:1022:G:O6	34:L:68:LYS:NZ	2.27	0.67
17:q:59:VAL:HG12	17:q:78:VAL:HA	1.77	0.66
24:v:24:ALA:O	24:v:142:ASN:ND2	2.27	0.66
24:v:324:LYS:NZ	24:v:358:GLU:OE1	2.27	0.66
1:a:1180:A:OP2	9:i:99:ARG:NH2	2.28	0.66
24:v:124:ARG:HH12	24:v:161:GLU:HG2	1.59	0.66
1:a:1137:C:O2	1:a:1138:G:N2	2.28	0.66
22:x:20:U:H2'	22:x:21:A:H4'	1.78	0.66
1:a:80:A:N7	1:a:82:G:N2	2.41	0.66
1:a:427:U:OP1	4:d:13:ARG:NH2	2.29	0.66
1:a:1012:A:N6	1:a:1018:G:O6	2.29	0.66
1:a:1218:C:H2'	1:a:1219:A:C8	2.30	0.66
35:M:105:ARG:HH22	40:R:34:GLU:HB3	1.61	0.66
12:l:5:ASN:HD21	17:q:36:LYS:HE2	1.60	0.66
24:v:108:CYS:SG	24:v:109:CYS:N	2.68	0.66
25:A:1869:G:O2'	25:A:1871:A:N6	2.28	0.66
30:F:119:ALA:O	30:F:167:ARG:NH2	2.29	0.66
9:i:123:ARG:NH1	9:i:124:ARG:O	2.29	0.66
25:A:2749:A:OP1	31:G:2:SER:N	2.29	0.66
7:g:53:ARG:HH12	7:g:121:ALA:HB1	1.61	0.65
24:v:20:SER:OG	24:v:21:HIS:N	2.28	0.65
25:A:276:U:O2'	25:A:278:A:N7	2.29	0.65
1:a:1017:U:H2'	1:a:1018:G:H8	1.61	0.65
6:f:43:GLY:HA2	6:f:58:HIS:CE1	2.31	0.65
51:3:41:HIS:HB3	51:3:44:PHE:HB2	1.77	0.65
5:e:164:ILE:HG13	5:e:165:LEU:HD12	1.79	0.65
25:A:1250:G:N7	36:N:18:ARG:NH2	2.44	0.65
25:A:1065:U:O2	25:A:1073:A:N6	2.30	0.65
1:a:1355:G:H2'	1:a:1356:G:H8	1.61	0.65
1:a:1356:G:H2'	1:a:1357:A:C8	2.32	0.65
30:F:126:GLY:O	30:F:158:THR:OG1	2.15	0.65
15:o:33:THR:HG23	15:o:63:ARG:HH12	1.62	0.64
22:x:15:G:N1	22:x:48:C:N3	2.45	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:A:2392:A:H2	36:N:55:MET:HE2	1.61	0.64
28:D:133:THR:HG22	28:D:134:HIS:H	1.62	0.64
1:a:1297:G:N2	7:g:114:LYS:O	2.30	0.64
24:v:26:LYS:HZ1	24:v:91:GLY:HA3	1.62	0.64
33:J:12:GLN:HE22	33:J:57:VAL:H	1.45	0.64
4:d:116:GLN:HE21	4:d:154:ARG:HH12	1.45	0.64
22:x:15:G:N2	22:x:48:C:N4	2.45	0.64
25:A:276:U:H1'	25:A:278:A:H62	1.63	0.64
25:A:1665:A:H1'	35:M:1:MET:HG2	1.80	0.64
25:A:1582:C:O2'	25:A:1585:C:N3	2.30	0.64
25:A:1802:A:H2'	25:A:1803:A:C8	2.33	0.64
27:C:107:PRO:HD2	27:C:110:LEU:HD22	1.78	0.64
1:a:429:U:H3'	4:d:9:LEU:HD12	1.80	0.64
32:I:38:MET:SD	33:J:118:THR:OG1	2.56	0.64
32:I:87:GLU:HG2	32:I:95:LEU:HB2	1.79	0.64
3:c:114:LYS:HD3	3:c:185:ASN:ND2	2.12	0.63
25:A:515:A:N7	60:A:3610:HOH:O	2.30	0.63
25:A:1048:A:OP2	25:A:1110:G:N2	2.31	0.63
25:A:1223:G:OP2	42:T:90:ARG:NH1	2.31	0.63
24:v:306:ASN:ND2	24:v:456:GLU:OE1	2.32	0.63
24:v:492:GLN:N	24:v:492:GLN:OE1	2.30	0.63
38:P:2:ARG:O	38:P:2:ARG:HD2	1.98	0.63
25:A:2135:A:N6	25:A:2156:G:O2'	2.30	0.63
33:J:39:CYS:O	33:J:43:ASN:ND2	2.31	0.63
1:a:1305:G:N2	1:a:1331:G:O2'	2.30	0.63
26:B:1:U:H2'	26:B:2:G:C8	2.33	0.63
25:A:2478:A:H5'	56:8:32:LYS:HE2	1.79	0.63
27:C:232:HIS:HA	27:C:242:LYS:HE3	1.80	0.63
49:1:2:LYS:NZ	49:1:5:GLU:OE2	2.22	0.63
5:e:61:GLN:O	5:e:65:GLU:HG3	1.99	0.63
24:v:190:THR:HG21	24:v:215:LEU:HD12	1.79	0.63
24:v:259:ALA:CB	58:v:601:GCP:O6	2.43	0.63
25:A:2134:A:N6	25:A:2157:G:N3	2.45	0.63
31:G:42:GLU:OE1	31:G:55:ARG:NH1	2.32	0.63
1:a:517:G:N2	1:a:530:G:OP1	2.31	0.63
1:a:1328:C:OP1	13:m:28:THR:OG1	2.13	0.63
24:v:446:PHE:HB3	24:v:463:TYR:HE2	1.64	0.63
25:A:870:U:O2	25:A:907:G:O6	2.16	0.63
20:t:35:VAL:HG21	20:t:54:MET:HG2	1.80	0.63
25:A:885:C:H2'	25:A:886:A:C8	2.33	0.63
25:A:1590:A:H2'	25:A:1591:A:H8	1.64	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:b:42:ASN:HD22	2:b:45:LYS:HG2	1.64	0.62
13:m:49:SER:N	13:m:52:GLN:OE1	2.32	0.62
25:A:463:G:N2	25:A:466:A:OP2	2.27	0.62
1:a:208:U:HO2'	1:a:211:G:H1	1.47	0.62
1:a:1317:C:OP1	14:n:24:ARG:NH2	2.32	0.62
13:m:28:THR:HA	13:m:31:LYS:HE2	1.80	0.62
25:A:2114:A:H3'	25:A:2115:G:H8	1.64	0.62
8:h:103:VAL:HG23	8:h:126:ILE:HB	1.80	0.62
24:v:478:ASP:OD1	24:v:479:ALA:N	2.33	0.62
5:e:56:VAL:HG23	5:e:57:PRO:HD3	1.81	0.62
1:a:1163:A:H2'	1:a:1164:G:C8	2.34	0.62
17:q:64:CYS:SG	17:q:74:THR:OG1	2.57	0.62
33:J:13:VAL:HG13	33:J:24:VAL:HG21	1.82	0.62
32:I:28:ALA:HA	32:I:82:ILE:HA	1.80	0.62
37:O:17:ASN:O	37:O:38:ARG:NH1	2.31	0.62
1:a:1191:A:OP2	3:c:2:GLY:N	2.33	0.62
4:d:62:ARG:NH1	4:d:69:GLU:OE1	2.33	0.62
1:a:993:G:O2'	1:a:995:C:N4	2.33	0.62
24:v:396:GLU:OE1	24:v:396:GLU:N	2.33	0.62
25:A:856:G:H2'	25:A:857:G:C8	2.34	0.61
12:l:68:GLY:O	12:l:99:ARG:NH1	2.31	0.61
22:x:11:C:N3	22:x:24:G:N2	2.43	0.61
1:a:1249:C:O2'	9:i:71:GLY:O	2.17	0.61
2:b:15:HIS:HB3	2:b:43:LEU:HD21	1.83	0.61
14:n:54:ASP:OD1	14:n:59:ARG:NH2	2.33	0.61
22:x:51:U:O2	22:x:63:G:N2	2.22	0.61
25:A:1054:A:H2'	25:A:1055:G:C8	2.36	0.61
1:a:673:A:H2'	1:a:674:G:C8	2.35	0.61
9:i:21:ILE:HG12	9:i:63:LEU:HD22	1.82	0.61
30:F:103:LEU:HA	30:F:107:ALA:HB3	1.80	0.61
3:c:55:ILE:HD12	3:c:68:ILE:HG22	1.81	0.61
25:A:495:G:H21	43:U:61:ASN:HD21	1.49	0.61
32:I:94:ARG:HH21	32:I:101:LYS:HE3	1.64	0.61
19:s:9:PRO:HB2	19:s:39:THR:HG21	1.80	0.61
24:v:21:HIS:ND1	24:v:22:PRO:CD	2.57	0.61
1:a:1124:G:N2	1:a:1125:U:O4	2.19	0.61
1:a:1319:A:O2'	1:a:1323:G:N7	2.30	0.61
1:a:1321:U:H3	19:s:36:ARG:HH22	1.47	0.61
25:A:1083:U:OP1	32:I:53:ARG:NH1	2.34	0.61
27:C:3:VAL:HG22	27:C:19:VAL:HG12	1.82	0.61
28:D:35:THR:HG22	28:D:73:VAL:HG21	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:a:958:A:N6	19:s:77:THR:O	2.34	0.61
5:e:83:HIS:HE1	5:e:148:ASN:H	1.47	0.61
1:a:1130:A:OP1	9:i:18:ARG:NH2	2.34	0.60
6:f:10:VAL:HG12	6:f:58:HIS:HB3	1.82	0.60
1:a:137:U:H3	1:a:226:G:H1	1.47	0.60
22:x:11:C:N4	22:x:24:G:N1	2.49	0.60
51:3:15:SER:O	51:3:33:ASN:ND2	2.31	0.60
22:x:76:A:O2'	25:A:2394:C:N3	2.33	0.60
25:A:2126:A:H1'	25:A:2173:A:H61	1.67	0.60
30:F:108:VAL:HG11	30:F:176:PRO:HG3	1.83	0.60
24:v:365:LEU:HD23	24:v:365:LEU:H	1.66	0.60
25:A:138:U:O2'	25:A:141:G:O6	2.17	0.60
25:A:848:C:H2'	25:A:849:A:H8	1.66	0.60
25:A:1040:A:N1	25:A:1115:G:O6	2.34	0.60
25:A:1153:C:OP1	41:S:92:ARG:NH2	2.34	0.60
25:A:2031:A:N3	25:A:2455:G:O2'	2.29	0.60
25:A:2467:C:OP1	56:8:8:LYS:NZ	2.30	0.60
9:i:46:MET:O	9:i:50:GLN:HG3	2.02	0.60
25:A:476:G:N1	25:A:479:A:OP2	2.34	0.60
46:X:42:LEU:HD13	46:X:47:VAL:HG21	1.84	0.60
55:7:17:THR:HG21	55:7:49:MET:HE1	1.83	0.60
13:m:107:ARG:NH1	13:m:111:GLY:O	2.35	0.60
24:v:314:ARG:NH1	24:v:418:GLN:OE1	2.34	0.60
11:k:37:ARG:NH1	11:k:83:GLU:OE2	2.35	0.60
22:x:18:G:O2'	22:x:57:G:N2	2.35	0.60
24:v:110:LEU:HD11	24:v:268:MET:HE2	1.81	0.60
26:B:2:G:H2'	26:B:3:C:H6	1.66	0.60
26:B:5:U:OP1	26:B:61:G:O2'	2.16	0.60
1:a:1355:G:H2'	1:a:1356:G:C8	2.36	0.60
13:m:20:THR:HA	13:m:25:VAL:HG13	1.84	0.60
31:G:25:THR:HG23	31:G:34:THR:HG22	1.83	0.60
1:a:1022:A:H2'	1:a:1023:U:O4'	2.01	0.60
25:A:500:G:N1	25:A:503:A:OP2	2.33	0.60
26:B:1:U:H2'	26:B:2:G:H8	1.67	0.60
30:F:135:GLN:OE1	30:F:150:ARG:N	2.30	0.60
1:a:1058:G:OP1	3:c:199:LYS:NZ	2.34	0.59
1:a:544:G:OP1	4:d:56:ARG:NH2	2.34	0.59
25:A:1649:G:O2'	38:P:106:ASP:OD2	2.19	0.59
1:a:337:G:H2'	1:a:338:A:C8	2.37	0.59
25:A:203:A:N7	60:A:3614:HOH:O	2.32	0.59
35:M:76:VAL:HG22	40:R:73:VAL:HB	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:A:2278:A:OP1	37:O:10:ARG:NH1	2.35	0.59
25:A:2312:U:H5'	30:F:85:ILE:HD11	1.84	0.59
17:q:58:VAL:HG22	17:q:80:GLU:HB2	1.84	0.59
27:C:30:PHE:HD2	27:C:33:LEU:HD12	1.66	0.59
44:V:3:ARG:NH1	44:V:5:GLU:OE2	2.31	0.59
1:a:1207:2MG:H2'	1:a:1208:C:C6	2.37	0.59
5:e:36:LEU:HD22	5:e:134:ILE:HD13	1.84	0.59
1:a:147:G:H2'	1:a:148:G:C8	2.38	0.59
1:a:637:C:OP1	17:q:4:LYS:NZ	2.27	0.59
25:A:1095:A:H2'	25:A:1096:A:C8	2.38	0.59
43:U:86:MET:HE3	43:U:88:ARG:HD3	1.85	0.59
25:A:1077:A:H4'	33:J:93:PRO:HG2	1.84	0.59
25:A:1730:C:H1'	25:A:1731:G:C6	2.38	0.59
25:A:2139:U:O2	25:A:2152:G:O6	2.21	0.59
24:v:31:GLU:HG3	24:v:55:ALA:O	2.02	0.58
25:A:1645:G:H5''	25:A:1646:C:H5'	1.83	0.58
1:a:28:A:O2'	1:a:296:U:OP1	2.20	0.58
24:v:296:PHE:HE2	24:v:325:TYR:HB2	1.68	0.58
24:v:517:ARG:HG3	24:v:518:TYR:CD2	2.38	0.58
25:A:2099:U:O2	25:A:2190:G:O6	2.21	0.58
32:I:51:TYR:HE2	32:I:88:HIS:HB2	1.67	0.58
24:v:114:ASP:OD1	24:v:115:ALA:N	2.35	0.58
1:a:1378:C:OP1	7:g:5:ARG:NH1	2.36	0.58
10:j:57:VAL:HG12	10:j:58:ASN:N	2.11	0.58
13:m:29:ARG:O	13:m:33:ILE:HG12	2.04	0.58
24:v:29:ILE:HD11	24:v:112:VAL:HG21	1.85	0.58
24:v:396:GLU:O	24:v:397:LEU:HG	2.03	0.58
25:A:729:G:H5''	25:A:730:A:H5''	1.85	0.58
1:a:408:A:O3'	4:d:23:SER:OG	2.22	0.58
1:a:1175:G:H2'	1:a:1176:A:H8	1.69	0.58
8:h:67:GLN:HE21	8:h:69:LYS:HB3	1.69	0.58
24:v:124:ARG:NH2	24:v:161:GLU:OE2	2.29	0.58
25:A:2246:G:H2'	25:A:2247:A:C8	2.39	0.58
43:U:92:ARG:NH2	43:U:94:ASP:OD1	2.37	0.58
7:g:15:ASP:OD1	7:g:20:SER:N	2.26	0.58
7:g:72:THR:HG23	7:g:73:VAL:HG12	1.85	0.58
14:n:17:ALA:HA	14:n:55:SER:HA	1.85	0.58
19:s:11:ILE:HG21	19:s:16:LEU:HD22	1.84	0.58
1:a:1081:A:OP2	5:e:52:LYS:NZ	2.31	0.58
3:c:59:ARG:HG2	3:c:64:ILE:HG22	1.85	0.58
25:A:881:G:O6	25:A:895:U:O4	2.22	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:m:20:THR:OG1	13:m:25:VAL:O	2.19	0.58
24:v:304:GLN:HG3	24:v:305:ALA:H	1.67	0.58
25:A:1528:A:N6	25:A:1543:G:O2'	2.37	0.58
1:a:324:G:N1	1:a:327:A:OP2	2.37	0.58
22:x:41:C:H2'	22:x:42:C:C6	2.39	0.58
25:A:2302:U:O2'	30:F:123:ASP:O	2.19	0.58
28:D:85:ALA:HB3	28:D:88:GLU:HG3	1.85	0.58
32:I:73:LYS:HB2	32:I:117:LEU:HB2	1.86	0.58
9:i:30:ILE:HG12	9:i:65:ILE:HD11	1.85	0.57
24:v:20:SER:HB3	24:v:26:LYS:CE	2.33	0.57
2:b:66:LYS:HB2	2:b:90:PHE:HE2	1.68	0.57
25:A:993:G:H1'	42:T:91:GLN:HE21	1.69	0.57
32:I:78:GLY:N	32:I:79:PRO:HD2	2.18	0.57
1:a:35:G:H2'	1:a:36:C:C6	2.39	0.57
15:o:76:ALA:O	15:o:80:GLN:HG3	2.03	0.57
25:A:2134:A:N6	25:A:2156:G:C2	2.71	0.57
37:O:74:THR:HG21	37:O:86:LYS:HE3	1.85	0.57
7:g:111:ARG:NH1	7:g:119:ARG:O	2.38	0.57
25:A:2312:U:O2	30:F:37:ASN:ND2	2.37	0.57
28:D:14:ILE:HA	40:R:12:GLN:HE22	1.69	0.57
56:8:6:SER:OG	56:8:8:LYS:NZ	2.37	0.57
17:q:10:GLY:HA3	17:q:25:ILE:HD13	1.87	0.57
25:A:848:C:H2'	25:A:849:A:C8	2.40	0.57
25:A:1103:A:H3'	25:A:1104:C:H5''	1.87	0.57
6:f:11:HIS:CD2	6:f:12:PRO:HD2	2.39	0.57
30:F:134:GLU:OE1	30:F:136:ILE:N	2.34	0.57
24:v:474:VAL:HG23	24:v:475:GLU:H	1.70	0.57
25:A:1817:G:OP1	27:C:87:ARG:NH2	2.38	0.57
27:C:180:GLU:OE2	27:C:267:ILE:HG23	2.05	0.57
29:E:48:THR:HG22	29:E:86:ALA:HB3	1.86	0.57
6:f:18:VAL:HG11	6:f:58:HIS:CD2	2.40	0.57
1:a:946:A:H2'	1:a:947:G:C8	2.39	0.57
6:f:5:GLU:CD	6:f:63:ASN:HD21	2.13	0.57
22:x:23:A:H62	22:x:46:G7M:H21	1.53	0.57
29:E:6:LYS:O	29:E:9:GLN:NE2	2.37	0.57
45:W:24:LYS:H	45:W:37:GLU:HG2	1.70	0.57
1:a:17:U:H2'	1:a:18:C:C6	2.40	0.56
1:a:356:A:N3	1:a:368:U:O2'	2.34	0.56
1:a:1314:C:H2'	1:a:1315:U:C6	2.39	0.56
10:j:45:ARG:HB3	10:j:69:THR:HB	1.87	0.56
20:t:9:LYS:O	20:t:13:GLN:HG2	2.04	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:v:104:THR:O	24:v:282:ARG:NH1	2.38	0.56
25:A:1084:A:OP2	32:I:55:VAL:HA	2.05	0.56
25:A:1434:A:H2'	25:A:1435:G:H8	1.70	0.56
1:a:908:A:H2'	1:a:909:A:C8	2.41	0.56
24:v:88:ASP:OD1	24:v:89:THR:N	2.37	0.56
26:B:42:C:C5	30:F:66:LEU:HD12	2.40	0.56
53:5:5:ILE:HD12	53:5:28:ARG:HH12	1.69	0.56
25:A:1094:U:H1'	25:A:1097:U:H5	1.70	0.56
25:A:1597:A:H5''	25:A:1598:A:H5'	1.86	0.56
25:A:1927:A:H2'	25:A:1928:A:C8	2.40	0.56
25:A:2246:G:H2'	25:A:2247:A:H8	1.70	0.56
25:A:2328:A:H2'	25:A:2329:U:C6	2.40	0.56
1:a:618:C:H1'	16:p:14:ARG:HH12	1.70	0.56
1:a:746:A:H2'	1:a:747:A:C8	2.41	0.56
1:a:1250:A:N3	1:a:1370:G:O2'	2.36	0.56
3:c:73:PRO:HA	3:c:76:VAL:HG12	1.86	0.56
24:v:281:PRO:HB3	24:v:290:GLU:HA	1.88	0.56
25:A:2064:C:H2'	25:A:2065:C:H6	1.70	0.56
4:d:196:ASN:HB3	4:d:199:LEU:HD23	1.88	0.56
8:h:10:MET:HG3	8:h:27:MET:SD	2.46	0.56
25:A:2183:A:H2'	25:A:2184:A:C8	2.39	0.56
43:U:82:MET:HB2	43:U:98:LYS:HB2	1.87	0.56
7:g:150:ALA:HB2	11:k:61:PHE:HB2	1.88	0.56
24:v:60:MET:SD	24:v:452:ARG:NH2	2.78	0.56
25:A:597:G:O2'	36:N:11:GLY:O	2.21	0.56
25:A:887:U:O2'	25:A:889:C:OP2	2.24	0.56
8:h:64:LYS:HG3	8:h:71:VAL:HG21	1.88	0.56
10:j:11:LYS:HG2	10:j:71:LEU:HD23	1.88	0.56
24:v:284:THR:HG21	24:v:385:MET:HE1	1.87	0.56
25:A:2064:C:H2'	25:A:2065:C:C6	2.40	0.56
1:a:210:C:H4'	1:a:211:G:C6	2.40	0.56
1:a:1265:C:H2'	1:a:1266:G:C8	2.41	0.56
25:A:885:C:H2'	25:A:886:A:H8	1.70	0.56
25:A:1138:G:H21	34:L:108:MET:HE2	1.69	0.56
29:E:108:ILE:HG23	36:N:1:MET:HE1	1.87	0.56
1:a:1297:G:C5	7:g:114:LYS:HE3	2.41	0.56
3:c:185:ASN:OD1	3:c:186:THR:N	2.39	0.56
41:S:66:ASN:O	41:S:70:ARG:HG3	2.06	0.56
1:a:1015:G:H2'	1:a:1016:A:C8	2.41	0.55
25:A:1059:G:OP1	33:J:72:LYS:NZ	2.34	0.55
25:A:2128:G:N3	25:A:2173:A:O2'	2.39	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:J:79:LEU:HA	33:J:82:LYS:HZ1	1.70	0.55
40:R:89:ARG:NH2	40:R:115:ASN:O	2.39	0.55
1:a:1157:A:N6	1:a:1180:A:N7	2.54	0.55
10:j:37:ARG:NH2	10:j:75:ASP:OD2	2.39	0.55
24:v:102:THR:HG22	24:v:301:PHE:HE1	1.70	0.55
25:A:893:C:H2'	25:A:894:U:C6	2.41	0.55
25:A:2502:G:H5''	25:A:2503:2MA:H5''	1.88	0.55
30:F:123:ASP:OD1	30:F:124:GLY:N	2.38	0.55
1:a:1402:4OC:HM22	1:a:1403:C:H5'	1.88	0.55
9:i:33:ARG:HH21	9:i:38:TYR:HA	1.71	0.55
25:A:728:G:H4'	27:C:13:ARG:HD3	1.88	0.55
25:A:1071:G:H2'	25:A:1072:C:C6	2.42	0.55
25:A:1590:A:H2'	25:A:1591:A:C8	2.41	0.55
26:B:2:G:H2'	26:B:3:C:C6	2.41	0.55
31:G:52:PHE:HZ	31:G:72:LEU:HD23	1.72	0.55
24:v:20:SER:H	24:v:26:LYS:HE2	1.70	0.55
25:A:5:A:H2'	25:A:6:A:C8	2.42	0.55
1:a:1013:G:N2	1:a:1016:A:OP2	2.33	0.55
7:g:53:ARG:NH1	7:g:121:ALA:HB1	2.22	0.55
24:v:20:SER:HB3	24:v:26:LYS:HE3	1.89	0.55
25:A:877:A:O2'	25:A:900:A:N6	2.38	0.55
29:E:58:LYS:HG3	29:E:71:GLY:HA2	1.89	0.55
32:I:38:MET:HG2	32:I:42:ARG:HE	1.70	0.55
2:b:83:ALA:O	2:b:87:CYS:CB	2.53	0.55
21:u:40:LYS:O	21:u:43:THR:OG1	2.21	0.55
25:A:1469:A:H2'	25:A:1470:A:C8	2.42	0.55
26:B:42:C:OP2	51:3:2:LYS:NZ	2.39	0.55
30:F:4:LEU:O	30:F:8:TYR:N	2.39	0.55
30:F:135:GLN:HE22	30:F:149:VAL:HA	1.72	0.55
31:G:17:VAL:HA	31:G:26:ILE:HG22	1.88	0.55
33:J:83:ALA:O	33:J:105:GLN:NE2	2.39	0.55
1:a:674:G:H2'	1:a:675:A:H8	1.72	0.55
21:u:4:ILE:HG13	21:u:19:PHE:HA	1.88	0.55
25:A:5:A:H2'	25:A:6:A:H8	1.70	0.55
28:D:1:MET:HE3	28:D:205:PRO:HG2	1.89	0.55
32:I:37:LYS:NZ	32:I:56:ARG:HH21	2.04	0.55
33:J:134:ARG:HD3	33:J:139:VAL:HG13	1.89	0.55
38:P:24:MET:HE3	38:P:36:THR:HG21	1.89	0.55
24:v:119:VAL:HG11	24:v:162:LEU:HD22	1.87	0.55
24:v:127:MET:HE1	24:v:137:ILE:HD13	1.87	0.55
24:v:493:LEU:HD12	24:v:501:LEU:HD11	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:F:20:PHE:CE2	30:F:165:GLU:HG3	2.42	0.55
31:G:18:LYS:HE3	31:G:25:THR:HB	1.89	0.55
46:X:61:LEU:HD11	46:X:91:PHE:HE1	1.72	0.55
25:A:1106:G:H1'	32:I:81:LEU:HD12	1.88	0.55
25:A:1433:A:H2'	25:A:1434:A:C8	2.42	0.55
25:A:1918:A:O2'	25:A:1920:C:N4	2.40	0.55
1:a:1149:C:OP2	9:i:11:ARG:NH2	2.38	0.54
7:g:113:ASP:HB2	7:g:119:ARG:HG3	1.89	0.54
24:v:182:VAL:HG13	24:v:263:PHE:HE1	1.72	0.54
25:A:1:G:H2'	25:A:2:G:H8	1.73	0.54
1:a:460:A:H2'	1:a:461:A:H8	1.72	0.54
1:a:745:G:H5'	1:a:851:G:H21	1.72	0.54
16:p:44:SER:OG	16:p:47:GLU:OE2	2.22	0.54
22:x:51:U:O4	22:x:63:G:O6	2.25	0.54
25:A:1434:A:H2'	25:A:1435:G:C8	2.42	0.54
30:F:102:ARG:NH1	51:3:9:TYR:OH	2.40	0.54
33:J:90:SER:HB3	33:J:136:MET:HA	1.88	0.54
37:O:21:ALA:HB2	37:O:97:GLN:HB2	1.88	0.54
39:Q:90:VAL:HG23	39:Q:117:PHE:HB3	1.90	0.54
1:a:1228:C:H2'	1:a:1229:A:H8	1.72	0.54
4:d:121:LYS:O	4:d:146:ARG:NH2	2.41	0.54
25:A:1495:A:N3	25:A:1578:U:O2'	2.39	0.54
1:a:1071:C:H2'	1:a:1072:G:C8	2.42	0.54
1:a:1407:5MC:H2'	1:a:1408:A:H8	1.73	0.54
25:A:1868:C:N4	25:A:1869:G:O6	2.41	0.54
25:A:1141:U:OP2	34:L:65:THR:OG1	2.22	0.54
29:E:176:ASP:OD1	29:E:179:SER:OG	2.24	0.54
6:f:21:MET:O	6:f:24:ARG:HG2	2.08	0.54
24:v:150:ASP:N	24:v:150:ASP:OD1	2.40	0.54
24:v:296:PHE:CE2	24:v:325:TYR:HB2	2.43	0.54
25:A:547:A:H1'	25:A:548:G:C8	2.42	0.54
25:A:2189:U:H2'	25:A:2190:G:C8	2.43	0.54
27:C:100:GLU:OE2	27:C:102:ARG:NE	2.32	0.54
6:f:2:ARG:HE	6:f:91:ARG:NH2	2.06	0.54
25:A:284:U:O4	25:A:356:G:O6	2.25	0.54
25:A:2106:U:H2'	25:A:2107:G:H8	1.73	0.54
12:l:55:VAL:HG11	12:l:80:ILE:HD11	1.88	0.54
25:A:2099:U:H2'	25:A:2100:G:C8	2.43	0.54
10:j:57:VAL:CG1	10:j:58:ASN:H	2.11	0.54
24:v:105:ALA:HB1	24:v:319:ARG:HG3	1.89	0.54
25:A:2392:A:C2	36:N:55:MET:HE2	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:F:121:SER:OG	30:F:129:SER:O	2.21	0.54
30:F:134:GLU:OE2	30:F:137:ILE:HG13	2.08	0.54
1:a:1258:G:N2	1:a:1279:G:H22	2.05	0.54
25:A:2636:C:O2'	28:D:45:TYR:OH	2.25	0.54
31:G:85:LYS:N	31:G:133:LEU:O	2.34	0.54
3:c:10:ILE:HG23	3:c:11:ARG:HG3	1.89	0.53
10:j:40:ILE:HD11	10:j:73:LEU:HD23	1.90	0.53
16:p:55:ASP:OD1	16:p:56:ARG:N	2.40	0.53
1:a:187:G:N2	1:a:190:A:OP2	2.39	0.53
1:a:404:G:OP2	4:d:115:ARG:NH2	2.36	0.53
25:A:784:G:H5'	25:A:785:G:OP1	2.08	0.53
25:A:1528:A:H2'	25:A:1529:G:O4'	2.08	0.53
31:G:12:PRO:HD2	31:G:15:VAL:HG11	1.90	0.53
1:a:750:C:O2'	15:o:21:ASP:OD1	2.26	0.53
1:a:1250:A:H2'	1:a:1251:A:C8	2.43	0.53
1:a:1346:A:OP1	9:i:122:ARG:NH1	2.31	0.53
14:n:24:ARG:NH1	14:n:55:SER:OG	2.41	0.53
26:B:106:G:H2'	26:B:107:G:O4'	2.08	0.53
24:v:110:LEU:CD1	24:v:268:MET:HE2	2.39	0.53
37:O:106:ASP:OD1	37:O:107:GLY:N	2.41	0.53
1:a:1115:U:H2'	1:a:1116:U:C6	2.43	0.53
1:a:1216:A:H5''	14:n:5:SER:OG	2.09	0.53
5:e:115:LEU:HD13	5:e:123:VAL:HG11	1.90	0.53
25:A:549:G:H2'	25:A:550:C:H6	1.73	0.53
25:A:1141:U:H4'	25:A:1142:A:O4'	2.09	0.53
30:F:134:GLU:OE1	30:F:135:GLN:N	2.41	0.53
1:a:459:A:H2'	1:a:460:A:C8	2.44	0.53
25:A:1054:A:H2'	25:A:1055:G:H8	1.74	0.53
25:A:1527:G:N1	25:A:1544:A:OP2	2.36	0.53
25:A:2191:A:H2'	25:A:2192:U:C6	2.44	0.53
1:a:1518:MA6:H92	1:a:1519:MA6:N6	2.24	0.53
22:x:15:G:H22	22:x:48:C:N4	2.05	0.53
25:A:1889:A:N3	25:A:2086:U:O2'	2.40	0.53
1:a:472:U:H2'	1:a:473:U:C6	2.44	0.53
1:a:613:C:OP1	4:d:81:ARG:NH1	2.42	0.53
4:d:49:SER:OG	4:d:50:ASP:N	2.42	0.53
6:f:21:MET:HG2	6:f:25:TYR:CE2	2.44	0.53
11:k:36:ASP:OD1	11:k:37:ARG:N	2.42	0.53
24:v:46:VAL:HG11	58:v:601:GCP:H5'2	1.91	0.53
44:V:38:ALA:O	44:V:81:LYS:NZ	2.39	0.53
1:a:1367:C:H5'	10:j:62:ARG:NH1	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:v:17:ALA:HB2	24:v:106:VAL:HG21	1.90	0.53
25:A:2305:U:O2	30:F:151:GLY:HA3	2.09	0.53
32:I:33:VAL:HG12	32:I:35:VAL:HG22	1.91	0.53
37:O:47:GLU:OE2	37:O:50:ARG:NH2	2.25	0.53
43:U:48:LYS:O	43:U:52:GLU:HG3	2.09	0.53
1:a:946:A:H2'	1:a:947:G:H8	1.74	0.53
3:c:85:GLU:OE1	3:c:89:LYS:NZ	2.41	0.53
13:m:25:VAL:HG23	13:m:29:ARG:HB2	1.90	0.53
25:A:1138:G:N2	34:L:108:MET:HE2	2.23	0.53
25:A:1179:G:H2'	25:A:1180:U:C6	2.44	0.53
26:B:31:C:O2'	26:B:53:A:N6	2.40	0.53
24:v:126:LEU:O	24:v:130:THR:HG22	2.09	0.52
25:A:895:U:H2'	25:A:897:C:C5	2.44	0.52
26:B:52:A:O2'	26:B:53:A:H8	1.92	0.52
1:a:1149:C:H2'	1:a:1150:A:C8	2.45	0.52
14:n:24:ARG:HD2	14:n:55:SER:HB2	1.92	0.52
25:A:871:U:H2'	25:A:872:U:C6	2.44	0.52
25:A:2111:U:N3	25:A:2147:A:N3	2.57	0.52
25:A:2498:OMC:HM22	25:A:2499:C:H5'	1.90	0.52
1:a:475:C:H2'	1:a:476:U:C6	2.45	0.52
1:a:652:U:O4	1:a:752:G:O2'	2.27	0.52
10:j:8:ILE:HG13	10:j:100:ILE:HG23	1.91	0.52
13:m:4:ILE:HG13	13:m:57:ARG:HG2	1.91	0.52
24:v:21:HIS:CD2	24:v:122:ARG:H	2.23	0.52
24:v:126:LEU:HA	24:v:129:VAL:HG12	1.92	0.52
25:A:2106:U:H2'	25:A:2107:G:C8	2.43	0.52
1:a:139:A:H2'	1:a:140:U:C6	2.45	0.52
1:a:527:G7M:HN73	12:l:46:ASN:HD21	1.74	0.52
24:v:454:LYS:HA	24:v:458:ASN:O	2.08	0.52
24:v:476:CYS:SG	24:v:521:VAL:HG22	2.50	0.52
25:A:1814:G:OP2	25:A:1815:A:O2'	2.25	0.52
25:A:2187:U:H2'	25:A:2188:U:C5	2.45	0.52
26:B:54:G:N2	30:F:26:MET:HE1	2.21	0.52
1:a:1017:U:H2'	1:a:1018:G:C8	2.42	0.52
7:g:47:LEU:HD12	7:g:50:LEU:HD11	1.92	0.52
10:j:11:LYS:HA	10:j:70:HIS:O	2.10	0.52
24:v:34:LEU:HD12	24:v:39:ALA:HB3	1.91	0.52
25:A:1060:U:H4'	25:A:1062:G:H5''	1.92	0.52
1:a:458:U:H2'	1:a:459:A:H8	1.75	0.52
6:f:19:PRO:O	6:f:22:ILE:HB	2.09	0.52
24:v:119:VAL:HG13	24:v:123:THR:HG23	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:A:2141:G:H2'	25:A:2142:A:C8	2.45	0.52
26:B:39:A:N1	26:B:44:G:C6	2.77	0.52
32:I:3:LEU:HG	32:I:5:LEU:H	1.75	0.52
1:a:459:A:H2'	1:a:460:A:H8	1.75	0.52
1:a:528:C:N4	12:l:46:ASN:OD1	2.42	0.52
1:a:1320:C:N3	19:s:72:GLY:HA3	2.25	0.52
25:A:636:G:N7	36:N:109:LYS:HE2	2.25	0.52
32:I:73:LYS:HE3	32:I:120:ALA:HB2	1.92	0.52
4:d:60:LYS:O	4:d:64:ILE:HG12	2.10	0.52
25:A:1315:C:O2'	25:A:1392:A:N3	2.37	0.52
10:j:10:LEU:HB3	10:j:98:VAL:HG13	1.91	0.52
25:A:851:C:H2'	25:A:852:U:C6	2.45	0.52
33:J:92:LYS:HB3	33:J:95:LYS:HB2	1.92	0.52
33:J:103:ARG:O	33:J:106:LEU:HG	2.10	0.52
1:a:216:U:H2'	1:a:217:C:C6	2.45	0.52
1:a:390:U:H2'	1:a:391:G:C8	2.45	0.52
1:a:929:G:O2'	1:a:930:C:OP1	2.26	0.52
5:e:85:VAL:HG11	5:e:147:MET:HG2	1.92	0.52
24:v:482:PHE:O	24:v:486:LYS:NZ	2.40	0.52
25:A:721:A:H2'	25:A:722:A:C8	2.45	0.52
25:A:1069:A:H1'	25:A:1096:A:H4'	1.92	0.52
51:3:26:SER:OG	51:3:27:THR:N	2.42	0.52
56:8:18:LYS:HE2	56:8:21:GLY:HA2	1.91	0.52
1:a:375:U:OP1	16:p:70:ARG:NH1	2.43	0.51
1:a:685:G:H2'	1:a:686:U:C6	2.45	0.51
17:q:28:PHE:CZ	17:q:37:PHE:HB3	2.44	0.51
24:v:21:HIS:HD2	24:v:122:ARG:N	2.08	0.51
25:A:1798:U:OP2	27:C:271:ARG:NH1	2.39	0.51
31:G:84:THR:HG22	31:G:134:LYS:HB3	1.92	0.51
32:I:47:GLU:HG3	32:I:48:ALA:H	1.74	0.51
1:a:1324:A:H2'	1:a:1325:C:C6	2.45	0.51
5:e:153:VAL:HG21	8:h:99:LEU:HD22	1.93	0.51
9:i:130:ARG:NH2	22:x:33:U:OP2	2.43	0.51
10:j:12:ALA:HB3	10:j:18:ILE:HD11	1.93	0.51
19:s:3:ARG:HH21	19:s:7:LYS:HD3	1.75	0.51
19:s:41:PHE:HB2	19:s:44:MET:HE1	1.93	0.51
20:t:31:PHE:O	20:t:35:VAL:HG23	2.09	0.51
31:G:98:VAL:HG11	31:G:124:GLU:HA	1.92	0.51
33:J:14:ALA:HB2	33:J:54:PRO:HG3	1.91	0.51
33:J:101:ILE:O	33:J:141:GLU:N	2.36	0.51
9:i:67:VAL:HG11	9:i:75:GLN:HG2	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:m:17:ILE:O	13:m:20:THR:HG22	2.11	0.51
22:x:26:A:N6	22:x:44:G:N1	2.38	0.51
24:v:390:ILE:HD12	24:v:390:ILE:H	1.75	0.51
28:D:131:ASP:OD1	28:D:132:ALA:N	2.43	0.51
1:a:523:A:N6	12:l:89:D2T:OD1	2.31	0.51
1:a:1014:A:H2'	1:a:1015:G:C4	2.45	0.51
6:f:15:SER:HA	6:f:18:VAL:HG23	1.91	0.51
25:A:881:G:O6	25:A:897:C:N4	2.43	0.51
25:A:1670:C:O2	28:D:134:HIS:HE1	1.94	0.51
25:A:2134:A:N1	25:A:2156:G:H2'	2.25	0.51
30:F:94:GLU:HA	30:F:97:TRP:HB2	1.92	0.51
46:X:86:LEU:HD13	46:X:89:ILE:HD11	1.91	0.51
1:a:1218:C:H2'	1:a:1219:A:H8	1.75	0.51
1:a:1287:A:H2'	1:a:1288:A:C8	2.45	0.51
18:r:32:TYR:CG	18:r:55:LEU:HD21	2.45	0.51
24:v:19:ILE:HD12	24:v:126:LEU:HD12	1.92	0.51
24:v:396:GLU:O	24:v:396:GLU:HG2	2.09	0.51
25:A:1794:A:H2'	25:A:1795:C:C6	2.45	0.51
25:A:2306:C:N4	30:F:39:GLY:O	2.41	0.51
30:F:46:ASP:HB3	30:F:49:LEU:HD23	1.92	0.51
24:v:401:ILE:HG12	24:v:463:TYR:CE1	2.45	0.51
25:A:263:G:O2'	25:A:429:A:N3	2.44	0.51
25:A:546:U:H3'	25:A:547:A:H8	1.76	0.51
25:A:1028:A:H2'	25:A:1029:A:C8	2.46	0.51
25:A:2204:G:OP2	27:C:147:LYS:NZ	2.44	0.51
33:J:49:ILE:HG13	33:J:50:GLU:N	2.24	0.51
1:a:1005:A:OP2	1:a:1024:G:N2	2.28	0.51
1:a:1120:C:H2'	1:a:1121:U:H6	1.76	0.51
7:g:66:LEU:HD13	7:g:70:ARG:HH21	1.76	0.51
7:g:70:ARG:NH2	7:g:97:ASN:OD1	2.44	0.51
8:h:43:GLU:OE1	8:h:43:GLU:N	2.44	0.51
24:v:204:VAL:HG23	24:v:206:ILE:HD11	1.93	0.51
25:A:2804:U:H2'	25:A:2805:C:C6	2.46	0.51
1:a:923:A:O2'	1:a:1399:C:OP2	2.26	0.51
1:a:1147:C:H4'	9:i:7:TYR:CE1	2.46	0.51
12:l:114:ARG:HG3	12:l:119:VAL:HG13	1.93	0.51
24:v:101:ARG:O	24:v:104:THR:HG22	2.09	0.51
1:a:1001:C:N4	1:a:1002:G:O6	2.44	0.51
1:a:1346:A:H61	1:a:1374:A:H3'	1.76	0.51
25:A:1216:G:OP1	41:S:11:ARG:NH1	2.43	0.51
25:A:2154:A:H2'	25:A:2155:U:C6	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
56:8:2:LYS:HG3	56:8:35:GLN:HG2	1.92	0.51
1:a:555:U:H2'	1:a:556:C:C6	2.45	0.51
25:A:1005:C:O2'	34:L:30:THR:HG21	2.10	0.51
25:A:2561:U:O2'	35:M:23:LYS:HG3	2.11	0.51
1:a:56:U:H2'	1:a:57:G:H8	1.76	0.50
9:i:23:PRO:HA	9:i:61:LEU:HD23	1.93	0.50
13:m:5:ALA:HB2	13:m:60:VAL:HG11	1.93	0.50
24:v:250:GLU:OE1	24:v:250:GLU:N	2.44	0.50
25:A:279:A:OP2	25:A:361:G:N2	2.44	0.50
1:a:908:A:H2'	1:a:909:A:H8	1.74	0.50
1:a:1268:G:H2'	1:a:1269:A:C8	2.46	0.50
2:b:101:LEU:HB3	2:b:179:LEU:HD12	1.93	0.50
7:g:13:LEU:HG	7:g:14:PRO:HD2	1.93	0.50
24:v:77:PHE:CG	24:v:77:PHE:O	2.64	0.50
24:v:325:TYR:CG	24:v:326:GLU:N	2.79	0.50
24:v:425:VAL:HG12	24:v:440:ALA:HB2	1.94	0.50
25:A:2184:A:H2'	25:A:2185:U:C6	2.46	0.50
25:A:2484:G:H1'	37:O:123:LYS:HD2	1.93	0.50
1:a:89:U:H2'	1:a:90:C:C6	2.46	0.50
24:v:160:ASN:OD1	24:v:161:GLU:N	2.45	0.50
24:v:260:LEU:H	58:v:601:GCP:C6	2.25	0.50
24:v:317:PHE:CZ	24:v:348:PHE:CZ	2.99	0.50
25:A:1000:A:H2'	25:A:1001:A:C8	2.46	0.50
25:A:1386:C:H2'	25:A:1387:A:C8	2.46	0.50
31:G:127:THR:HB	31:G:130:GLU:HB2	1.91	0.50
33:J:54:PRO:HG2	33:J:78:VAL:HG21	1.93	0.50
1:a:8:A:N6	4:d:202:GLU:O	2.45	0.50
25:A:549:G:H2'	25:A:550:C:C6	2.47	0.50
25:A:2157:G:O2'	25:A:2158:A:O4'	2.25	0.50
1:a:212:G:H2'	1:a:213:G:H8	1.76	0.50
1:a:1412:C:H2'	1:a:1413:A:C8	2.46	0.50
9:i:10:GLY:HA3	9:i:78:ALA:O	2.11	0.50
24:v:319:ARG:O	24:v:320:VAL:C	2.54	0.50
25:A:1061:U:H1'	33:J:11:LEU:HA	1.94	0.50
29:E:146:VAL:HG12	29:E:167:VAL:HG12	1.92	0.50
47:Y:11:ARG:O	47:Y:14:ARG:NH1	2.43	0.50
1:a:219:U:H2'	1:a:220:G:H8	1.77	0.50
24:v:284:THR:OG1	24:v:287:ARG:O	2.24	0.50
25:A:639:U:H2'	25:A:640:C:C6	2.47	0.50
25:A:2071:A:H2'	25:A:2072:C:C6	2.47	0.50
27:C:258:ARG:NH2	27:C:264:ASP:OD1	2.38	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:E:149:ILE:HB	29:E:188:MET:HG2	1.94	0.50
32:I:6:GLN:HA	32:I:9:GLN:HG3	1.94	0.50
3:c:180:ALA:HB1	3:c:203:PHE:HE1	1.76	0.50
25:A:1536:C:H4'	25:A:1537:G:C2	2.47	0.50
2:b:151:ILE:HG13	2:b:154:MET:SD	2.52	0.50
9:i:75:GLN:O	9:i:79:ILE:HG12	2.11	0.50
10:j:10:LEU:HB2	10:j:98:VAL:HG22	1.94	0.50
25:A:624:C:O2'	25:A:657:U:OP1	2.24	0.50
25:A:2114:A:H3'	25:A:2115:G:C8	2.47	0.50
1:a:1217:C:OP2	14:n:9:ARG:NH2	2.45	0.50
4:d:10:LYS:HG2	4:d:38:PRO:HB3	1.93	0.50
7:g:47:LEU:O	7:g:50:LEU:HG	2.12	0.50
11:k:64:GLN:HG2	11:k:95:SER:HB2	1.94	0.50
20:t:28:MET:O	20:t:32:ILE:HG12	2.12	0.50
25:A:1800:C:O2'	27:C:153:GLN:NE2	2.41	0.50
1:a:56:U:H2'	1:a:57:G:C8	2.47	0.49
1:a:1004:A:H8	1:a:1024:G:H21	1.57	0.49
1:a:1006:G:H2'	1:a:1007:U:C6	2.46	0.49
1:a:1530:G:H2'	1:a:1531:A:H8	1.77	0.49
24:v:194:GLN:OE1	24:v:194:GLN:HA	2.11	0.49
25:A:495:G:H5''	43:U:4:ILE:HG23	1.94	0.49
25:A:1111:A:O2'	25:A:1112:G:OP1	2.29	0.49
26:B:42:C:H5	30:F:66:LEU:HD12	1.76	0.49
34:L:96:ARG:NH1	34:L:98:GLU:OE1	2.44	0.49
1:a:216:U:H2'	1:a:217:C:H6	1.77	0.49
6:f:42:TRP:HB2	6:f:59:TYR:HB2	1.95	0.49
7:g:9:GLN:HE22	7:g:11:LYS:HA	1.76	0.49
7:g:116:MET:HA	7:g:119:ARG:HD2	1.94	0.49
11:k:21:ALA:HB3	11:k:84:VAL:HG12	1.93	0.49
22:x:23:A:H2'	22:x:24:G:C8	2.47	0.49
25:A:219:A:N3	25:A:234:U:O2'	2.33	0.49
25:A:1857:G:N2	25:A:1884:G:O2'	2.37	0.49
25:A:2135:A:H5''	25:A:2136:G:N7	2.27	0.49
1:a:1321:U:OP2	1:a:1322:C:O2'	2.23	0.49
5:e:148:ASN:HD22	8:h:96:MET:HE1	1.77	0.49
10:j:21:ALA:O	10:j:24:GLU:HG3	2.12	0.49
20:t:44:LYS:HD2	20:t:87:ALA:HA	1.94	0.49
25:A:700:G:O2'	25:A:1632:A:N3	2.38	0.49
25:A:1992:G:N2	25:A:1996:C:O2'	2.45	0.49
26:B:34:A:N6	26:B:44:G:O2'	2.45	0.49
32:I:38:MET:O	32:I:42:ARG:N	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:W:14:LEU:HD11	45:W:71:ALA:HB2	1.94	0.49
1:a:76:G:H2'	1:a:77:A:C8	2.47	0.49
1:a:401:C:O2'	1:a:621:A:N3	2.42	0.49
1:a:974:A:OP1	14:n:69:ARG:NH2	2.45	0.49
1:a:1002:G:C2	1:a:1003:G:H1'	2.47	0.49
5:e:19:ASN:OD1	5:e:20:ARG:N	2.45	0.49
14:n:49:GLN:NE2	19:s:11:ILE:O	2.43	0.49
25:A:2130:U:H1'	25:A:2159:G:C6	2.46	0.49
47:Y:46:HIS:CE1	47:Y:77:ARG:HD3	2.47	0.49
1:a:208:U:O2'	1:a:211:G:N1	2.36	0.49
25:A:1808:A:H3'	25:A:1809:A:C8	2.48	0.49
25:A:2303:G:H2'	25:A:2304:G:C8	2.48	0.49
1:a:424:G:H2'	1:a:425:G:H8	1.78	0.49
1:a:979:C:O2	14:n:59:ARG:HD2	2.12	0.49
4:d:45:LYS:HG3	4:d:47:ARG:NH2	2.27	0.49
25:A:1263:U:O2'	52:4:8:PRO:HD2	2.13	0.49
40:R:25:THR:HG23	40:R:87:LYS:HB2	1.95	0.49
45:W:50:PRO:HA	45:W:54:GLN:HG3	1.94	0.49
1:a:1004:A:H8	1:a:1024:G:N2	2.11	0.49
1:a:1147:C:H2'	1:a:1148:U:C6	2.48	0.49
2:b:203:ASN:OD1	2:b:205:ASP:N	2.42	0.49
11:k:90:GLY:O	11:k:93:ARG:HG3	2.12	0.49
25:A:491:G:O6	43:U:49:LYS:NZ	2.38	0.49
25:A:2134:A:C6	25:A:2157:G:H1'	2.47	0.49
26:B:41:G:OP1	26:B:43:C:N4	2.39	0.49
1:a:82:G:O2'	1:a:83:C:O4'	2.29	0.49
1:a:1288:A:H2'	1:a:1289:A:H8	1.77	0.49
2:b:90:PHE:HE1	2:b:153:ASP:HB2	1.78	0.49
2:b:109:GLN:HA	2:b:112:LYS:HG2	1.94	0.49
25:A:1009:A:N3	25:A:1153:C:O2'	2.44	0.49
25:A:1111:A:C2	25:A:1112:G:H1'	2.47	0.49
25:A:2134:A:N6	25:A:2156:G:C5	2.81	0.49
30:F:73:SER:OG	30:F:81:GLN:N	2.45	0.49
1:a:80:A:N6	1:a:86:G:H21	2.11	0.49
1:a:1288:A:H2'	1:a:1289:A:C8	2.48	0.49
24:v:14:ARG:HB2	24:v:84:VAL:HG12	1.95	0.49
25:A:898:C:H2'	25:A:899:A:H8	1.78	0.49
25:A:1112:G:H2'	25:A:1113:U:C6	2.48	0.49
36:N:81:ASP:HB3	36:N:100:ILE:HD12	1.94	0.49
48:Z:11:ARG:HD3	48:Z:12:PRO:HD2	1.94	0.49
9:i:51:PRO:HG3	9:i:80:ARG:HG3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:A:612:G:H2'	25:A:614:A:OP2	2.13	0.48
25:A:2303:G:H2'	25:A:2304:G:H8	1.77	0.48
32:I:42:ARG:HH12	33:J:120:ALA:HA	1.78	0.48
32:I:48:ALA:HB3	32:I:51:TYR:CE1	2.48	0.48
4:d:34:ILE:H	4:d:34:ILE:HD12	1.78	0.48
7:g:97:ASN:HB3	7:g:101:MET:HE1	1.95	0.48
13:m:2:ALA:HA	13:m:9:ILE:H	1.78	0.48
24:v:430:PRO:HA	24:v:494:ALA:HB2	1.95	0.48
24:v:433:ASN:OD1	24:v:434:ASN:N	2.47	0.48
25:A:1032:A:OP1	56:8:8:LYS:HG2	2.13	0.48
25:A:1746:A:H2'	25:A:1747:U:C6	2.47	0.48
25:A:2547:A:H2'	25:A:2548:U:C6	2.48	0.48
1:a:1223:C:P	19:s:78:ARG:HH21	2.35	0.48
22:x:55:PSU:HN3	22:x:58:A:P	2.36	0.48
24:v:403:LEU:HD23	24:v:403:LEU:H	1.76	0.48
25:A:2030:6MZ:C2	25:A:2499:C:H5''	2.44	0.48
25:A:2127:G:H1'	25:A:2173:A:H2	1.79	0.48
25:A:2162:G:O2'	25:A:2173:A:N6	2.45	0.48
36:N:132:ARG:HG3	36:N:142:ILE:HD12	1.94	0.48
50:2:10:THR:HG21	50:2:56:LYS:HG3	1.96	0.48
1:a:78:A:H2'	1:a:79:G:C8	2.49	0.48
1:a:562:U:O2'	12:l:13:ALA:O	2.32	0.48
1:a:1141:C:H2'	1:a:1142:G:C8	2.49	0.48
3:c:83:ASP:O	3:c:86:LYS:HG2	2.13	0.48
22:x:60:U:OP1	22:x:61:C:N4	2.41	0.48
25:A:1083:U:H5'	25:A:1084:A:OP2	2.13	0.48
25:A:2804:U:H2'	25:A:2805:C:H6	1.79	0.48
37:O:42:THR:HG22	37:O:45:GLN:HE21	1.78	0.48
1:a:1191:A:OP1	3:c:4:LYS:NZ	2.35	0.48
25:A:1102:C:H2'	25:A:1103:A:C8	2.41	0.48
25:A:1667:G:O2'	25:A:1991:U:O4	2.27	0.48
25:A:2134:A:N6	25:A:2156:G:N3	2.58	0.48
25:A:2154:A:H2'	25:A:2155:U:H6	1.78	0.48
25:A:2444:G:OP2	29:E:63:LYS:HD2	2.12	0.48
46:X:9:ARG:NH2	46:X:17:SER:OG	2.36	0.48
49:1:10:SER:OG	49:1:13:GLU:OE1	2.20	0.48
1:a:1157:A:H5'	1:a:1158:C:C6	2.49	0.48
1:a:1297:G:C6	7:g:114:LYS:HE3	2.47	0.48
7:g:47:LEU:HD12	7:g:50:LEU:HD21	1.95	0.48
25:A:1871:A:OP2	25:A:1871:A:H8	1.96	0.48
31:G:42:GLU:N	31:G:53:GLY:O	2.27	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:M:109:SER:O	35:M:113:MET:HG2	2.14	0.48
45:W:7:ARG:NH1	45:W:26:LYS:O	2.46	0.48
1:a:714:G:H2'	1:a:715:A:C8	2.49	0.48
3:c:86:LYS:HA	3:c:89:LYS:HG2	1.94	0.48
11:k:100:LEU:HD22	11:k:107:ILE:HD11	1.94	0.48
25:A:279:A:N6	25:A:361:G:H1'	2.28	0.48
25:A:593:U:H2'	25:A:594:U:C6	2.49	0.48
25:A:2484:G:O2'	37:O:123:LYS:O	2.30	0.48
26:B:117:G:OP2	39:Q:56:LYS:NZ	2.46	0.48
35:M:65:THR:HG22	35:M:67:LYS:H	1.78	0.48
1:a:90:C:H2'	1:a:91:U:C6	2.49	0.48
1:a:1147:C:H2'	1:a:1148:U:H6	1.78	0.48
1:a:1530:G:H2'	1:a:1531:A:C8	2.48	0.48
24:v:149:ARG:NH2	24:v:157:GLU:OE1	2.47	0.48
28:D:5:VAL:H	28:D:32:ASN:HD21	1.62	0.48
30:F:64:LYS:HE3	51:3:5:ILE:HD12	1.95	0.48
1:a:1125:U:O2'	1:a:1126:U:O5'	2.29	0.48
1:a:1323:G:H2'	1:a:1324:A:C8	2.49	0.48
2:b:134:ALA:O	2:b:138:THR:HG23	2.14	0.48
7:g:69:VAL:HG13	7:g:100:ALA:HB1	1.95	0.48
10:j:10:LEU:HD11	10:j:22:THR:HG23	1.95	0.48
25:A:364:C:H2'	25:A:365:U:C6	2.48	0.48
25:A:644:A:H2'	25:A:645:C:O4'	2.14	0.48
25:A:1476:U:H2'	25:A:1477:A:H8	1.78	0.48
1:a:1372:U:H2'	1:a:1373:G:O4'	2.14	0.48
25:A:1386:C:H2'	25:A:1387:A:H8	1.79	0.48
25:A:2172:U:O4'	25:A:2175:C:N4	2.33	0.48
1:a:475:C:H2'	1:a:476:U:H6	1.79	0.47
1:a:500:G:H2'	1:a:501:C:C6	2.49	0.47
2:b:58:ASN:HB3	2:b:220:THR:HG22	1.96	0.47
2:b:207:ILE:O	2:b:211:THR:HG23	2.14	0.47
6:f:2:ARG:NH1	6:f:68:GLN:OE1	2.45	0.47
9:i:42:GLU:HA	9:i:45:ARG:HD2	1.96	0.47
22:x:15:G:N1	22:x:48:C:C2	2.82	0.47
24:v:378:THR:O	24:v:378:THR:OG1	2.31	0.47
25:A:191:A:H2'	25:A:192:C:H6	1.79	0.47
25:A:1203:U:O2'	36:N:4:ASN:ND2	2.44	0.47
1:a:84:U:N3	1:a:87:C:O2	2.47	0.47
1:a:1254:A:H2'	1:a:1255:G:C8	2.50	0.47
7:g:115:SER:H	7:g:118:LEU:HD21	1.79	0.47
24:v:458:ASN:O	24:v:459:VAL:HG22	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:F:142:ASP:OD1	30:F:142:ASP:N	2.46	0.47
37:O:42:THR:HG23	37:O:45:GLN:HG3	1.95	0.47
49:1:39:GLN:HB3	49:1:41:HIS:CE1	2.49	0.47
1:a:954:G:H2'	1:a:955:U:C6	2.48	0.47
1:a:1226:C:H2'	13:m:102:THR:HG22	1.96	0.47
1:a:1348:U:H2'	1:a:1349:A:H8	1.80	0.47
3:c:112:ASP:OD1	3:c:115:LEU:HB2	2.14	0.47
3:c:123:GLN:HE22	3:c:136:ARG:HD2	1.79	0.47
4:d:146:ARG:HG2	4:d:147:GLU:H	1.78	0.47
8:h:5:ASP:OD1	8:h:77:ARG:NH1	2.46	0.47
22:x:47:U:N3	22:x:50:U:OP1	2.46	0.47
24:v:284:THR:HG22	24:v:387:PHE:CD2	2.49	0.47
24:v:363:ASP:OD1	24:v:364:ILE:N	2.47	0.47
25:A:895:U:H2'	25:A:897:C:C6	2.49	0.47
25:A:1060:U:H3	25:A:1088:A:H1'	1.79	0.47
37:O:36:VAL:HG12	46:X:82:TYR:HB2	1.96	0.47
1:a:460:A:H2'	1:a:461:A:C8	2.49	0.47
1:a:1179:A:H5''	9:i:99:ARG:HH12	1.79	0.47
2:b:60:ILE:HA	2:b:63:ARG:NH2	2.28	0.47
5:e:55:GLU:CD	5:e:57:PRO:HD2	2.40	0.47
25:A:545:U:H4'	25:A:546:U:OP2	2.14	0.47
25:A:1071:G:H8	25:A:1071:G:OP1	1.96	0.47
29:E:165:HIS:CE1	29:E:166:LYS:HG3	2.49	0.47
33:J:99:GLY:HA3	33:J:138:LEU:HD23	1.95	0.47
2:b:7:ARG:HH12	2:b:11:LYS:HB2	1.79	0.47
3:c:180:ALA:HB1	3:c:203:PHE:CE1	2.50	0.47
15:o:15:PHE:HE2	15:o:26:GLU:HB3	1.80	0.47
25:A:621:A:OP2	36:N:99:ASN:ND2	2.48	0.47
25:A:1734:G:H2'	25:A:1735:A:H8	1.80	0.47
1:a:1239:A:N7	1:a:1298:U:N3	2.62	0.47
3:c:55:ILE:HG23	3:c:68:ILE:HG22	1.96	0.47
6:f:10:VAL:CG1	6:f:58:HIS:HB3	2.44	0.47
17:q:7:THR:OG1	17:q:60:GLU:OE2	2.31	0.47
24:v:102:THR:HG22	24:v:301:PHE:CE1	2.49	0.47
25:A:1040:A:H2	25:A:1115:G:H1	1.62	0.47
37:O:2:LEU:HD12	37:O:68:PHE:CE2	2.49	0.47
1:a:1292:G:H2'	1:a:1293:C:C6	2.49	0.47
2:b:111:ILE:HD11	2:b:151:ILE:HG23	1.96	0.47
4:d:188:ARG:NH2	4:d:197:GLU:OE2	2.48	0.47
19:s:4:SER:HB2	19:s:7:LYS:HD2	1.95	0.47
24:v:18:ILE:HG23	24:v:18:ILE:O	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:v:19:ILE:HG22	24:v:89:THR:OG1	2.15	0.47
24:v:316:ALA:O	24:v:366:GLY:HA2	2.14	0.47
24:v:419:LEU:HD11	24:v:452:ARG:HD2	1.95	0.47
25:A:875:G:H2'	25:A:876:C:C6	2.50	0.47
25:A:2030:6MZ:H2	25:A:2499:C:H5''	1.97	0.47
25:A:2126:A:O2'	25:A:2173:A:N1	2.46	0.47
25:A:2172:U:O2'	25:A:2174:C:OP1	2.22	0.47
25:A:2756:U:OP2	56:8:19:ARG:NE	2.47	0.47
26:B:52:A:N7	39:Q:33:ARG:NH1	2.63	0.47
30:F:134:GLU:OE1	30:F:136:ILE:HG12	2.15	0.47
30:F:148:ARG:HD3	30:F:149:VAL:N	2.29	0.47
1:a:1151:A:O2'	1:a:1152:A:H8	1.98	0.47
24:v:482:PHE:O	24:v:486:LYS:HG2	2.15	0.47
25:A:1063:G:H21	33:J:77:ALA:HA	1.79	0.47
25:A:2554:U:H2'	25:A:2555:U:H6	1.79	0.47
56:8:14:CYS:HA	56:8:27:CYS:HB3	1.96	0.47
1:a:1263:C:H2'	1:a:1264:U:C6	2.50	0.47
2:b:60:ILE:HD13	2:b:160:ALA:HB2	1.96	0.47
24:v:93:GLU:OE1	24:v:122:ARG:NH2	2.48	0.47
25:A:709:U:H2'	25:A:710:U:C6	2.49	0.47
25:A:2303:G:O2'	30:F:121:SER:O	2.32	0.47
50:2:12:SER:HB2	50:2:32:ILE:HD11	1.97	0.47
50:2:40:ASP:OD1	50:2:45:ARG:NE	2.38	0.47
1:a:1141:C:H2'	1:a:1142:G:H8	1.80	0.47
24:v:433:ASN:CG	24:v:434:ASN:N	2.72	0.47
24:v:473:TRP:NE1	24:v:496:ASP:OD2	2.46	0.47
25:A:277:G:O2'	25:A:278:A:O5'	2.25	0.47
25:A:851:C:H2'	25:A:852:U:H6	1.79	0.47
25:A:2728:U:HO2'	25:A:2729:G:H8	1.61	0.47
32:I:36:ASP:O	32:I:39:THR:OG1	2.26	0.47
45:W:98:SER:O	45:W:98:SER:OG	2.33	0.47
1:a:634:C:H2'	1:a:635:A:H8	1.80	0.46
1:a:821:G:H2'	1:a:822:U:C6	2.50	0.46
1:a:1228:C:OP1	13:m:107:ARG:NH2	2.48	0.46
2:b:119:THR:O	2:b:123:ASP:CB	2.55	0.46
24:v:331:LEU:HD23	24:v:379:PHE:HB3	1.96	0.46
25:A:645:C:H2'	25:A:647:G:N7	2.30	0.46
25:A:2189:U:H2'	25:A:2190:G:H8	1.80	0.46
25:A:2192:U:H2'	25:A:2193:G:H8	1.80	0.46
25:A:2849:U:H4'	25:A:2868:A:C2	2.50	0.46
27:C:117:GLN:H	27:C:128:ASN:ND2	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:J:93:PRO:HB3	33:J:135:SER:O	2.15	0.46
36:N:78:ARG:HG2	36:N:113:ALA:HB3	1.97	0.46
39:Q:28:VAL:HG12	39:Q:93:ASP:O	2.14	0.46
51:3:34:LEU:HD23	51:3:34:LEU:H	1.79	0.46
1:a:405:U:OP2	4:d:3:ARG:NH1	2.48	0.46
1:a:1006:G:H5'	14:n:19:LYS:NZ	2.30	0.46
1:a:1019:A:H2'	1:a:1020:G:C8	2.51	0.46
24:v:492:GLN:O	24:v:492:GLN:HG2	2.15	0.46
25:A:225:C:H2'	25:A:226:A:O4'	2.14	0.46
25:A:1432:G:H2'	25:A:1433:A:C8	2.50	0.46
27:C:68:LYS:HD2	27:C:70:ASN:HD21	1.79	0.46
50:2:24:LEU:HD21	50:2:54:MET:HE3	1.96	0.46
1:a:21:G:H2'	1:a:22:G:C8	2.51	0.46
1:a:476:U:H2'	1:a:477:C:C6	2.50	0.46
1:a:1524:C:H2'	1:a:1525:G:C8	2.51	0.46
4:d:105:MET:CE	4:d:180:GLY:HA3	2.46	0.46
7:g:17:LYS:HE2	7:g:44:TYR:CD1	2.50	0.46
11:k:60:PRO:O	11:k:64:GLN:HG3	2.15	0.46
13:m:40:ALA:HB3	13:m:43:VAL:HG23	1.96	0.46
25:A:102:U:O4	49:1:4:LYS:HG3	2.15	0.46
25:A:1281:G:H2'	25:A:1282:U:C6	2.51	0.46
25:A:2100:G:C6	25:A:2190:G:C6	3.04	0.46
25:A:2514:U:H2'	25:A:2515:C:C6	2.50	0.46
45:W:5:ILE:O	45:W:5:ILE:HG13	2.16	0.46
1:a:4:U:H2'	1:a:5:U:H5''	1.96	0.46
1:a:539:A:H2'	1:a:540:G:C8	2.50	0.46
1:a:674:G:H2'	1:a:675:A:C8	2.49	0.46
1:a:1277:C:O2'	1:a:1279:G:N3	2.32	0.46
2:b:87:CYS:SG	2:b:221:VAL:HG11	2.55	0.46
4:d:161:LEU:HA	4:d:164:GLN:NE2	2.30	0.46
8:h:47:GLU:HG2	8:h:64:LYS:HB3	1.98	0.46
24:v:127:MET:O	24:v:131:ARG:HB2	2.15	0.46
24:v:186:TYR:N	24:v:186:TYR:CD1	2.79	0.46
24:v:347:THR:N	24:v:354:SER:O	2.47	0.46
33:J:106:LEU:HB2	33:J:129:ILE:CG2	2.45	0.46
33:J:114:ALA:HA	33:J:117:MET:HB3	1.98	0.46
11:k:19:GLY:O	11:k:37:ARG:NH2	2.48	0.46
11:k:107:ILE:HG21	11:k:110:ILE:HD11	1.98	0.46
24:v:130:THR:OG1	24:v:135:THR:OG1	2.33	0.46
25:A:1084:A:OP1	32:I:56:ARG:N	2.37	0.46
35:M:63:VAL:HG12	35:M:107:LEU:HD11	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:a:269:C:N4	1:a:270:A:H62	2.14	0.46
1:a:1313:U:H2'	1:a:1314:C:H6	1.80	0.46
5:e:55:GLU:OE1	5:e:57:PRO:HD2	2.15	0.46
6:f:37:HIS:HB3	6:f:97:THR:HG22	1.97	0.46
6:f:38:ARG:HB3	6:f:63:ASN:OD1	2.15	0.46
13:m:69:LEU:O	13:m:73:ILE:HD12	2.16	0.46
25:A:568:U:H1'	25:A:2030:6MZ:H9C1	1.97	0.46
25:A:642:U:O2'	25:A:644:A:N7	2.38	0.46
25:A:1412:U:H2'	25:A:1413:A:C8	2.51	0.46
25:A:1469:A:H2'	25:A:1470:A:H8	1.80	0.46
33:J:102:SER:OG	33:J:105:GLN:HB2	2.16	0.46
46:X:76:ASP:OD1	46:X:77:VAL:N	2.49	0.46
1:a:139:A:H2'	1:a:140:U:H6	1.81	0.46
1:a:555:U:H2'	1:a:556:C:H6	1.80	0.46
1:a:728:A:H2'	1:a:729:A:C8	2.51	0.46
1:a:1376:U:O4	7:g:10:ARG:NH1	2.49	0.46
2:b:43:LEU:HA	2:b:46:THR:HG22	1.97	0.46
3:c:36:ASP:O	3:c:39:VAL:HG12	2.16	0.46
24:v:307:MET:O	24:v:307:MET:HG2	2.15	0.46
25:A:574:A:N6	25:A:2034:U:OP1	2.49	0.46
25:A:1067:A:H2'	25:A:1067:A:N3	2.30	0.46
25:A:2820:A:O2'	25:A:2821:A:OP1	2.27	0.46
26:B:39:A:C2	26:B:44:G:C2	3.03	0.46
32:I:70:GLU:HG3	32:I:71:CYS:H	1.80	0.46
1:a:409:U:H2'	1:a:410:G:O4'	2.15	0.46
1:a:947:G:H2'	1:a:948:C:C6	2.51	0.46
1:a:987:G:H21	1:a:1015:G:N2	2.13	0.46
1:a:1055:A:O2'	3:c:161:GLU:O	2.31	0.46
6:f:29:ILE:HG12	6:f:64:VAL:HG21	1.98	0.46
25:A:589:U:H2'	25:A:590:A:C8	2.51	0.46
25:A:2136:G:H22	25:A:2155:U:H3	1.64	0.46
1:a:246:A:C2	1:a:282:A:C5	3.04	0.46
6:f:90:MET:HE2	6:f:90:MET:HB3	1.72	0.46
24:v:83:LEU:O	24:v:83:LEU:HD23	2.15	0.46
24:v:419:LEU:HD23	24:v:419:LEU:O	2.15	0.46
25:A:359:G:H2'	25:A:360:U:H6	1.81	0.46
25:A:373:U:H2'	25:A:374:A:H8	1.81	0.46
25:A:1077:A:O2'	25:A:1088:A:N6	2.47	0.46
25:A:1180:U:H2'	25:A:1181:U:O4'	2.15	0.46
25:A:1794:A:H2'	25:A:1795:C:H6	1.81	0.46
25:A:2176:A:O2'	25:A:2177:C:O4'	2.21	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:F:46:ASP:OD2	30:F:49:LEU:N	2.44	0.46
40:R:25:THR:HG22	40:R:88:ARG:HG2	1.96	0.46
1:a:380:G:N2	1:a:383:A:OP2	2.47	0.46
1:a:1002:G:N2	1:a:1003:G:H1'	2.31	0.46
1:a:1219:A:H5''	14:n:53:ARG:NH1	2.31	0.46
3:c:57:ILE:HD12	3:c:66:VAL:HG22	1.97	0.46
25:A:586:A:N1	25:A:809:G:O2'	2.47	0.46
32:I:103:ASN:HA	32:I:107:GLU:HB3	1.98	0.46
2:b:9:MET:HE2	2:b:14:VAL:HG11	1.98	0.45
9:i:28:ILE:HG12	9:i:63:LEU:HD12	1.97	0.45
24:v:430:PRO:HG2	24:v:435:ASP:HB3	1.98	0.45
25:A:993:G:H1'	42:T:91:GLN:NE2	2.31	0.45
25:A:2532:G:N2	25:A:2663:G:O2'	2.49	0.45
25:A:2683:C:H4'	28:D:13:ARG:NH2	2.31	0.45
28:D:2:ILE:HD11	28:D:84:LEU:HD12	1.98	0.45
30:F:106:ILE:HG22	51:3:34:LEU:HD13	1.98	0.45
32:I:42:ARG:NH1	33:J:120:ALA:HA	2.31	0.45
37:O:1:MET:HE1	37:O:47:GLU:HG2	1.97	0.45
1:a:55:A:C6	24:v:311:HIS:CD2	3.04	0.45
1:a:883:C:O2'	1:a:884:U:H5'	2.16	0.45
1:a:1315:U:H2'	1:a:1316:G:C8	2.51	0.45
4:d:65:TYR:OH	4:d:95:GLU:OE2	2.20	0.45
11:k:87:LYS:HG3	11:k:113:VAL:HG23	1.98	0.45
22:x:45:U:H5'	22:x:46:G7M:OP2	2.16	0.45
24:v:295:LYS:HA	24:v:381:GLN:HE21	1.81	0.45
25:A:1353:A:H2'	25:A:1354:A:C8	2.52	0.45
25:A:2126:A:O3'	25:A:2162:G:N2	2.41	0.45
27:C:245:VAL:HG12	27:C:251:GLN:HA	1.98	0.45
51:3:41:HIS:CD2	51:3:43:PHE:HB3	2.51	0.45
54:6:3:ARG:HD3	54:6:3:ARG:HA	1.75	0.45
1:a:333:U:OP1	20:t:2:ALA:N	2.49	0.45
9:i:8:GLY:HA3	9:i:86:ALA:HB2	1.98	0.45
11:k:33:THR:HG22	11:k:44:TRP:HB3	1.98	0.45
24:v:476:CYS:SG	24:v:520:ASP:HA	2.56	0.45
25:A:1060:U:C2	25:A:1088:A:H8	2.34	0.45
25:A:2820:A:N3	25:A:2820:A:H2'	2.32	0.45
56:8:16:ILE:HD13	56:8:25:VAL:HG22	1.98	0.45
1:a:1021:A:N1	1:a:1022:A:N6	2.64	0.45
1:a:1221:G:H2'	1:a:1222:G:C8	2.51	0.45
22:x:42:C:H2'	22:x:43:C:C6	2.52	0.45
25:A:1101:U:H2'	25:A:1102:C:O4'	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:A:2097:A:O2'	25:A:2098:U:OP1	2.32	0.45
25:A:2139:U:C2	25:A:2152:G:O6	2.70	0.45
25:A:2680:U:O2'	25:A:2681:C:H5'	2.16	0.45
36:N:10:GLU:OE2	36:N:10:GLU:N	2.45	0.45
37:O:136:MET:HB3	46:X:57:TYR:CD2	2.51	0.45
1:a:492:C:H2'	1:a:493:A:C8	2.52	0.45
2:b:68:LEU:HD12	2:b:90:PHE:HB2	1.98	0.45
2:b:144:LEU:O	2:b:148:LEU:HD23	2.16	0.45
7:g:51:ALA:HB2	7:g:58:GLU:HB3	1.99	0.45
11:k:126:LYS:HA	21:u:37:PHE:CD1	2.51	0.45
1:a:519:C:OP2	12:l:47:SER:OG	2.25	0.45
2:b:42:ASN:ND2	2:b:45:LYS:HG2	2.32	0.45
2:b:54:LEU:HD12	2:b:220:THR:OG1	2.16	0.45
6:f:2:ARG:HG2	6:f:91:ARG:HH21	1.82	0.45
25:A:1077:A:H5'	33:J:93:PRO:HD2	1.98	0.45
31:G:52:PHE:CE2	31:G:69:ARG:HA	2.51	0.45
1:a:137:U:O2	1:a:226:G:N2	2.41	0.45
1:a:359:G:H5''	24:v:312:ARG:NH2	2.32	0.45
1:a:390:U:H2'	1:a:391:G:H8	1.81	0.45
5:e:153:VAL:HA	5:e:156:LYS:HG2	1.99	0.45
25:A:248:G:H5'	25:A:250:G:N7	2.31	0.45
26:B:44:G:H5''	26:B:45:A:OP1	2.17	0.45
26:B:49:C:OP1	39:Q:102:ARG:HG2	2.16	0.45
37:O:63:ILE:HD13	37:O:105:MET:HG3	1.98	0.45
1:a:1144:G:N2	1:a:1146:A:H62	2.15	0.45
1:a:1435:G:H2'	1:a:1436:U:C6	2.52	0.45
6:f:14:GLN:NE2	6:f:17:GLN:HE22	2.14	0.45
7:g:68:ASN:ND2	7:g:130:ASN:HD21	2.15	0.45
22:x:15:G:N2	22:x:48:C:C4	2.85	0.45
24:v:360:TYR:HB3	24:v:361:PRO:HD2	1.98	0.45
24:v:400:ARG:NE	24:v:435:ASP:OD1	2.49	0.45
25:A:1060:U:H1'	25:A:1062:G:N3	2.32	0.45
25:A:1062:G:N7	25:A:1064:C:N4	2.64	0.45
25:A:1730:C:O2'	25:A:1731:G:O5'	2.32	0.45
25:A:2020:A:H5'	52:4:9:THR:HG21	1.99	0.45
25:A:2898:U:H2'	25:A:2899:A:H8	1.81	0.45
30:F:60:ILE:HA	30:F:140:GLU:HG3	1.99	0.45
31:G:33:LEU:HB2	31:G:75:MET:HE3	1.99	0.45
31:G:44:LYS:HB2	31:G:51:THR:OG1	2.17	0.45
31:G:90:VAL:HG12	31:G:160:LYS:HA	1.98	0.45
33:J:56:PRO:HD3	33:J:75:PRO:HD3	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:a:509:A:N3	1:a:543:U:O2'	2.46	0.45
1:a:954:G:H2'	1:a:955:U:H6	1.82	0.45
1:a:994:A:C5	1:a:1216:A:H4'	2.52	0.45
1:a:1347:G:C8	9:i:109:ARG:HB3	2.51	0.45
2:b:7:ARG:NH1	2:b:11:LYS:HB2	2.32	0.45
3:c:79:LYS:O	3:c:80:LYS:HG2	2.17	0.45
4:d:172:GLU:HG3	4:d:183:LYS:HD2	1.98	0.45
10:j:59:LYS:HD2	10:j:62:ARG:NH2	2.25	0.45
13:m:90:ARG:HA	13:m:90:ARG:HD2	1.77	0.45
25:A:2131:U:H4'	25:A:2133:G:H4'	1.99	0.45
25:A:2230:G:H2'	25:A:2231:U:C6	2.52	0.45
27:C:30:PHE:CD2	27:C:33:LEU:HD12	2.50	0.45
30:F:144:ASP:OD1	30:F:144:ASP:N	2.50	0.45
33:J:76:ALA:HB2	33:J:113:LYS:HE3	1.98	0.45
42:T:14:VAL:HG11	42:T:20:VAL:HG21	1.98	0.45
1:a:33:A:H2'	1:a:34:C:C6	2.52	0.45
1:a:412:A:O2'	1:a:413:G:H4'	2.16	0.45
1:a:1254:A:H2'	1:a:1255:G:H8	1.82	0.45
3:c:36:ASP:HA	3:c:39:VAL:HG12	1.97	0.45
3:c:58:GLU:OE2	10:j:94:ALA:HB1	2.17	0.45
3:c:91:VAL:HA	3:c:94:ILE:HG22	1.99	0.45
5:e:82:GLN:HE21	5:e:149:SER:HA	1.82	0.45
19:s:55:ARG:HG3	19:s:56:GLN:HG2	1.98	0.45
24:v:481:LYS:HA	24:v:481:LYS:HD2	1.72	0.45
25:A:2312:U:O3'	30:F:68:THR:HG21	2.17	0.45
27:C:121:ASP:OD1	27:C:121:ASP:N	2.48	0.45
8:h:54:ASP:OD1	8:h:55:THR:N	2.49	0.44
19:s:19:VAL:O	19:s:23:VAL:HG23	2.17	0.44
25:A:247:G:OP2	25:A:249:C:N4	2.47	0.44
25:A:357:C:H2'	25:A:358:U:C6	2.52	0.44
25:A:849:A:H2'	25:A:850:U:H6	1.81	0.44
25:A:1720:U:H2'	25:A:1721:G:O4'	2.16	0.44
25:A:2123:G:H1'	25:A:2176:A:C2	2.52	0.44
33:J:106:LEU:HD12	33:J:107:GLN:N	2.32	0.44
34:L:49:ASP:CG	34:L:121:LYS:HZ3	2.25	0.44
1:a:1243:C:H2'	1:a:1244:G:H8	1.82	0.44
2:b:57:LEU:O	2:b:60:ILE:HG22	2.17	0.44
6:f:29:ILE:HD12	6:f:29:ILE:H	1.81	0.44
21:u:39:GLU:HG2	21:u:43:THR:OG1	2.17	0.44
25:A:2074:U:H2'	25:A:2075:U:C6	2.53	0.44
39:Q:81:ARG:O	39:Q:84:GLU:HG2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
53:5:10:LYS:HE3	53:5:53:LYS:O	2.16	0.44
1:a:1220:G:H1'	19:s:52:HIS:CE1	2.53	0.44
5:e:34:THR:HG22	5:e:52:LYS:HG3	1.99	0.44
7:g:108:ALA:O	7:g:119:ARG:NH1	2.51	0.44
8:h:9:ASP:O	8:h:13:ARG:HG3	2.17	0.44
12:l:66:TYR:HB2	12:l:93:VAL:HG11	2.00	0.44
22:x:15:G:N2	22:x:48:C:H42	2.14	0.44
24:v:425:VAL:HG11	24:v:449:VAL:HG21	1.99	0.44
25:A:1038:G:H2'	25:A:1039:A:C8	2.52	0.44
25:A:1223:G:OP1	42:T:68:ARG:NH1	2.50	0.44
25:A:1857:G:O2'	25:A:1884:G:N2	2.45	0.44
33:J:79:LEU:HD23	33:J:80:LEU:HD23	1.99	0.44
38:P:9:GLN:HG3	38:P:17:ARG:HH12	1.83	0.44
38:P:38:LEU:HB3	38:P:39:PRO:HD3	1.99	0.44
1:a:26:A:H61	1:a:558:G:H1'	1.83	0.44
1:a:1107:C:C4	1:a:1108:G:C8	3.05	0.44
1:a:1411:C:H2'	1:a:1412:C:C6	2.53	0.44
5:e:41:ASP:OD1	5:e:41:ASP:N	2.48	0.44
8:h:96:MET:SD	8:h:130:ALA:HB1	2.58	0.44
24:v:37:GLY:HA3	24:v:77:PHE:HD2	1.82	0.44
25:A:1925:C:H2'	25:A:1926:U:H2'	1.99	0.44
25:A:2900:A:H2'	25:A:2901:C:C6	2.53	0.44
29:E:159:LEU:HD22	29:E:162:ARG:HH21	1.82	0.44
32:I:37:LYS:HZ3	32:I:56:ARG:HH21	1.63	0.44
43:U:88:ARG:HA	43:U:88:ARG:HD2	1.87	0.44
44:V:25:GLU:OE1	44:V:25:GLU:N	2.50	0.44
1:a:821:G:H2'	1:a:822:U:H6	1.83	0.44
24:v:300:VAL:HA	24:v:317:PHE:O	2.18	0.44
25:A:2149:U:H2'	25:A:2150:C:C6	2.53	0.44
30:F:135:GLN:NE2	30:F:149:VAL:HA	2.33	0.44
32:I:4:ASN:HA	32:I:7:ASP:HB2	1.98	0.44
45:W:81:ASP:OD2	45:W:98:SER:OG	2.34	0.44
1:a:1004:A:C2	1:a:1026:G:H1'	2.52	0.44
4:d:117:LEU:CD2	4:d:123:ILE:HD11	2.42	0.44
9:i:12:ARG:NE	9:i:107:ASP:OD2	2.51	0.44
22:x:15:G:O6	22:x:48:C:O2	2.36	0.44
25:A:878:A:N6	25:A:899:A:H1'	2.33	0.44
27:C:145:GLU:HB2	27:C:188:CYS:HB3	1.99	0.44
28:D:4:LEU:HD21	28:D:96:ILE:HG22	2.00	0.44
33:J:101:ILE:HG22	33:J:140:VAL:HA	2.00	0.44
39:Q:26:LEU:HD23	39:Q:92:PHE:HD1	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:U:38:TYR:CE2	52:4:28:LEU:HD21	2.53	0.44
1:a:35:G:H2'	1:a:36:C:H6	1.81	0.44
1:a:713:G:H2'	1:a:714:G:C8	2.53	0.44
1:a:753:A:OP1	15:o:69:TYR:OH	2.24	0.44
1:a:1265:C:N4	1:a:1271:A:H61	2.16	0.44
7:g:115:SER:OG	7:g:116:MET:N	2.51	0.44
22:x:2:C:H2'	22:x:3:C:C6	2.52	0.44
24:v:12:LYS:HG3	24:v:12:LYS:O	2.18	0.44
24:v:77:PHE:CE1	24:v:86:LEU:HD23	2.45	0.44
25:A:636:G:N1	36:N:76:GLU:OE1	2.42	0.44
25:A:645:C:H2'	25:A:647:G:C8	2.53	0.44
33:J:8:TYR:HB2	33:J:59:ILE:HG13	1.99	0.44
1:a:323:U:H2'	1:a:324:G:O4'	2.17	0.44
1:a:950:U:H2'	1:a:951:G:C8	2.53	0.44
2:b:49:MET:HE3	2:b:199:VAL:O	2.18	0.44
16:p:14:ARG:HE	16:p:42:ILE:HD12	1.83	0.44
25:A:155:A:H2'	25:A:156:A:C8	2.52	0.44
25:A:703:U:H2'	25:A:704:G:O4'	2.17	0.44
25:A:1465:G:H2'	25:A:1466:U:O4'	2.17	0.44
25:A:1790:C:H2'	25:A:1791:A:C8	2.52	0.44
25:A:2745:C:C4	25:A:2746:U:C4	3.05	0.44
33:J:57:VAL:HG23	33:J:71:THR:HB	2.00	0.44
40:R:13:MET:HE3	40:R:13:MET:HB3	1.78	0.44
1:a:131:A:H2'	1:a:132:C:C6	2.52	0.44
1:a:820:U:H4'	1:a:821:G:OP2	2.18	0.44
1:a:1121:U:C2	1:a:1122:U:C5	3.06	0.44
1:a:1316:G:N1	1:a:1319:A:OP2	2.50	0.44
9:i:19:VAL:HA	9:i:65:ILE:HG22	1.98	0.44
16:p:18:GLN:OE1	16:p:35:ARG:NE	2.50	0.44
24:v:329:MET:O	24:v:341:VAL:HA	2.18	0.44
24:v:429:ARG:O	24:v:494:ALA:HA	2.17	0.44
25:A:2285:C:OP2	53:5:6:ARG:HD3	2.18	0.44
25:A:2298:A:OP1	30:F:71:ARG:NH1	2.51	0.44
1:a:296:U:O2'	1:a:556:C:O2	2.34	0.43
1:a:337:G:H2'	1:a:338:A:H8	1.81	0.43
10:j:6:ILE:HD11	10:j:79:PRO:HB3	2.00	0.43
13:m:77:ILE:O	13:m:81:MET:HG2	2.18	0.43
24:v:101:ARG:C	24:v:103:LEU:N	2.75	0.43
24:v:185:LEU:C	24:v:186:TYR:HD1	2.26	0.43
24:v:520:ASP:CG	24:v:522:GLN:HG2	2.43	0.43
25:A:673:C:OP1	29:E:49:ARG:NH1	2.50	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:A:2139:U:H2'	25:A:2140:G:H8	1.84	0.43
25:A:2291:U:OP1	25:A:2380:C:O2'	2.36	0.43
25:A:2306:C:OP2	25:A:2307:G:O2'	2.27	0.43
25:A:2500:U:O2	25:A:2504:PSU:N1	2.51	0.43
25:A:2516:A:O2'	25:A:2517:C:H5'	2.18	0.43
46:X:1:MET:HE1	46:X:59:GLU:CD	2.43	0.43
1:a:83:C:O2	1:a:85:U:H5''	2.18	0.43
1:a:436:C:H2'	1:a:437:U:C6	2.52	0.43
1:a:472:U:H2'	1:a:473:U:H6	1.83	0.43
1:a:1126:U:OP1	10:j:7:ARG:NH2	2.47	0.43
1:a:1296:C:H4'	1:a:1302:C:H41	1.82	0.43
5:e:151:GLU:CD	5:e:151:GLU:H	2.26	0.43
7:g:84:THR:O	7:g:86:GLN:NE2	2.45	0.43
8:h:113:ASP:N	8:h:113:ASP:OD1	2.47	0.43
24:v:21:HIS:CD2	24:v:120:GLU:CG	2.96	0.43
48:Z:32:ASN:OD1	48:Z:34:HIS:NE2	2.51	0.43
1:a:74:A:H2'	1:a:75:G:C8	2.53	0.43
2:b:109:GLN:HB3	2:b:112:LYS:HZ3	1.84	0.43
5:e:94:VAL:HB	5:e:111:MET:HE2	2.00	0.43
8:h:25:VAL:HG12	8:h:27:MET:HE2	2.00	0.43
25:A:869:G:O3'	37:O:6:ARG:NH1	2.51	0.43
25:A:1363:C:O2'	25:A:1809:A:N3	2.43	0.43
25:A:2745:C:H2'	25:A:2746:U:C6	2.53	0.43
30:F:50:LEU:HD12	30:F:53:ALA:HB3	1.99	0.43
40:R:64:ILE:HA	40:R:69:GLY:HA2	1.99	0.43
1:a:107:G:N7	20:t:10:ARG:NH2	2.65	0.43
1:a:1143:G:C2	1:a:1144:G:C5	3.06	0.43
1:a:1409:C:H2'	1:a:1410:A:H8	1.83	0.43
2:b:31:ILE:HG21	2:b:39:HIS:HD2	1.82	0.43
17:q:65:ARG:HB2	17:q:66:PRO:HD2	2.00	0.43
22:x:15:G:N2	22:x:20:U:O4	2.48	0.43
25:A:1360:G:N7	25:A:1361:G:C8	2.87	0.43
25:A:2109:U:H1'	25:A:2181:U:O2	2.17	0.43
2:b:45:LYS:HA	2:b:45:LYS:HD2	1.84	0.43
22:x:55:PSU:N3	22:x:58:A:OP2	2.49	0.43
24:v:240:GLU:OE1	24:v:240:GLU:N	2.49	0.43
25:A:299:A:N3	25:A:319:G:O2'	2.40	0.43
25:A:1405:U:H2'	25:A:1406:U:C6	2.53	0.43
25:A:1724:G:O6	25:A:1737:G:N2	2.52	0.43
51:3:15:SER:HB2	51:3:33:ASN:HD21	1.83	0.43
53:5:37:LYS:HB2	53:5:37:LYS:HE3	1.85	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:a:254:G:H21	17:q:18:GLU:CD	2.26	0.43
1:a:1062:U:H2'	1:a:1063:C:C6	2.53	0.43
1:a:1401:G:H2'	1:a:1402:4OC:O4'	2.19	0.43
10:j:64:GLN:HE22	14:n:101:TRP:HH2	1.66	0.43
11:k:45:ALA:HB3	11:k:70:CYS:HB2	2.00	0.43
22:x:21:A:N6	22:x:46:G7M:H1'	2.32	0.43
24:v:180:LYS:HA	24:v:180:LYS:HD2	1.66	0.43
25:A:3:U:H2'	25:A:4:U:C6	2.54	0.43
25:A:359:G:H2'	25:A:360:U:C6	2.54	0.43
25:A:1054:A:O2'	32:I:31:ARG:HA	2.18	0.43
25:A:1673:G:O6	28:D:134:HIS:HD2	2.02	0.43
25:A:2788:C:H2'	25:A:2789:C:C6	2.53	0.43
29:E:21:ARG:HH11	29:E:106:LYS:HB2	1.84	0.43
29:E:189:THR:OG1	29:E:191:ASP:OD1	2.33	0.43
36:N:20:GLY:HA2	36:N:28:GLY:HA2	2.01	0.43
37:O:42:THR:CG2	37:O:45:GLN:HG3	2.49	0.43
1:a:1060:U:H2'	1:a:1061:G:H8	1.84	0.43
14:n:65:ARG:HG2	14:n:78:GLY:O	2.19	0.43
24:v:20:SER:N	24:v:26:LYS:HE2	2.33	0.43
24:v:207:VAL:O	24:v:207:VAL:HG12	2.19	0.43
25:A:1540:G:H2'	25:A:1541:C:C6	2.54	0.43
25:A:1921:G:H2'	25:A:1922:G:O4'	2.19	0.43
25:A:2101:A:H2'	25:A:2102:G:C8	2.54	0.43
25:A:2114:A:N6	25:A:2119:A:C6	2.81	0.43
25:A:2662:A:H8	25:A:2662:A:O5'	2.01	0.43
27:C:144:VAL:HB	27:C:154:LEU:HB2	2.00	0.43
1:a:299:G:H2'	1:a:300:A:C8	2.54	0.43
1:a:1149:C:P	9:i:11:ARG:HH21	2.40	0.43
4:d:126:ASN:ND2	4:d:141:ASP:OD1	2.52	0.43
8:h:79:SER:OG	8:h:124:GLU:OE2	2.20	0.43
14:n:88:ALA:HB2	14:n:96:LEU:HD23	2.01	0.43
21:u:39:GLU:HB3	21:u:44:GLU:OE2	2.19	0.43
24:v:107:ASP:OD1	24:v:319:ARG:NH2	2.52	0.43
26:B:66:A:H5''	26:B:67:G:OP1	2.19	0.43
39:Q:67:ASN:OD1	39:Q:70:ALA:N	2.46	0.43
51:3:14:ALA:HA	51:3:32:LEU:O	2.18	0.43
1:a:215:C:H2'	1:a:216:U:C6	2.53	0.43
1:a:634:C:H2'	1:a:635:A:C8	2.53	0.43
1:a:767:A:H2'	1:a:768:A:O4'	2.19	0.43
6:f:45:ARG:HD2	6:f:59:TYR:CE2	2.54	0.43
6:f:96:VAL:HG12	6:f:96:VAL:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:j:80:THR:O	10:j:83:THR:HG22	2.19	0.43
11:k:74:VAL:HG11	11:k:79:ILE:HD12	2.01	0.43
24:v:20:SER:N	24:v:26:LYS:CE	2.82	0.43
24:v:162:LEU:HD23	24:v:162:LEU:H	1.83	0.43
24:v:292:SER:OG	24:v:293:GLU:N	2.52	0.43
24:v:478:ASP:CG	24:v:481:LYS:HB2	2.44	0.43
25:A:2139:U:H2'	25:A:2140:G:C8	2.54	0.43
25:A:2194:U:H2'	25:A:2195:U:H6	1.84	0.43
25:A:2377:A:H8	25:A:2377:A:OP2	2.01	0.43
25:A:2646:C:OP2	25:A:2732:G:O2'	2.36	0.43
49:l:46:VAL:O	49:l:50:VAL:HG23	2.19	0.43
2:b:72:THR:O	2:b:72:THR:HG22	2.19	0.43
5:e:148:ASN:HB3	5:e:153:VAL:HG13	2.01	0.43
6:f:73:GLU:O	6:f:76:THR:OG1	2.31	0.43
13:m:33:ILE:HG23	13:m:59:GLU:HB3	2.01	0.43
14:n:6:MET:HB3	14:n:63:ARG:NH2	2.34	0.43
19:s:60:VAL:HG21	19:s:74:PHE:O	2.19	0.43
25:A:414:C:H2'	25:A:415:A:C8	2.54	0.43
25:A:2162:G:P	25:A:2162:G:H8	2.42	0.43
28:D:55:LYS:HB2	28:D:76:GLY:O	2.19	0.43
33:J:92:LYS:HB2	33:J:92:LYS:HE3	1.79	0.43
4:d:118:VAL:HA	4:d:123:ILE:CD1	2.48	0.42
6:f:10:VAL:HA	6:f:84:VAL:HA	2.00	0.42
11:k:74:VAL:O	11:k:74:VAL:HG12	2.19	0.42
24:v:120:GLU:N	24:v:120:GLU:OE1	2.52	0.42
25:A:1060:U:N3	25:A:1088:A:H8	2.17	0.42
25:A:1113:U:H2'	25:A:1114:C:C6	2.54	0.42
35:M:47:ILE:HD13	35:M:47:ILE:HA	1.89	0.42
54:6:25:LYS:HE3	54:6:25:LYS:HB3	1.70	0.42
1:a:1238:A:H2	1:a:1241:G:N3	2.17	0.42
13:m:75:MET:O	13:m:79:ARG:HG2	2.19	0.42
25:A:984:A:H2'	25:A:984:A:N3	2.33	0.42
25:A:1902:C:H4'	27:C:242:LYS:O	2.19	0.42
29:E:161:ALA:HB1	29:E:167:VAL:HG23	2.01	0.42
32:I:54:VAL:HG22	32:I:56:ARG:NH1	2.34	0.42
36:N:124:GLY:H	36:N:144:GLU:HA	1.83	0.42
40:R:51:ARG:NH1	40:R:53:ARG:HG3	2.35	0.42
1:a:996:A:H2'	1:a:997:U:C6	2.55	0.42
1:a:1207:2MG:H2'	1:a:1208:C:H6	1.79	0.42
1:a:1313:U:H2'	1:a:1314:C:C6	2.54	0.42
5:e:45:ARG:HA	5:e:72:ILE:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:v:13:ARG:O	24:v:14:ARG:HD2	2.18	0.42
24:v:118:GLY:HA2	24:v:154:LEU:HD11	2.00	0.42
24:v:259:ALA:CA	58:v:601:GCP:O6	2.67	0.42
24:v:304:GLN:HG3	24:v:305:ALA:N	2.34	0.42
25:A:1:G:H2'	25:A:2:G:C8	2.52	0.42
25:A:572:A:OP2	42:T:80:ARG:NH2	2.45	0.42
25:A:1412:U:H2'	25:A:1413:A:H8	1.84	0.42
25:A:2127:G:N2	25:A:2173:A:H1'	2.34	0.42
25:A:2147:A:C2	25:A:2148:G:H1'	2.54	0.42
31:G:31:GLY:HA3	31:G:79:VAL:HG22	2.01	0.42
32:I:94:ARG:HG3	32:I:129:LEU:HD13	2.02	0.42
33:J:26:PRO:C	33:J:29:GLY:H	2.27	0.42
51:3:28:VAL:HG21	51:3:32:LEU:HD11	2.01	0.42
1:a:359:G:H5''	24:v:312:ARG:CZ	2.49	0.42
1:a:721:G:H4'	1:a:722:G:O4'	2.20	0.42
1:a:1319:A:C8	1:a:1323:G:C6	3.07	0.42
4:d:147:GLU:OE1	4:d:150:LYS:NZ	2.53	0.42
24:v:425:VAL:CG1	24:v:449:VAL:HG21	2.50	0.42
24:v:492:GLN:O	24:v:493:LEU:HD22	2.18	0.42
25:A:1097:U:H3'	25:A:1098:A:H8	1.84	0.42
25:A:1730:C:HO2'	25:A:1731:G:P	2.42	0.42
25:A:2138:G:C2	25:A:2154:A:C2	3.07	0.42
30:F:136:ILE:HD13	30:F:143:TYR:HD2	1.84	0.42
31:G:77:ILE:H	31:G:77:ILE:HD12	1.84	0.42
33:J:42:PHE:O	33:J:46:THR:OG1	2.20	0.42
36:N:82:LEU:HD12	36:N:90:VAL:HG21	2.02	0.42
36:N:109:LYS:HG2	36:N:126:ARG:HB2	2.00	0.42
1:a:524:G:H2'	1:a:525:C:C6	2.54	0.42
1:a:1296:C:N4	1:a:1297:G:O6	2.52	0.42
7:g:50:LEU:HA	7:g:53:ARG:HG2	1.99	0.42
24:v:464:GLU:HG2	24:v:465:SER:H	1.83	0.42
25:A:635:C:O2'	25:A:639:U:OP1	2.35	0.42
25:A:2131:U:H5'	25:A:2132:U:H5''	2.01	0.42
33:J:134:ARG:HG2	33:J:139:VAL:HA	2.01	0.42
46:X:24:ASN:O	46:X:25:LYS:HG2	2.20	0.42
1:a:477:C:H2'	1:a:478:A:C8	2.53	0.42
1:a:1240:U:C4	7:g:116:MET:HE1	2.54	0.42
6:f:9:MET:HE3	6:f:59:TYR:CZ	2.55	0.42
21:u:21:ARG:HD2	21:u:21:ARG:HA	1.94	0.42
21:u:36:GLU:C	21:u:37:PHE:HD2	2.27	0.42
25:A:288:U:H2'	25:A:289:G:C8	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:A:548:G:H2'	25:A:549:G:H1'	2.00	0.42
25:A:1111:A:HO2'	25:A:1112:G:P	2.41	0.42
26:B:54:G:H21	30:F:26:MET:CE	2.27	0.42
26:B:117:G:H2'	26:B:118:C:C6	2.54	0.42
31:G:72:LEU:HA	31:G:75:MET:HG2	2.01	0.42
44:V:11:LEU:O	49:1:29:ARG:NH2	2.51	0.42
1:a:6:G:O2'	1:a:7:A:H8	2.03	0.42
1:a:939:G:H2'	1:a:940:C:C6	2.55	0.42
1:a:1161:C:H2'	1:a:1162:C:C6	2.54	0.42
2:b:203:ASN:OD1	2:b:204:ASP:N	2.52	0.42
14:n:14:VAL:HA	14:n:60:GLN:NE2	2.35	0.42
17:q:57:ASP:OD1	17:q:82:ALA:N	2.51	0.42
24:v:33:VAL:HG23	24:v:269:LEU:HD22	2.01	0.42
25:A:323:C:H2'	29:E:163:ASN:OD1	2.20	0.42
25:A:1682:G:H2'	25:A:1683:U:C6	2.55	0.42
25:A:1734:G:H2'	25:A:1735:A:C8	2.55	0.42
25:A:1796:U:H2'	25:A:1797:G:H8	1.85	0.42
25:A:1831:G:H2'	25:A:1832:C:C6	2.55	0.42
25:A:2261:C:OP1	47:Y:19:LYS:NZ	2.43	0.42
35:M:71:ARG:NH2	35:M:123:LEU:O	2.53	0.42
36:N:3:LEU:HD23	36:N:3:LEU:HA	1.86	0.42
40:R:14:LYS:HE2	40:R:77:HIS:HA	2.02	0.42
41:S:107:THR:O	41:S:111:GLU:HG2	2.19	0.42
46:X:61:LEU:HD11	46:X:91:PHE:CE1	2.54	0.42
1:a:165:G:H2'	1:a:166:U:C6	2.55	0.42
1:a:1271:A:H2'	1:a:1272:G:C8	2.54	0.42
7:g:66:LEU:HA	7:g:69:VAL:HG12	2.02	0.42
25:A:1070:A:H3'	25:A:1071:G:H5''	2.01	0.42
25:A:1494:A:H2'	25:A:1495:A:C8	2.55	0.42
26:B:45:A:C4	26:B:46:A:C8	3.07	0.42
34:L:3:THR:HG21	41:S:61:TRP:HE1	1.85	0.42
45:W:43:LYS:HE2	45:W:43:LYS:HB2	1.89	0.42
1:a:109:A:H5'	1:a:110:C:C5	2.55	0.42
1:a:204:G:C4	1:a:205:A:C8	3.07	0.42
1:a:1030:U:H5''	1:a:1032:G:O6	2.20	0.42
4:d:45:LYS:HG3	4:d:47:ARG:HH21	1.85	0.42
24:v:144:LEU:CD1	24:v:148:ILE:HD13	2.50	0.42
24:v:318:MET:HE3	24:v:365:LEU:HD21	2.02	0.42
25:A:2812:G:H2'	25:A:2813:A:C8	2.55	0.42
29:E:141:MET:HG3	29:E:143:LEU:HD12	2.01	0.42
1:a:553:A:H2'	1:a:554:A:H8	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:f:26:THR:HA	6:f:29:ILE:HD13	2.02	0.42
12:l:67:ILE:HA	12:l:97:THR:HG22	2.02	0.42
20:t:62:ALA:HA	20:t:67:ILE:HG13	2.02	0.42
22:x:18:G:N2	22:x:57:G:H2'	2.34	0.42
25:A:1078:U:H1'	25:A:1088:A:H2	1.84	0.42
25:A:1585:C:H2'	25:A:1586:A:O4'	2.20	0.42
25:A:2494:G:H2'	25:A:2495:G:H8	1.85	0.42
25:A:2497:A:N3	25:A:2498:OMC:N4	2.68	0.42
1:a:671:G:H4'	6:f:79:ARG:HH22	1.85	0.41
1:a:1324:A:H2'	1:a:1325:C:H6	1.83	0.41
7:g:140:ASP:CG	7:g:143:ARG:HH21	2.28	0.41
24:v:79:TYR:CE2	24:v:80:HIS:CD2	3.08	0.41
24:v:146:ARG:HB2	25:A:2657:A:OP1	2.20	0.41
24:v:220:GLY:C	24:v:222:ASP:N	2.78	0.41
25:A:1056:G:H4'	25:A:1086:A:H8	1.85	0.41
28:D:25:THR:OG1	28:D:191:GLY:O	2.37	0.41
30:F:148:ARG:HD3	30:F:149:VAL:O	2.19	0.41
31:G:87:LEU:HD12	31:G:162:VAL:HG12	2.01	0.41
33:J:22:PRO:N	33:J:23:PRO:HD2	2.34	0.41
33:J:130:GLU:O	33:J:134:ARG:N	2.53	0.41
46:X:14:LYS:HE3	46:X:18:ARG:HH21	1.84	0.41
1:a:109:A:H5'	1:a:110:C:H5	1.84	0.41
1:a:1032:G:H3'	1:a:1032:G:N3	2.35	0.41
4:d:35:GLU:OE1	4:d:35:GLU:HA	2.20	0.41
4:d:105:MET:HE3	4:d:180:GLY:HA3	2.02	0.41
16:p:67:ILE:HG22	16:p:68:SER:O	2.21	0.41
18:r:50:LYS:HE3	18:r:50:LYS:HB2	1.82	0.41
25:A:78:U:H2'	25:A:79:C:C6	2.55	0.41
25:A:240:C:OP2	25:A:241:A:O2'	2.30	0.41
25:A:1199:U:H1'	41:S:4:VAL:HG22	2.01	0.41
25:A:1534:U:O2'	25:A:1537:G:O6	2.37	0.41
25:A:1915:3TD:O5'	25:A:1915:3TD:H6	2.19	0.41
36:N:123:ARG:NH2	36:N:143:GLU:OE2	2.53	0.41
38:P:24:MET:HE2	38:P:44:LEU:HD22	2.03	0.41
1:a:493:A:H2'	1:a:494:G:C8	2.55	0.41
1:a:1101:A:H4'	1:a:1102:A:O5'	2.21	0.41
24:v:94:ASP:CB	24:v:444:LEU:HB2	2.51	0.41
25:A:30:G:O2'	25:A:1214:A:N3	2.47	0.41
25:A:84:A:H4'	25:A:85:G:O5'	2.20	0.41
25:A:2153:C:H2'	25:A:2154:A:C8	2.56	0.41
25:A:2305:U:C2	30:F:151:GLY:HA3	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:L:11:VAL:HG21	34:L:50:THR:HG22	2.02	0.41
36:N:51:GLU:OE1	36:N:56:PRO:HA	2.20	0.41
1:a:71:A:C2	1:a:72:A:C8	3.08	0.41
1:a:96:U:H2'	1:a:97:G:C8	2.54	0.41
1:a:1009:U:C2	1:a:1021:A:C6	3.09	0.41
1:a:1157:A:C2	1:a:1181:G:C4	3.09	0.41
1:a:1326:U:C2	1:a:1327:C:C5	3.08	0.41
3:c:35:SER:OG	3:c:59:ARG:NH2	2.53	0.41
18:r:25:ASP:OD2	18:r:25:ASP:N	2.51	0.41
24:v:174:GLY:O	24:v:179:PHE:HA	2.20	0.41
24:v:327:LYS:HG2	24:v:328:GLY:N	2.36	0.41
25:A:2144:G:O2'	25:A:2147:A:N6	2.53	0.41
25:A:2515:C:H2'	25:A:2516:A:H8	1.86	0.41
51:3:41:HIS:CD2	51:3:43:PHE:H	2.39	0.41
1:a:505:G:H2'	1:a:506:G:H8	1.83	0.41
1:a:1319:A:C8	1:a:1323:G:C5	3.08	0.41
1:a:1371:G:O3'	9:i:71:GLY:HA3	2.20	0.41
2:b:58:ASN:OD1	2:b:59:LYS:N	2.54	0.41
3:c:46:GLU:HG3	3:c:47:LEU:CD2	2.51	0.41
4:d:44:ARG:NH1	4:d:46:PRO:HB3	2.35	0.41
7:g:68:ASN:O	7:g:135:VAL:HG22	2.20	0.41
7:g:116:MET:O	7:g:119:ARG:HB2	2.21	0.41
9:i:19:VAL:HG13	9:i:65:ILE:HG22	2.01	0.41
9:i:108:ALA:O	9:i:110:GLN:HG2	2.20	0.41
10:j:29:ALA:O	10:j:32:THR:OG1	2.35	0.41
11:k:33:THR:HG22	11:k:44:TRP:CB	2.51	0.41
12:l:110:ARG:HB3	12:l:119:VAL:HG11	2.03	0.41
13:m:81:MET:CE	13:m:91:HIS:HB3	2.50	0.41
24:v:10:VAL:HG22	24:v:361:PRO:HD2	2.02	0.41
25:A:1665:A:H1'	35:M:1:MET:CG	2.49	0.41
25:A:2194:U:H2'	25:A:2195:U:C6	2.56	0.41
25:A:2698:U:H2'	25:A:2699:C:C6	2.56	0.41
39:Q:64:TYR:HB3	39:Q:67:ASN:ND2	2.35	0.41
39:Q:76:LYS:O	39:Q:80:GLU:HG3	2.20	0.41
1:a:96:U:H2'	1:a:97:G:H8	1.85	0.41
1:a:678:U:H2'	1:a:679:C:C6	2.55	0.41
1:a:1452:C:H4'	1:a:1453:G:C2	2.55	0.41
3:c:123:GLN:HG2	3:c:126:ARG:HH21	1.86	0.41
23:z:17:U:H2'	23:z:18:A:C8	2.55	0.41
24:v:484:GLU:O	24:v:488:LYS:HG2	2.20	0.41
25:A:52:A:H2'	25:A:53:A:C8	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:A:657:U:H2'	25:A:658:U:C6	2.55	0.41
25:A:2107:G:C2	25:A:2108:A:H1'	2.55	0.41
25:A:2243:U:H2'	25:A:2244:U:C6	2.55	0.41
25:A:2581:G:OP2	25:A:2581:G:N2	2.47	0.41
25:A:2846:G:H2'	25:A:2847:U:C6	2.56	0.41
26:B:116:G:H4'	39:Q:54:VAL:HG12	2.03	0.41
1:a:83:C:H1'	1:a:87:C:H42	1.86	0.41
1:a:142:G:H2'	1:a:143:A:O4'	2.21	0.41
3:c:130:PHE:HE2	3:c:131:ARG:HE	1.68	0.41
14:n:34:VAL:HG13	14:n:35:ASN:ND2	2.36	0.41
24:v:27:THR:HA	24:v:30:THR:HG22	2.03	0.41
24:v:35:LEU:HD23	24:v:35:LEU:HA	1.80	0.41
24:v:101:ARG:C	24:v:103:LEU:H	2.28	0.41
25:A:1112:G:H2'	25:A:1113:U:H6	1.85	0.41
25:A:2308:G:N3	25:A:2308:G:H2'	2.36	0.41
26:B:28:C:H2'	26:B:29:A:H8	1.85	0.41
28:D:74:GLU:OE1	28:D:74:GLU:N	2.51	0.41
33:J:12:GLN:NE2	33:J:57:VAL:HG12	2.35	0.41
39:Q:39:VAL:HB	39:Q:49:VAL:HG22	2.02	0.41
46:X:61:LEU:HD12	46:X:61:LEU:O	2.21	0.41
55:7:15:LYS:HB2	55:7:23:LYS:HE2	2.03	0.41
56:8:11:CYS:HB3	56:8:33:HIS:CE1	2.56	0.41
1:a:270:A:H2'	1:a:271:C:C6	2.55	0.41
1:a:1107:C:H4'	3:c:169:ARG:HH22	1.86	0.41
1:a:1161:C:H2'	1:a:1162:C:H6	1.86	0.41
1:a:1180:A:P	9:i:99:ARG:HH22	2.43	0.41
1:a:1256:A:H5'	1:a:1258:G:H1'	2.01	0.41
1:a:1465:A:H2'	1:a:1466:C:C6	2.56	0.41
4:d:28:ILE:HD12	4:d:34:ILE:HG12	2.02	0.41
6:f:90:MET:HE1	18:r:61:ARG:CZ	2.51	0.41
11:k:35:THR:HA	11:k:41:ALA:HA	2.03	0.41
19:s:34:TRP:HA	19:s:52:HIS:HB3	2.02	0.41
24:v:37:GLY:HA3	24:v:77:PHE:CD2	2.55	0.41
24:v:60:MET:HE3	24:v:70:ILE:HG21	2.02	0.41
24:v:242:ASP:OD1	24:v:242:ASP:N	2.51	0.41
25:A:306:U:H2'	25:A:307:G:O4'	2.20	0.41
32:I:25:ALA:HB2	32:I:96:PHE:CD1	2.55	0.41
37:O:45:GLN:OE1	37:O:125:PRO:HD3	2.20	0.41
1:a:929:G:HO2'	1:a:930:C:P	2.44	0.41
1:a:1074:G:O2'	1:a:1101:A:N1	2.49	0.41
1:a:1305:G:H22	1:a:1331:G:HO2'	1.62	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:a:1330:U:H2'	1:a:1331:G:C8	2.56	0.41
4:d:146:ARG:HG2	4:d:147:GLU:N	2.35	0.41
6:f:45:ARG:HB3	6:f:59:TYR:CE2	2.56	0.41
6:f:73:GLU:O	6:f:77:THR:HG23	2.20	0.41
11:k:20:VAL:HG11	11:k:85:MET:HE1	2.02	0.41
11:k:122:ARG:HA	11:k:123:PRO:HD3	1.92	0.41
16:p:56:ARG:HD2	16:p:56:ARG:HA	1.76	0.41
25:A:495:G:N2	43:U:61:ASN:HD21	2.15	0.41
25:A:713:G:C2	25:A:714:U:H1'	2.56	0.41
25:A:1036:G:H1	25:A:1119:U:H3	1.69	0.41
25:A:1441:G:H2'	25:A:1442:U:C6	2.55	0.41
27:C:43:ARG:HA	27:C:48:ARG:O	2.20	0.41
27:C:117:GLN:HG3	27:C:118:SER:N	2.36	0.41
27:C:119:GLY:O	27:C:130:LEU:HD23	2.20	0.41
30:F:29:PRO:HB3	30:F:160:ALA:HB2	2.02	0.41
30:F:46:ASP:O	30:F:49:LEU:HG	2.21	0.41
31:G:86:LYS:HE3	31:G:86:LYS:HB2	1.95	0.41
33:J:33:VAL:HG13	33:J:65:ARG:HE	1.86	0.41
34:L:13:ARG:HG2	34:L:51:GLY:O	2.20	0.41
47:Y:72:LYS:HB3	47:Y:72:LYS:HE3	1.94	0.41
50:2:3:LYS:HA	50:2:3:LYS:HD2	1.72	0.41
55:7:8:ARG:HD2	55:7:8:ARG:HA	1.77	0.41
1:a:425:G:H21	4:d:40:GLN:NE2	2.19	0.41
1:a:768:A:H4'	1:a:1523:G:N2	2.36	0.41
1:a:1223:C:OP2	19:s:78:ARG:NH2	2.54	0.41
7:g:95:ARG:HG3	7:g:99:LEU:HD23	2.03	0.41
9:i:15:SER:OG	9:i:70:GLY:HA3	2.21	0.41
14:n:65:ARG:HA	14:n:65:ARG:HD2	1.86	0.41
18:r:34:THR:HG22	18:r:35:GLU:H	1.86	0.41
19:s:53:ASN:HB2	19:s:77:THR:HG22	2.03	0.41
25:A:480:A:O2'	45:W:43:LYS:O	2.39	0.41
25:A:587:C:H5'	29:E:85:PHE:CE2	2.56	0.41
25:A:2262:U:H2'	25:A:2263:C:H6	1.86	0.41
26:B:28:C:H2'	26:B:29:A:C8	2.56	0.41
28:D:181:ASP:OD2	28:D:186:LEU:HB2	2.21	0.41
30:F:11:GLU:HB2	30:F:15:LYS:HZ3	1.86	0.41
31:G:141:ILE:HD12	31:G:141:ILE:HA	1.99	0.41
33:J:84:ALA:HB2	33:J:101:ILE:HD13	2.03	0.41
1:a:1160:G:C2	1:a:1161:C:C6	3.10	0.40
1:a:1219:A:H2'	1:a:1220:G:H8	1.86	0.40
1:a:1267:C:H2'	1:a:1268:G:O4'	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:u:51:SER:O	21:u:55:ARG:HG2	2.21	0.40
25:A:956:G:H5''	37:O:76:LYS:HE2	2.03	0.40
25:A:1924:C:H3'	25:A:1925:C:H6	1.85	0.40
44:V:46:ALA:O	44:V:50:LEU:HG	2.21	0.40
1:a:678:U:H2'	1:a:679:C:H6	1.85	0.40
1:a:1027:C:C2	1:a:1028:C:H5	2.39	0.40
7:g:138:ARG:HG3	7:g:139:GLU:N	2.36	0.40
14:n:97:LYS:HE3	14:n:97:LYS:HB2	1.92	0.40
22:x:2:C:H2'	22:x:3:C:H6	1.85	0.40
22:x:51:U:H2'	22:x:52:G:C8	2.56	0.40
24:v:21:HIS:CD2	24:v:122:ARG:HB3	2.56	0.40
24:v:127:MET:SD	24:v:137:ILE:HD13	2.61	0.40
24:v:220:GLY:C	24:v:222:ASP:H	2.28	0.40
24:v:453:LEU:HD12	24:v:453:LEU:HA	1.87	0.40
25:A:1081:U:H2'	25:A:1082:U:C2	2.56	0.40
25:A:1182:G:H2'	25:A:1183:U:O4'	2.21	0.40
25:A:1910:G:H2'	25:A:1911:PSU:O4'	2.21	0.40
25:A:2316:G:H2'	25:A:2317:A:H8	1.86	0.40
25:A:2774:C:H2'	25:A:2775:G:O4'	2.21	0.40
31:G:128:GLN:HG2	31:G:129:THR:HG23	2.02	0.40
33:J:116:ASP:OD1	33:J:116:ASP:N	2.43	0.40
1:a:219:U:H2'	1:a:220:G:C8	2.54	0.40
1:a:604:G:H2'	1:a:605:U:O4'	2.22	0.40
1:a:1277:C:H1'	1:a:1282:C:O2	2.22	0.40
1:a:1310:G:O6	1:a:1328:C:N4	2.54	0.40
5:e:37:THR:HG21	5:e:64:MET:HE2	2.02	0.40
10:j:8:ILE:O	10:j:73:LEU:HD12	2.21	0.40
17:q:64:CYS:SG	17:q:67:LEU:HD11	2.62	0.40
25:A:81:G:H2'	25:A:82:U:O4'	2.22	0.40
25:A:288:U:H2'	25:A:289:G:H8	1.86	0.40
25:A:1338:G:O2'	25:A:1393:A:N1	2.50	0.40
25:A:1538:G:H2'	25:A:1539:U:C6	2.56	0.40
25:A:1796:U:H2'	25:A:1797:G:C8	2.57	0.40
29:E:7:ASP:CG	29:E:122:GLU:H	2.30	0.40
29:E:191:ASP:OD1	29:E:192:ALA:N	2.54	0.40
33:J:99:GLY:O	33:J:100:LYS:HE2	2.22	0.40
46:X:62:THR:HB	46:X:71:LYS:HD3	2.04	0.40
1:a:1004:A:OP1	1:a:1025:U:O2'	2.32	0.40
1:a:1006:G:H5'	14:n:19:LYS:HZ1	1.85	0.40
1:a:1031:C:H4'	1:a:1032:G:C6	2.55	0.40
1:a:1476:A:C2	1:a:1477:U:C4	3.08	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:d:54:GLN:HG3	4:d:199:LEU:HD12	2.04	0.40
5:e:144:LEU:HD23	5:e:144:LEU:HA	1.88	0.40
7:g:35:LYS:HA	7:g:35:LYS:HD3	1.83	0.40
19:s:13:LEU:HD12	19:s:13:LEU:HA	1.94	0.40
25:A:1056:G:H4'	25:A:1086:A:C8	2.56	0.40
25:A:2061:G:H2'	25:A:2501:C:O2'	2.21	0.40
25:A:2102:G:N1	25:A:2188:U:C4	2.89	0.40
25:A:2277:G:OP2	47:Y:10:THR:OG1	2.33	0.40
25:A:2291:U:H2'	25:A:2292:U:C6	2.57	0.40
25:A:2720:U:C2	25:A:2721:A:C8	3.10	0.40
40:R:60:THR:OG1	40:R:73:VAL:HG22	2.21	0.40
1:a:1004:A:N1	1:a:1026:G:H1'	2.36	0.40
1:a:1309:G:OP1	13:m:87:ARG:NH2	2.54	0.40
14:n:6:MET:HB3	14:n:63:ARG:CZ	2.51	0.40
24:v:317:PHE:CZ	24:v:348:PHE:HZ	2.40	0.40
24:v:407:LEU:HD12	24:v:407:LEU:HA	1.93	0.40
25:A:849:A:H2'	25:A:850:U:C6	2.57	0.40
25:A:1282:U:H2'	25:A:1283:G:O4'	2.22	0.40
25:A:2052:A:H4'	28:D:148:GLN:O	2.21	0.40
29:E:7:ASP:OD2	29:E:122:GLU:HG3	2.22	0.40
31:G:52:PHE:CZ	31:G:72:LEU:HD23	2.53	0.40
32:I:3:LEU:HD23	32:I:3:LEU:H	1.86	0.40
32:I:23:LEU:HA	32:I:116:GLU:OE2	2.22	0.40
32:I:71:CYS:HB3	32:I:74:ASP:HB2	2.02	0.40
33:J:34:ASN:HB3	33:J:37:GLU:HB2	2.03	0.40
33:J:51:LYS:HB2	33:J:51:LYS:HE2	1.83	0.40
41:S:15:LYS:HE2	41:S:15:LYS:HB3	1.79	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	
2	b	222/241 (92%)	203 (91%)	19 (9%)	0	100	100
3	c	204/233 (88%)	193 (95%)	11 (5%)	0	100	100
4	d	203/206 (98%)	194 (96%)	9 (4%)	0	100	100
5	e	154/167 (92%)	147 (96%)	7 (4%)	0	100	100
6	f	101/131 (77%)	95 (94%)	6 (6%)	0	100	100
7	g	150/156 (96%)	142 (95%)	8 (5%)	0	100	100
8	h	127/130 (98%)	122 (96%)	5 (4%)	0	100	100
9	i	125/130 (96%)	120 (96%)	5 (4%)	0	100	100
10	j	96/103 (93%)	89 (93%)	6 (6%)	1 (1%)	12	41
11	k	113/129 (88%)	108 (96%)	5 (4%)	0	100	100
12	l	118/124 (95%)	110 (93%)	8 (7%)	0	100	100
13	m	113/118 (96%)	102 (90%)	11 (10%)	0	100	100
14	n	98/101 (97%)	96 (98%)	2 (2%)	0	100	100
15	o	86/89 (97%)	82 (95%)	4 (5%)	0	100	100
16	p	79/82 (96%)	74 (94%)	5 (6%)	0	100	100
17	q	77/84 (92%)	74 (96%)	3 (4%)	0	100	100
18	r	52/75 (69%)	50 (96%)	2 (4%)	0	100	100
19	s	81/92 (88%)	77 (95%)	4 (5%)	0	100	100
20	t	84/87 (97%)	82 (98%)	2 (2%)	0	100	100
21	u	53/71 (75%)	51 (96%)	2 (4%)	0	100	100
24	v	523/529 (99%)	414 (79%)	104 (20%)	5 (1%)	12	41
27	C	269/273 (98%)	255 (95%)	14 (5%)	0	100	100
28	D	206/209 (99%)	196 (95%)	9 (4%)	1 (0%)	24	57
29	E	199/201 (99%)	196 (98%)	3 (2%)	0	100	100
30	F	175/179 (98%)	160 (91%)	15 (9%)	0	100	100
31	G	174/177 (98%)	167 (96%)	7 (4%)	0	100	100
32	I	129/165 (78%)	104 (81%)	25 (19%)	0	100	100
33	J	139/142 (98%)	123 (88%)	15 (11%)	1 (1%)	18	49
34	L	140/142 (99%)	136 (97%)	4 (3%)	0	100	100
35	M	121/123 (98%)	115 (95%)	6 (5%)	0	100	100
36	N	142/144 (99%)	138 (97%)	4 (3%)	0	100	100
37	O	132/136 (97%)	127 (96%)	5 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
38	P	116/127 (91%)	113 (97%)	3 (3%)	0	100	100
39	Q	114/117 (97%)	108 (95%)	6 (5%)	0	100	100
40	R	112/115 (97%)	109 (97%)	3 (3%)	0	100	100
41	S	115/118 (98%)	113 (98%)	2 (2%)	0	100	100
42	T	101/103 (98%)	99 (98%)	2 (2%)	0	100	100
43	U	108/110 (98%)	105 (97%)	3 (3%)	0	100	100
44	V	91/100 (91%)	86 (94%)	5 (6%)	0	100	100
45	W	100/104 (96%)	93 (93%)	7 (7%)	0	100	100
46	X	92/94 (98%)	88 (96%)	4 (4%)	0	100	100
47	Y	74/85 (87%)	71 (96%)	3 (4%)	0	100	100
48	Z	75/78 (96%)	71 (95%)	4 (5%)	0	100	100
49	1	59/63 (94%)	57 (97%)	2 (3%)	0	100	100
50	2	56/59 (95%)	54 (96%)	2 (4%)	0	100	100
51	3	46/70 (66%)	44 (96%)	2 (4%)	0	100	100
52	4	53/57 (93%)	51 (96%)	2 (4%)	0	100	100
53	5	49/55 (89%)	46 (94%)	3 (6%)	0	100	100
54	6	44/46 (96%)	42 (96%)	2 (4%)	0	100	100
55	7	62/65 (95%)	57 (92%)	5 (8%)	0	100	100
56	8	36/38 (95%)	36 (100%)	0	0	100	100
All	All	6188/6573 (94%)	5785 (94%)	395 (6%)	8 (0%)	49	78

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
24	v	18	ILE
24	v	293	GLU
24	v	459	VAL
10	j	57	VAL
28	D	149	ASN
33	J	33	VAL
24	v	494	ALA
24	v	306	ASN

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	
2	b	186/199 (94%)	186 (100%)	0	100	100
3	c	170/190 (90%)	170 (100%)	0	100	100
4	d	172/173 (99%)	172 (100%)	0	100	100
5	e	119/126 (94%)	119 (100%)	0	100	100
6	f	90/112 (80%)	90 (100%)	0	100	100
7	g	125/129 (97%)	125 (100%)	0	100	100
8	h	104/105 (99%)	104 (100%)	0	100	100
9	i	105/107 (98%)	105 (100%)	0	100	100
10	j	86/90 (96%)	86 (100%)	0	100	100
11	k	89/98 (91%)	89 (100%)	0	100	100
12	l	101/103 (98%)	101 (100%)	0	100	100
13	m	93/96 (97%)	93 (100%)	0	100	100
14	n	83/84 (99%)	83 (100%)	0	100	100
15	o	76/77 (99%)	76 (100%)	0	100	100
16	p	65/65 (100%)	65 (100%)	0	100	100
17	q	73/78 (94%)	73 (100%)	0	100	100
18	r	47/65 (72%)	47 (100%)	0	100	100
19	s	72/79 (91%)	72 (100%)	0	100	100
20	t	65/66 (98%)	65 (100%)	0	100	100
21	u	48/61 (79%)	48 (100%)	0	100	100
24	v	437/453 (96%)	432 (99%)	5 (1%)	65	78
27	C	216/218 (99%)	216 (100%)	0	100	100
28	D	163/163 (100%)	163 (100%)	0	100	100
29	E	165/165 (100%)	165 (100%)	0	100	100
30	F	148/150 (99%)	148 (100%)	0	100	100
31	G	137/138 (99%)	137 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
32	I	100/123 (81%)	100 (100%)	0	100	100
33	J	109/110 (99%)	109 (100%)	0	100	100
34	L	116/116 (100%)	116 (100%)	0	100	100
35	M	104/104 (100%)	104 (100%)	0	100	100
36	N	103/103 (100%)	103 (100%)	0	100	100
37	O	107/107 (100%)	107 (100%)	0	100	100
38	P	98/103 (95%)	98 (100%)	0	100	100
39	Q	86/87 (99%)	86 (100%)	0	100	100
40	R	99/100 (99%)	99 (100%)	0	100	100
41	S	89/90 (99%)	89 (100%)	0	100	100
42	T	84/84 (100%)	84 (100%)	0	100	100
43	U	93/93 (100%)	93 (100%)	0	100	100
44	V	80/84 (95%)	80 (100%)	0	100	100
45	W	83/85 (98%)	83 (100%)	0	100	100
46	X	78/78 (100%)	78 (100%)	0	100	100
47	Y	58/63 (92%)	58 (100%)	0	100	100
48	Z	67/68 (98%)	67 (100%)	0	100	100
49	1	54/55 (98%)	54 (100%)	0	100	100
50	2	48/49 (98%)	48 (100%)	0	100	100
51	3	44/62 (71%)	44 (100%)	0	100	100
52	4	46/48 (96%)	46 (100%)	0	100	100
53	5	46/49 (94%)	46 (100%)	0	100	100
54	6	38/38 (100%)	38 (100%)	0	100	100
55	7	51/52 (98%)	51 (100%)	0	100	100
56	8	34/34 (100%)	34 (100%)	0	100	100
All	All	5150/5375 (96%)	5145 (100%)	5 (0%)	87	90

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
24	v	26	LYS
24	v	122	ARG
24	v	189	GLU

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Mol	Chain	Res	Type
24	v	435	ASP
24	v	436	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (100) such sidechains are listed below:

Mol	Chain	Res	Type
2	b	18	HIS
2	b	39	HIS
2	b	42	ASN
2	b	94	HIS
2	b	109	GLN
2	b	170	HIS
3	c	123	GLN
3	c	140	ASN
3	c	176	HIS
4	d	40	GLN
4	d	41	HIS
4	d	54	GLN
4	d	116	GLN
4	d	136	GLN
4	d	198	HIS
5	e	70	ASN
5	e	77	ASN
5	e	82	GLN
5	e	83	HIS
5	e	148	ASN
6	f	11	HIS
6	f	14	GLN
6	f	58	HIS
7	g	9	GLN
7	g	130	ASN
8	h	67	GLN
8	h	118	GLN
9	i	75	GLN
10	j	15	HIS
10	j	64	GLN
11	k	22	HIS
11	k	28	ASN
11	k	101	ASN
12	l	5	ASN
12	l	29	GLN
12	l	112	GLN

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Mol	Chain	Res	Type
14	n	66	GLN
15	o	38	HIS
15	o	40	GLN
15	o	51	HIS
16	p	26	ASN
16	p	29	ASN
20	t	48	GLN
20	t	52	ASN
20	t	55	GLN
20	t	68	HIS
20	t	70	ASN
20	t	78	ASN
24	v	21	HIS
24	v	85	ASN
24	v	211	ASN
24	v	225	GLN
27	C	25	HIS
27	C	70	ASN
27	C	90	ASN
27	C	128	ASN
27	C	153	GLN
27	C	239	ASN
27	C	243	HIS
28	D	49	GLN
28	D	58	ASN
28	D	134	HIS
28	D	136	ASN
28	D	173	GLN
28	D	185	ASN
29	E	97	ASN
29	E	165	HIS
30	F	5	HIS
30	F	37	ASN
31	G	22	GLN
31	G	111	HIS
32	I	88	HIS
33	J	31	GLN
34	L	128	ASN
35	M	3	GLN
36	N	4	ASN
38	P	9	GLN
38	P	23	ASN

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Mol	Chain	Res	Type
40	R	12	GLN
40	R	15	GLN
41	S	72	ASN
42	T	82	HIS
42	T	91	GLN
43	U	9	HIS
43	U	57	ASN
43	U	60	HIS
43	U	61	ASN
45	W	54	GLN
45	W	66	GLN
46	X	49	ASN
46	X	75	GLN
47	Y	46	HIS
47	Y	50	ASN
47	Y	57	HIS
48	Z	16	ASN
49	1	45	GLN
51	3	6	HIS
51	3	41	HIS
54	6	6	GLN
55	7	28	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	a	1526/1542 (98%)	212 (13%)	0
22	x	72/76 (94%)	18 (25%)	0
23	z	7/21 (33%)	2 (28%)	0
25	A	2897/2904 (99%)	388 (13%)	14 (0%)
26	B	119/120 (99%)	18 (15%)	2 (1%)
All	All	4621/4663 (99%)	638 (13%)	16 (0%)

All (638) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	a	4	U
1	a	5	U
1	a	6	G
1	a	9	G
1	a	32	A

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Mol	Chain	Res	Type
1	a	39	G
1	a	44	A
1	a	47	C
1	a	48	C
1	a	50	A
1	a	51	A
1	a	59	A
1	a	69	G
1	a	71	A
1	a	78	A
1	a	80	A
1	a	81	A
1	a	82	G
1	a	85	U
1	a	86	G
1	a	87	C
1	a	92	U
1	a	95	C
1	a	116	A
1	a	121	U
1	a	122	G
1	a	130	A
1	a	131	A
1	a	144	G
1	a	163	C
1	a	164	G
1	a	174	A
1	a	177	G
1	a	184	G
1	a	197	A
1	a	199	A
1	a	209	U
1	a	210	C
1	a	211	G
1	a	212	G
1	a	240	G
1	a	245	U
1	a	247	G
1	a	251	G
1	a	266	G
1	a	267	C
1	a	280	C

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Mol	Chain	Res	Type
1	a	281	G
1	a	289	G
1	a	321	A
1	a	328	C
1	a	339	C
1	a	344	A
1	a	345	C
1	a	347	G
1	a	352	C
1	a	354	G
1	a	367	U
1	a	369	G
1	a	372	C
1	a	384	G
1	a	398	U
1	a	406	G
1	a	411	A
1	a	412	A
1	a	413	G
1	a	414	A
1	a	415	A
1	a	421	U
1	a	422	C
1	a	423	G
1	a	429	U
1	a	462	G
1	a	467	U
1	a	468	A
1	a	478	A
1	a	479	U
1	a	484	G
1	a	486	U
1	a	495	A
1	a	497	G
1	a	511	C
1	a	512	U
1	a	514	C
1	a	518	C
1	a	527	G7M
1	a	532	A
1	a	547	A
1	a	564	C

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Mol	Chain	Res	Type
1	a	572	A
1	a	573	A
1	a	576	C
1	a	577	G
1	a	593	U
1	a	633	G
1	a	650	G
1	a	653	U
1	a	665	A
1	a	683	G
1	a	684	U
1	a	705	G
1	a	723	U
1	a	724	G
1	a	742	G
1	a	747	A
1	a	755	G
1	a	760	G
1	a	777	A
1	a	793	U
1	a	794	A
1	a	802	A
1	a	815	A
1	a	817	C
1	a	819	A
1	a	821	G
1	a	851	G
1	a	890	G
1	a	914	A
1	a	926	G
1	a	929	G
1	a	930	C
1	a	934	C
1	a	935	A
1	a	958	A
1	a	960	U
1	a	966	2MG
1	a	969	A
1	a	975	A
1	a	976	G
1	a	977	A
1	a	982	U

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Mol	Chain	Res	Type
1	a	984	C
1	a	992	U
1	a	993	G
1	a	1003	G
1	a	1004	A
1	a	1006	G
1	a	1008	U
1	a	1009	U
1	a	1022	A
1	a	1024	G
1	a	1027	C
1	a	1029	U
1	a	1031	C
1	a	1032	G
1	a	1033	G
1	a	1034	G
1	a	1035	A
1	a	1043	G
1	a	1044	A
1	a	1046	A
1	a	1053	G
1	a	1065	U
1	a	1094	G
1	a	1095	U
1	a	1101	A
1	a	1136	C
1	a	1137	C
1	a	1139	G
1	a	1140	C
1	a	1145	A
1	a	1152	A
1	a	1159	U
1	a	1160	G
1	a	1167	A
1	a	1184	G
1	a	1196	A
1	a	1197	A
1	a	1212	U
1	a	1213	A
1	a	1225	A
1	a	1227	A
1	a	1238	A

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Mol	Chain	Res	Type
1	a	1257	A
1	a	1258	G
1	a	1260	G
1	a	1274	A
1	a	1275	A
1	a	1280	A
1	a	1287	A
1	a	1300	G
1	a	1302	C
1	a	1305	G
1	a	1308	U
1	a	1317	C
1	a	1320	C
1	a	1338	G
1	a	1346	A
1	a	1363	A
1	a	1378	C
1	a	1379	G
1	a	1388	C
1	a	1397	C
1	a	1398	A
1	a	1419	G
1	a	1441	A
1	a	1442	G
1	a	1446	A
1	a	1448	C
1	a	1452	C
1	a	1453	G
1	a	1473	G
1	a	1479	C
1	a	1492	A
1	a	1493	A
1	a	1499	A
1	a	1503	A
1	a	1506	U
1	a	1517	G
1	a	1520	C
1	a	1529	G
1	a	1530	G
22	x	9	A
22	x	14	A
22	x	19	G

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Mol	Chain	Res	Type
22	x	21	A
22	x	25	C
22	x	26	A
22	x	27	G
22	x	29	G
22	x	35	A
22	x	45	U
22	x	47	U
22	x	48	C
22	x	53	G
22	x	57	G
22	x	58	A
22	x	59	U
22	x	73	A
22	x	76	A
23	z	17	U
23	z	19	A
25	A	3	U
25	A	4	U
25	A	6	A
25	A	10	A
25	A	12	U
25	A	34	U
25	A	35	G
25	A	71	A
25	A	74	A
25	A	75	G
25	A	84	A
25	A	100	U
25	A	101	A
25	A	118	A
25	A	119	A
25	A	120	U
25	A	139	U
25	A	142	A
25	A	181	A
25	A	186	G
25	A	196	A
25	A	199	A
25	A	215	G
25	A	216	A
25	A	221	A

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Mol	Chain	Res	Type
25	A	222	A
25	A	233	A
25	A	248	G
25	A	269	C
25	A	271	G
25	A	272	A
25	A	277	G
25	A	278	A
25	A	279	A
25	A	281	C
25	A	311	A
25	A	329	G
25	A	330	A
25	A	357	C
25	A	361	G
25	A	362	A
25	A	367	G
25	A	386	G
25	A	393	C
25	A	396	G
25	A	404	A
25	A	405	U
25	A	406	G
25	A	411	G
25	A	412	A
25	A	428	A
25	A	481	G
25	A	491	G
25	A	504	A
25	A	505	A
25	A	508	A
25	A	509	C
25	A	532	A
25	A	545	U
25	A	546	U
25	A	547	A
25	A	549	G
25	A	563	A
25	A	573	U
25	A	575	A
25	A	603	A
25	A	614	A

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Mol	Chain	Res	Type
25	A	615	U
25	A	621	A
25	A	627	A
25	A	637	A
25	A	645	C
25	A	647	G
25	A	654	A
25	A	655	A
25	A	686	U
25	A	714	U
25	A	715	A
25	A	716	A
25	A	717	C
25	A	723	C
25	A	729	G
25	A	730	A
25	A	740	C
25	A	747	5MU
25	A	764	A
25	A	775	G
25	A	776	G
25	A	782	A
25	A	784	G
25	A	785	G
25	A	792	A
25	A	805	G
25	A	812	C
25	A	827	U
25	A	828	U
25	A	830	G
25	A	846	U
25	A	859	G
25	A	866	A
25	A	880	G
25	A	883	G
25	A	889	C
25	A	890	C
25	A	891	G
25	A	895	U
25	A	896	A
25	A	897	C
25	A	907	G

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Mol	Chain	Res	Type
25	A	910	A
25	A	914	G
25	A	915	C
25	A	931	U
25	A	945	A
25	A	946	C
25	A	961	C
25	A	974	G
25	A	983	A
25	A	990	A
25	A	996	A
25	A	1005	C
25	A	1012	U
25	A	1013	C
25	A	1026	G
25	A	1032	A
25	A	1033	U
25	A	1039	A
25	A	1045	C
25	A	1046	A
25	A	1047	G
25	A	1057	A
25	A	1059	G
25	A	1061	U
25	A	1062	G
25	A	1064	C
25	A	1065	U
25	A	1067	A
25	A	1068	G
25	A	1070	A
25	A	1071	G
25	A	1076	C
25	A	1081	U
25	A	1082	U
25	A	1084	A
25	A	1087	G
25	A	1088	A
25	A	1090	A
25	A	1095	A
25	A	1096	A
25	A	1101	U
25	A	1103	A

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Mol	Chain	Res	Type
25	A	1104	C
25	A	1106	G
25	A	1112	G
25	A	1119	U
25	A	1130	U
25	A	1132	U
25	A	1135	C
25	A	1136	G
25	A	1142	A
25	A	1206	G
25	A	1236	G
25	A	1253	A
25	A	1256	G
25	A	1268	A
25	A	1271	G
25	A	1272	A
25	A	1273	U
25	A	1275	A
25	A	1288	G
25	A	1289	C
25	A	1301	A
25	A	1313	U
25	A	1365	A
25	A	1379	U
25	A	1380	G
25	A	1383	A
25	A	1416	G
25	A	1428	C
25	A	1452	G
25	A	1458	U
25	A	1482	G
25	A	1493	C
25	A	1508	A
25	A	1509	A
25	A	1510	G
25	A	1515	A
25	A	1522	A
25	A	1524	G
25	A	1529	G
25	A	1535	A
25	A	1536	C
25	A	1537	G

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Mol	Chain	Res	Type
25	A	1539	U
25	A	1554	U
25	A	1566	A
25	A	1569	A
25	A	1584	U
25	A	1585	C
25	A	1606	C
25	A	1607	C
25	A	1647	U
25	A	1648	U
25	A	1674	G
25	A	1675	C
25	A	1676	A
25	A	1693	U
25	A	1698	A
25	A	1715	G
25	A	1723	G
25	A	1729	U
25	A	1730	C
25	A	1731	G
25	A	1732	C
25	A	1733	G
25	A	1738	G
25	A	1739	A
25	A	1744	A
25	A	1758	U
25	A	1764	C
25	A	1773	A
25	A	1800	C
25	A	1801	A
25	A	1808	A
25	A	1816	C
25	A	1829	A
25	A	1848	A
25	A	1858	A
25	A	1871	A
25	A	1872	A
25	A	1873	G
25	A	1879	C
25	A	1907	G
25	A	1913	A
25	A	1914	C

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Mol	Chain	Res	Type
25	A	1924	C
25	A	1925	C
25	A	1926	U
25	A	1927	A
25	A	1930	G
25	A	1936	A
25	A	1937	A
25	A	1938	A
25	A	1955	U
25	A	1967	C
25	A	1970	A
25	A	1971	U
25	A	1972	G
25	A	1991	U
25	A	1993	U
25	A	2006	C
25	A	2021	C
25	A	2023	C
25	A	2031	A
25	A	2033	A
25	A	2036	C
25	A	2043	C
25	A	2055	C
25	A	2056	G
25	A	2060	A
25	A	2061	G
25	A	2062	A
25	A	2069	G7M
25	A	2093	G
25	A	2098	U
25	A	2100	G
25	A	2108	A
25	A	2110	G
25	A	2112	G
25	A	2113	U
25	A	2115	G
25	A	2116	G
25	A	2118	U
25	A	2119	A
25	A	2124	G
25	A	2132	U
25	A	2133	G

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Mol	Chain	Res	Type
25	A	2136	G
25	A	2140	G
25	A	2146	C
25	A	2147	A
25	A	2157	G
25	A	2163	A
25	A	2164	C
25	A	2165	C
25	A	2167	U
25	A	2168	G
25	A	2172	U
25	A	2173	A
25	A	2174	C
25	A	2176	A
25	A	2178	C
25	A	2188	U
25	A	2190	G
25	A	2192	U
25	A	2198	A
25	A	2203	U
25	A	2204	G
25	A	2211	A
25	A	2212	A
25	A	2225	A
25	A	2238	G
25	A	2239	G
25	A	2279	G
25	A	2283	C
25	A	2287	A
25	A	2288	A
25	A	2294	G
25	A	2305	U
25	A	2308	G
25	A	2322	A
25	A	2324	U
25	A	2325	G
25	A	2333	A
25	A	2335	A
25	A	2350	C
25	A	2361	G
25	A	2377	A
25	A	2383	G

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Mol	Chain	Res	Type
25	A	2385	C
25	A	2401	U
25	A	2402	U
25	A	2406	A
25	A	2419	U
25	A	2424	C
25	A	2425	A
25	A	2429	G
25	A	2431	U
25	A	2441	U
25	A	2447	G
25	A	2448	A
25	A	2459	A
25	A	2469	A
25	A	2470	G
25	A	2474	U
25	A	2475	C
25	A	2476	A
25	A	2491	U
25	A	2492	U
25	A	2498	OMC
25	A	2502	G
25	A	2505	G
25	A	2517	C
25	A	2518	A
25	A	2520	C
25	A	2529	G
25	A	2547	A
25	A	2554	U
25	A	2566	A
25	A	2567	G
25	A	2573	C
25	A	2582	G
25	A	2602	A
25	A	2609	U
25	A	2613	U
25	A	2615	U
25	A	2624	G
25	A	2629	U
25	A	2636	C
25	A	2646	C
25	A	2649	C

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Mol	Chain	Res	Type
25	A	2663	G
25	A	2669	G
25	A	2689	U
25	A	2690	U
25	A	2724	U
25	A	2726	A
25	A	2733	A
25	A	2744	G
25	A	2748	A
25	A	2749	A
25	A	2752	C
25	A	2765	A
25	A	2778	A
25	A	2791	G
25	A	2792	A
25	A	2798	U
25	A	2800	A
25	A	2820	A
25	A	2821	A
25	A	2835	A
25	A	2849	U
25	A	2861	U
25	A	2873	A
25	A	2880	C
25	A	2884	U
25	A	2901	C
25	A	2902	C
26	B	4	C
26	B	9	G
26	B	24	G
26	B	35	C
26	B	36	C
26	B	44	G
26	B	45	A
26	B	52	A
26	B	56	G
26	B	65	U
26	B	67	G
26	B	88	C
26	B	89	U
26	B	90	C
26	B	91	C

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Mol	Chain	Res	Type
26	B	109	A
26	B	111	U
26	B	116	G

All (16) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
25	A	3	U
25	A	271	G
25	A	277	G
25	A	784	G
25	A	1026	G
25	A	1031	G
25	A	1111	A
25	A	1583	A
25	A	1730	C
25	A	2097	A
25	A	2189	U
25	A	2191	A
25	A	2473	U
25	A	2799	A
26	B	3	C
26	B	66	A

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

47 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
25	OMG	A	2251	25	23,26,27	1.21	3 (13%)	32,38,41	1.99	5 (15%)
25	H2U	A	2449	25	18,21,22	1.05	2 (11%)	19,30,33	1.06	3 (15%)
22	MIA	x	37	22	28,31,32	2.32	7 (25%)	38,44,47	2.77	14 (36%)
25	PSU	A	1911	25	18,21,22	1.41	3 (16%)	21,30,33	2.12	4 (19%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
22	PSU	x	39	22	18,21,22	1.34	2 (11%)	21,30,33	2.15	4 (19%)
1	MA6	a	1519	1	23,26,27	1.48	5 (21%)	33,38,41	2.43	15 (45%)
25	6MZ	A	2030	25	22,25,26	1.44	5 (22%)	29,36,39	2.28	9 (31%)
25	OMC	A	2498	25,57	19,22,23	0.83	0	25,31,34	1.17	2 (8%)
37	4D4	O	81	37	9,11,12	2.69	3 (33%)	7,13,15	0.97	1 (14%)
25	5MC	A	1962	25,57	19,22,23	1.31	3 (15%)	26,32,35	1.05	2 (7%)
1	5MC	a	1407	1	19,22,23	1.49	3 (15%)	26,32,35	1.12	3 (11%)
25	PSU	A	2604	25	18,21,22	1.47	4 (22%)	21,30,33	2.11	4 (19%)
1	2MG	a	966	1	23,26,27	1.28	4 (17%)	33,38,41	2.22	8 (24%)
1	2MG	a	1516	1	23,26,27	1.25	3 (13%)	33,38,41	2.24	7 (21%)
25	PSU	A	2457	25	18,21,22	1.40	4 (22%)	21,30,33	2.06	4 (19%)
25	PSU	A	2580	25	18,21,22	1.41	5 (27%)	21,30,33	2.09	4 (19%)
1	PSU	a	516	1,57	18,21,22	1.38	3 (16%)	21,30,33	2.10	4 (19%)
1	4OC	a	1402	1	20,23,24	0.76	0	25,32,35	0.92	1 (4%)
25	PSU	A	2605	25	18,21,22	1.46	4 (22%)	21,30,33	2.13	4 (19%)
25	2MA	A	2503	25	22,25,26	1.48	5 (22%)	32,37,40	2.26	9 (28%)
25	2MG	A	1835	25	23,26,27	1.26	3 (13%)	33,38,41	2.15	7 (21%)
1	2MG	a	1207	1	23,26,27	1.26	4 (17%)	33,38,41	2.18	7 (21%)
25	OMU	A	2552	25	19,22,23	1.34	4 (21%)	25,31,34	1.95	6 (24%)
22	PSU	x	55	22	18,21,22	1.38	2 (11%)	21,30,33	2.03	4 (19%)
28	MEQ	D	150	28	8,9,10	0.48	0	5,10,12	0.35	0
25	6MZ	A	1618	25	22,25,26	1.41	4 (18%)	29,36,39	2.28	9 (31%)
25	G7M	A	2069	25	23,26,27	1.57	4 (17%)	34,39,42	1.73	5 (14%)
25	2MG	A	2445	25	23,26,27	1.27	4 (17%)	33,38,41	2.18	10 (30%)
1	5MC	a	967	1	19,22,23	1.48	3 (15%)	26,32,35	1.14	3 (11%)
1	G7M	a	527	1	23,26,27	1.57	3 (13%)	34,39,42	1.75	5 (14%)
1	UR3	a	1498	1	19,22,23	0.94	1 (5%)	26,32,35	1.75	4 (15%)
22	5MU	x	54	22	19,22,23	1.44	6 (31%)	27,32,35	1.84	5 (18%)
12	D2T	l	89	12	8,9,10	2.82	1 (12%)	6,11,13	2.58	2 (33%)
25	PSU	A	746	25,57	18,21,22	1.45	4 (22%)	21,30,33	2.10	3 (14%)
37	MS6	O	82	37	5,7,8	0.57	0	2,7,9	1.29	0
25	PSU	A	2504	25,57	18,21,22	1.40	4 (22%)	21,30,33	1.95	3 (14%)
25	5MU	A	747	25	19,22,23	1.40	4 (21%)	27,32,35	2.18	7 (25%)
22	PSU	x	32	22	18,21,22	1.42	3 (16%)	21,30,33	2.08	4 (19%)
11	IAS	k	119	11	6,7,8	0.97	0	3,8,10	1.45	1 (33%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
25	PSU	A	955	25	18,21,22	1.43	4 (22%)	21,30,33	2.20	4 (19%)
22	4SU	x	8	22	18,21,22	1.85	4 (22%)	25,30,33	2.49	5 (20%)
25	3TD	A	1915	25	19,22,23	4.23	7 (36%)	23,32,35	1.78	3 (13%)
25	5MU	A	1939	25	19,22,23	1.37	4 (21%)	27,32,35	2.29	6 (22%)
25	1MG	A	745	25	23,26,27	1.15	2 (8%)	33,39,42	1.80	5 (15%)
1	MA6	a	1518	1	23,26,27	1.46	4 (17%)	33,38,41	2.46	13 (39%)
25	PSU	A	1917	25	18,21,22	1.44	4 (22%)	21,30,33	2.00	4 (19%)
22	G7M	x	46	22	23,26,27	2.35	5 (21%)	34,39,42	3.09	10 (29%)

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	OMG	A	2251	25	-	0/9/27/28	0/3/3/3
25	H2U	A	2449	25	-	0/7/38/39	0/2/2/2
22	MIA	x	37	22	-	4/15/33/34	0/3/3/3
25	PSU	A	1911	25	-	0/7/25/26	0/2/2/2
22	PSU	x	39	22	-	1/7/25/26	0/2/2/2
1	MA6	a	1519	1	-	4/11/29/30	0/3/3/3
25	6MZ	A	2030	25	-	1/9/27/28	0/3/3/3
25	OMC	A	2498	25,57	-	3/9/27/28	0/2/2/2
37	4D4	O	81	37	-	6/11/12/14	-
25	5MC	A	1962	25,57	-	0/7/25/26	0/2/2/2
1	5MC	a	1407	1	-	0/7/25/26	0/2/2/2
25	PSU	A	2604	25	-	0/7/25/26	0/2/2/2
1	2MG	a	966	1	-	2/9/27/28	0/3/3/3
1	2MG	a	1516	1	-	0/9/27/28	0/3/3/3
25	PSU	A	2457	25	-	0/7/25/26	0/2/2/2
25	PSU	A	2580	25	-	0/7/25/26	0/2/2/2
1	PSU	a	516	1,57	-	0/7/25/26	0/2/2/2
1	4OC	a	1402	1	-	3/9/29/30	0/2/2/2
25	PSU	A	2605	25	-	0/7/25/26	0/2/2/2
25	2MA	A	2503	25	-	1/7/25/26	0/3/3/3
25	2MG	A	1835	25	-	2/9/27/28	0/3/3/3
1	2MG	a	1207	1	-	0/9/27/28	0/3/3/3
25	OMU	A	2552	25	-	1/9/27/28	0/2/2/2
22	PSU	x	55	22	-	0/7/25/26	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
28	MEQ	D	150	28	-	2/8/9/11	-
25	6MZ	A	1618	25	-	0/9/27/28	0/3/3/3
25	G7M	A	2069	25	-	0/7/25/26	0/3/3/3
25	2MG	A	2445	25	-	1/9/27/28	0/3/3/3
1	5MC	a	967	1	-	0/7/25/26	0/2/2/2
1	G7M	a	527	1	-	1/7/25/26	0/3/3/3
1	UR3	a	1498	1	-	0/7/25/26	0/2/2/2
22	5MU	x	54	22	-	0/7/25/26	0/2/2/2
12	D2T	l	89	12	-	1/7/12/14	-
25	PSU	A	746	25,57	-	3/7/25/26	0/2/2/2
37	MS6	O	82	37	-	0/4/6/8	-
25	PSU	A	2504	25,57	-	0/7/25/26	0/2/2/2
25	5MU	A	747	25	-	0/7/25/26	0/2/2/2
22	PSU	x	32	22	-	0/7/25/26	0/2/2/2
11	IAS	k	119	11	-	1/7/7/8	-
25	PSU	A	955	25	-	0/7/25/26	0/2/2/2
22	4SU	x	8	22	-	0/7/25/26	0/2/2/2
25	3TD	A	1915	25	-	0/7/25/26	0/2/2/2
25	5MU	A	1939	25	-	0/7/25/26	0/2/2/2
25	1MG	A	745	25	-	0/7/25/26	0/3/3/3
1	MA6	a	1518	1	-	2/11/29/30	0/3/3/3
25	PSU	A	1917	25	-	2/7/25/26	0/2/2/2
22	G7M	x	46	22	-	3/7/25/26	0/3/3/3

All (156) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
25	A	1915	3TD	C6-C5	12.58	1.49	1.35
25	A	1915	3TD	C2-N1	9.75	1.49	1.37
12	l	89	D2T	CB-CA	-7.44	1.52	1.54
22	x	46	G7M	C8-N7	7.40	1.45	1.33
22	x	37	MIA	C2-S10	-7.24	1.69	1.75
22	x	37	MIA	C13-C14	6.79	1.52	1.32
37	O	81	4D4	CZ-NE	6.23	1.45	1.33
25	A	1915	3TD	C6-N1	5.87	1.45	1.36
1	a	527	G7M	C5-N7	-5.31	1.33	1.39
25	A	2069	G7M	C5-N7	-5.24	1.33	1.39
1	a	1407	5MC	C5-C4	5.18	1.48	1.44
22	x	8	4SU	C4-S4	-5.18	1.59	1.68
25	A	1915	3TD	C2-N3	5.12	1.49	1.38
1	a	967	5MC	C5-C4	5.12	1.48	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
22	x	46	G7M	C5-N7	-4.90	1.33	1.39
1	a	1518	MA6	C5-C4	4.47	1.47	1.39
1	a	1519	MA6	C5-C4	4.39	1.46	1.39
22	x	37	MIA	C5-C4	4.31	1.46	1.39
25	A	2503	2MA	C5-C4	4.30	1.46	1.39
25	A	1962	5MC	C5-C4	4.29	1.47	1.44
25	A	2030	6MZ	C5-C4	4.24	1.46	1.39
25	A	1618	6MZ	C5-C4	4.10	1.46	1.39
22	x	46	G7M	C8-N9	4.08	1.46	1.35
22	x	46	G7M	C5-C4	3.79	1.47	1.38
37	O	81	4D4	CZ-NH2	3.78	1.45	1.32
22	x	32	PSU	C6-C5	3.32	1.39	1.35
22	x	8	4SU	C4-N3	-3.28	1.34	1.37
25	A	1917	PSU	C6-C5	3.27	1.38	1.35
25	A	2552	OMU	C4-N3	-3.21	1.33	1.38
22	x	55	PSU	C6-C5	3.21	1.38	1.35
25	A	1911	PSU	C6-C5	3.18	1.38	1.35
25	A	747	5MU	C4-N3	-3.17	1.32	1.38
25	A	2604	PSU	C4-N3	-3.13	1.33	1.38
25	A	1939	5MU	C4-N3	-3.12	1.33	1.38
25	A	2605	PSU	C4-N3	-3.11	1.33	1.38
25	A	746	PSU	C4-N3	-3.09	1.33	1.38
25	A	955	PSU	C4-N3	-3.09	1.33	1.38
1	a	966	2MG	C5-C4	3.08	1.47	1.38
25	A	2251	OMG	C5-C4	3.05	1.47	1.38
1	a	1207	2MG	C5-C4	3.04	1.47	1.38
25	A	745	1MG	C5-C4	2.99	1.47	1.38
25	A	2552	OMU	C2-N3	-2.98	1.32	1.38
25	A	2504	PSU	C6-C5	2.92	1.38	1.35
1	a	1516	2MG	C5-C4	2.90	1.46	1.38
25	A	1835	2MG	C6-N1	-2.85	1.33	1.38
1	a	527	G7M	C5-C4	2.85	1.45	1.38
22	x	8	4SU	C5-C4	-2.85	1.39	1.42
25	A	2069	G7M	C5-C4	2.84	1.45	1.38
25	A	2457	PSU	C4-N3	-2.84	1.33	1.38
25	A	1835	2MG	C5-C4	2.84	1.46	1.38
22	x	54	5MU	C4-N3	-2.83	1.33	1.38
25	A	2445	2MG	C5-C4	2.83	1.46	1.38
22	x	54	5MU	C6-C5	2.83	1.39	1.34
25	A	2069	G7M	C6-N1	-2.82	1.33	1.38
25	A	2605	PSU	C6-C5	2.82	1.38	1.35
1	a	516	PSU	C6-C5	2.82	1.38	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
22	x	32	PSU	C4-N3	-2.79	1.33	1.38
25	A	1915	3TD	C4-N3	2.79	1.46	1.40
1	a	527	G7M	C6-N1	-2.79	1.33	1.38
25	A	2503	2MA	C5-C6	2.78	1.48	1.41
1	a	966	2MG	C6-N1	-2.78	1.33	1.38
25	A	1911	PSU	C4-N3	-2.78	1.33	1.38
1	a	1516	2MG	C6-N1	-2.77	1.33	1.38
25	A	2457	PSU	C6-C5	2.77	1.38	1.35
1	a	1519	MA6	C5-C6	2.77	1.48	1.41
25	A	2445	2MG	C6-N1	-2.76	1.33	1.38
25	A	2504	PSU	C4-N3	-2.75	1.33	1.38
22	x	39	PSU	C6-C5	2.75	1.38	1.35
25	A	2580	PSU	C4-N3	-2.73	1.33	1.38
1	a	516	PSU	C4-N3	-2.71	1.33	1.38
25	A	1917	PSU	C4-N3	-2.71	1.33	1.38
22	x	39	PSU	C4-N3	-2.71	1.33	1.38
25	A	2449	H2U	C4-N3	-2.69	1.33	1.37
22	x	55	PSU	C4-N3	-2.68	1.33	1.38
25	A	1962	5MC	C6-N1	-2.67	1.33	1.38
25	A	1939	5MU	C2-N3	-2.66	1.33	1.38
25	A	746	PSU	C6-C5	2.66	1.38	1.35
25	A	2251	OMG	C6-N1	-2.66	1.33	1.38
25	A	747	5MU	C6-N1	-2.65	1.33	1.38
25	A	2449	H2U	C2-N3	-2.65	1.33	1.38
25	A	1939	5MU	C6-N1	-2.65	1.33	1.38
25	A	2030	6MZ	C5-C6	2.62	1.48	1.41
25	A	1618	6MZ	C5-C6	2.62	1.48	1.41
25	A	1915	3TD	O2-C2	-2.62	1.18	1.23
25	A	2604	PSU	C6-C5	2.61	1.38	1.35
22	x	37	MIA	C5-C6	2.58	1.48	1.41
25	A	955	PSU	C6-C5	2.57	1.38	1.35
1	a	1518	MA6	C5-N7	-2.56	1.34	1.39
25	A	2030	6MZ	C5-N7	-2.55	1.34	1.39
1	a	967	5MC	C6-C5	2.53	1.38	1.34
1	a	1407	5MC	C6-C5	2.53	1.38	1.34
25	A	1618	6MZ	C5-N7	-2.51	1.34	1.39
1	a	1518	MA6	C5-C6	2.50	1.47	1.41
1	a	1207	2MG	C6-N1	-2.49	1.34	1.38
25	A	2605	PSU	C2-N3	-2.49	1.33	1.37
25	A	2580	PSU	C6-C5	2.48	1.38	1.35
25	A	2580	PSU	O4'-C1'	-2.48	1.40	1.43
22	x	46	G7M	C6-N1	-2.46	1.34	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
25	A	745	1MG	C5-N7	-2.45	1.34	1.39
25	A	747	5MU	C2-N3	-2.44	1.33	1.38
1	a	1407	5MC	C6-N1	-2.43	1.33	1.38
25	A	2604	PSU	C2-N3	-2.43	1.33	1.37
1	a	967	5MC	C6-N1	-2.42	1.33	1.38
25	A	955	PSU	C2-N3	-2.42	1.33	1.37
25	A	747	5MU	C6-C5	2.42	1.38	1.34
22	x	37	MIA	C5-N7	-2.41	1.34	1.39
25	A	2445	2MG	C5-N7	-2.38	1.34	1.39
1	a	1519	MA6	C5-N7	-2.37	1.34	1.39
1	a	1519	MA6	C8-N7	2.36	1.36	1.31
25	A	2503	2MA	C8-N7	2.35	1.36	1.31
25	A	2503	2MA	C5-N7	-2.34	1.34	1.39
37	O	81	4D4	CZ-NH1	-2.33	1.25	1.34
1	a	966	2MG	C5-N7	-2.32	1.34	1.39
25	A	2580	PSU	C2-N1	-2.32	1.33	1.36
25	A	2251	OMG	C5-N7	-2.31	1.34	1.39
22	x	54	5MU	C2-N3	-2.31	1.33	1.38
25	A	1835	2MG	C5-N7	-2.30	1.34	1.39
25	A	2604	PSU	C2-N1	-2.27	1.33	1.36
25	A	2552	OMU	C6-N1	-2.27	1.32	1.38
22	x	37	MIA	C8-N7	2.26	1.36	1.31
25	A	746	PSU	C2-N3	-2.24	1.33	1.37
22	x	54	5MU	C2-N1	2.24	1.42	1.38
25	A	2552	OMU	C5-C4	-2.22	1.38	1.43
25	A	746	PSU	C2-N1	-2.22	1.33	1.36
25	A	1939	5MU	C6-C5	2.22	1.38	1.34
25	A	1618	6MZ	C4-N9	-2.22	1.33	1.37
25	A	2605	PSU	C2-N1	-2.21	1.33	1.36
22	x	54	5MU	C4-C5	2.19	1.48	1.44
25	A	1962	5MC	C6-C5	2.18	1.38	1.34
25	A	2503	2MA	C4-N9	-2.17	1.33	1.37
25	A	2030	6MZ	C4-N9	-2.16	1.33	1.37
25	A	2457	PSU	C2-N3	-2.16	1.33	1.37
25	A	2030	6MZ	C8-N7	2.16	1.35	1.31
25	A	2504	PSU	C2-N3	-2.15	1.33	1.37
25	A	1917	PSU	C2-N3	-2.14	1.33	1.37
1	a	1207	2MG	C2-N3	2.14	1.36	1.32
1	a	966	2MG	C2-N3	2.13	1.36	1.32
1	a	1207	2MG	C5-N7	-2.12	1.34	1.39
1	a	1519	MA6	C4-N9	-2.12	1.33	1.37
22	x	8	4SU	C2-N3	-2.11	1.34	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
25	A	1917	PSU	C2-N1	-2.10	1.33	1.36
25	A	1915	3TD	O4-C4	-2.10	1.18	1.23
25	A	2580	PSU	C2-N3	-2.08	1.34	1.37
1	a	1516	2MG	C5-N7	-2.08	1.34	1.39
25	A	1911	PSU	C2-N3	-2.07	1.34	1.37
22	x	37	MIA	C4-N9	-2.07	1.33	1.37
25	A	955	PSU	C2-N1	-2.05	1.34	1.36
22	x	54	5MU	C6-N1	-2.05	1.34	1.38
1	a	1498	UR3	C5-C4	-2.04	1.38	1.43
1	a	1518	MA6	C8-N7	2.04	1.35	1.31
1	a	516	PSU	C2-N3	-2.04	1.34	1.37
22	x	32	PSU	C2-N3	-2.03	1.34	1.37
25	A	2445	2MG	C4-N9	-2.03	1.32	1.38
25	A	2457	PSU	C2-N1	-2.02	1.34	1.36
25	A	2069	G7M	C4-N9	-2.02	1.32	1.38
25	A	2504	PSU	C2-N1	-2.01	1.34	1.36

All (242) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
22	x	37	MIA	C12-C13-C14	-8.79	111.23	127.01
22	x	46	G7M	CN7-N7-C8	-8.50	111.91	124.79
25	A	2503	2MA	C5-C4-N3	-7.87	118.89	127.18
22	x	8	4SU	C4-N3-C2	-7.83	119.81	127.31
22	x	37	MIA	C5-C4-N3	-7.69	119.08	127.18
1	a	1516	2MG	C2-N3-C4	7.24	121.06	112.00
1	a	1207	2MG	C2-N3-C4	7.13	120.91	112.00
1	a	966	2MG	C2-N3-C4	7.04	120.81	112.00
25	A	1835	2MG	C2-N3-C4	6.99	120.75	112.00
25	A	2445	2MG	C2-N3-C4	6.93	120.67	112.00
1	a	1498	UR3	C4-N3-C2	-6.87	119.05	124.58
25	A	955	PSU	N1-C2-N3	6.83	122.37	115.17
22	x	46	G7M	N9-C8-N7	-6.71	96.18	112.48
25	A	746	PSU	N1-C2-N3	6.69	122.23	115.17
22	x	46	G7M	C8-N7-C5	6.63	116.07	107.78
25	A	1911	PSU	N1-C2-N3	6.63	122.16	115.17
25	A	2605	PSU	N1-C2-N3	6.61	122.14	115.17
22	x	39	PSU	N1-C2-N3	6.60	122.13	115.17
25	A	2604	PSU	N1-C2-N3	6.55	122.08	115.17
22	x	46	G7M	N9-C4-N3	6.54	139.03	125.95
22	x	32	PSU	N1-C2-N3	6.47	122.00	115.17
1	a	516	PSU	N1-C2-N3	6.47	121.99	115.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	a	966	2MG	C5-C4-N3	-6.43	118.15	128.39
22	x	8	4SU	C5-C4-N3	6.43	120.73	114.75
25	A	2457	PSU	N1-C2-N3	6.41	121.93	115.17
25	A	1917	PSU	N1-C2-N3	6.39	121.91	115.17
22	x	55	PSU	N1-C2-N3	6.38	121.90	115.17
25	A	2251	OMG	C5-C4-N3	-6.34	118.29	128.39
25	A	2580	PSU	N1-C2-N3	6.23	121.73	115.17
1	a	1207	2MG	C5-C4-N3	-6.21	118.51	128.39
25	A	2504	PSU	N1-C2-N3	6.14	121.64	115.17
1	a	1516	2MG	C5-C4-N3	-6.08	118.72	128.39
25	A	2445	2MG	C5-C4-N3	-6.03	118.79	128.39
25	A	1835	2MG	C5-C4-N3	-6.00	118.83	128.39
1	a	1518	MA6	C5-C4-N3	-5.92	118.56	126.72
22	x	37	MIA	N3-C4-N9	5.91	134.49	126.99
25	A	1618	6MZ	C5-C4-N3	-5.90	118.59	126.72
25	A	2030	6MZ	C5-C4-N3	-5.78	118.76	126.72
25	A	745	1MG	C5-C4-N3	-5.76	119.23	128.39
25	A	2503	2MA	N3-C4-N9	5.65	134.16	126.99
25	A	1939	5MU	C4-N3-C2	-5.65	119.93	127.34
1	a	527	G7M	N9-C4-N3	5.64	137.23	125.95
22	x	46	G7M	C5-C4-N3	-5.64	117.50	128.15
25	A	1915	3TD	N1-C2-N3	5.63	120.23	116.13
1	a	1519	MA6	C5-C4-N3	-5.58	119.03	126.72
25	A	1939	5MU	N3-C2-N1	5.55	122.12	114.89
25	A	2552	OMU	C4-N3-C2	-5.48	119.81	126.61
1	a	527	G7M	C5-C4-N3	-5.45	117.86	128.15
25	A	2069	G7M	C5-C4-N3	-5.27	118.19	128.15
25	A	2069	G7M	N9-C4-N3	5.27	136.49	125.95
25	A	2251	OMG	C2-N3-C4	5.17	121.21	112.30
25	A	747	5MU	C4-N3-C2	-5.15	120.58	127.34
25	A	747	5MU	N3-C2-N1	5.05	121.47	114.89
1	a	1518	MA6	N3-C4-N9	4.85	135.41	127.17
1	a	527	G7M	C2-N3-C4	4.82	120.59	112.30
25	A	2069	G7M	C2-N3-C4	4.81	120.58	112.30
1	a	966	2MG	N9-C4-N3	4.75	135.46	125.95
22	x	54	5MU	N3-C2-N1	4.71	121.02	114.89
25	A	1835	2MG	N9-C4-N3	4.70	135.35	125.95
25	A	1939	5MU	O4-C4-C5	-4.69	119.55	124.92
25	A	747	5MU	O4-C4-C5	-4.67	119.58	124.92
25	A	747	5MU	C5-C4-N3	4.64	119.35	115.32
25	A	1939	5MU	C5-C4-N3	4.62	119.34	115.32
1	a	1207	2MG	N9-C4-N3	4.60	135.16	125.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
25	A	2251	OMG	N9-C4-N3	4.54	135.03	125.95
25	A	955	PSU	C4-N3-C2	-4.53	120.13	126.37
25	A	2552	OMU	N3-C2-N1	4.53	120.78	114.89
25	A	2030	6MZ	C9-N6-C6	-4.52	118.66	122.85
25	A	1618	6MZ	N3-C4-N9	4.51	134.83	127.17
25	A	2445	2MG	N9-C4-N3	4.50	134.94	125.95
12	l	89	D2T	CB1-SB-CB	4.47	110.41	102.36
22	x	39	PSU	C4-N3-C2	-4.45	120.25	126.37
22	x	46	G7M	C2-N3-C4	4.43	119.93	112.30
25	A	745	1MG	N9-C4-N3	4.41	134.78	125.95
1	a	1519	MA6	C9-N6-C6	-4.38	109.37	120.52
25	A	745	1MG	C2-N3-C4	4.38	121.83	111.98
1	a	1518	MA6	C10-N6-C6	-4.38	109.38	120.52
1	a	1519	MA6	C10-N6-C6	-4.36	109.44	120.52
22	x	8	4SU	N3-C2-N1	4.34	120.55	114.89
25	A	746	PSU	C4-N3-C2	-4.34	120.40	126.37
25	A	2605	PSU	C4-N3-C2	-4.31	120.44	126.37
22	x	54	5MU	C4-N3-C2	-4.30	121.70	127.34
25	A	1618	6MZ	C9-N6-C6	-4.30	118.86	122.85
22	x	8	4SU	C5-C4-S4	-4.29	119.40	124.31
25	A	1915	3TD	C4-N3-C2	-4.29	120.08	124.61
25	A	2030	6MZ	N3-C4-N9	4.28	134.44	127.17
25	A	1911	PSU	C4-N3-C2	-4.26	120.50	126.37
25	A	2604	PSU	C4-N3-C2	-4.24	120.53	126.37
22	x	46	G7M	C8-N9-C4	4.20	117.49	107.09
1	a	1516	2MG	N9-C4-N3	4.20	134.35	125.95
1	a	1519	MA6	C2-N1-C6	4.20	122.08	111.83
1	a	516	PSU	C4-N3-C2	-4.19	120.60	126.37
22	x	46	G7M	CN7-N7-C5	4.18	132.01	126.80
22	x	32	PSU	C4-N3-C2	-4.16	120.64	126.37
25	A	2552	OMU	C5-C4-N3	4.16	120.62	114.80
25	A	2580	PSU	O2-C2-N1	-4.15	118.51	122.79
1	a	1518	MA6	N1-C6-N6	4.12	121.87	116.86
25	A	2457	PSU	C4-N3-C2	-4.11	120.71	126.37
25	A	1939	5MU	C5-C6-N1	-4.10	118.86	123.31
1	a	1519	MA6	N3-C4-N9	4.09	134.12	127.17
1	a	1518	MA6	C2-N1-C6	4.08	121.80	111.83
25	A	2580	PSU	C4-N3-C2	-4.04	120.81	126.37
1	a	1518	MA6	C9-N6-C6	-4.02	110.30	120.52
22	x	55	PSU	C4-N3-C2	-4.01	120.85	126.37
22	x	37	MIA	C16-C14-C13	-3.99	110.67	122.66
1	a	516	PSU	O2-C2-N1	-3.99	118.67	122.79

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
22	x	39	PSU	O2-C2-N1	-3.96	118.71	122.79
25	A	2030	6MZ	C6-C5-N7	3.95	136.74	132.43
25	A	746	PSU	O2-C2-N1	-3.93	118.73	122.79
25	A	2503	2MA	C4-C5-N7	-3.91	106.11	110.58
22	x	46	G7M	C1'-N9-C8	-3.88	113.65	126.74
25	A	1618	6MZ	C2-N3-C4	3.80	121.12	111.83
1	a	1519	MA6	C4-C5-N7	-3.80	106.23	110.58
25	A	955	PSU	O2-C2-N1	-3.78	118.89	122.79
25	A	1911	PSU	O2-C2-N1	-3.77	118.90	122.79
22	x	37	MIA	C15-C14-C13	-3.76	111.38	122.66
25	A	2030	6MZ	C4-C5-N7	-3.75	106.30	110.58
25	A	2030	6MZ	C2-N3-C4	3.72	120.92	111.83
25	A	2604	PSU	O2-C2-N1	-3.72	118.95	122.79
22	x	55	PSU	O2-C2-N1	-3.72	118.95	122.79
25	A	2457	PSU	O2-C2-N1	-3.70	118.97	122.79
25	A	1917	PSU	C4-N3-C2	-3.70	121.28	126.37
25	A	1618	6MZ	C6-C5-N7	3.70	136.46	132.43
1	a	1519	MA6	C2-N3-C4	3.69	120.85	111.83
22	x	37	MIA	C6-C5-N7	3.69	136.45	132.43
25	A	2504	PSU	C4-N3-C2	-3.69	121.29	126.37
22	x	32	PSU	O2-C2-N1	-3.67	119.00	122.79
1	a	1516	2MG	C6-C5-N7	3.67	136.97	130.29
22	x	54	5MU	C5-C4-N3	3.67	118.52	115.32
1	a	1518	MA6	C2-N3-C4	3.63	120.71	111.83
25	A	1618	6MZ	C4-C5-N7	-3.62	106.44	110.58
25	A	2504	PSU	O2-C2-N1	-3.55	119.13	122.79
22	x	37	MIA	C4-C5-N7	-3.50	106.58	110.58
25	A	1917	PSU	O2-C2-N1	-3.50	119.18	122.79
22	x	54	5MU	O4-C4-C5	-3.50	120.92	124.92
25	A	747	5MU	C5-C6-N1	-3.47	119.55	123.31
22	x	37	MIA	C2-N3-C4	3.42	121.24	112.29
12	l	89	D2T	OD2-CG-CB	3.39	120.48	113.15
22	x	37	MIA	C11-S10-C2	-3.39	99.71	102.25
22	x	54	5MU	C5-C6-N1	-3.37	119.65	123.31
25	A	2605	PSU	O2-C2-N1	-3.36	119.33	122.79
1	a	1519	MA6	N1-C2-N3	-3.34	123.53	128.58
1	a	1407	5MC	C5-C6-N1	-3.31	119.71	123.31
25	A	2498	OMC	O2-C2-N3	-3.29	117.14	122.33
1	a	967	5MC	C5-C6-N1	-3.28	119.75	123.31
1	a	1518	MA6	C4-C5-N7	-3.26	106.85	110.58
1	a	1498	UR3	C5-C4-N3	3.26	119.33	115.04
25	A	1618	6MZ	N1-C2-N3	-3.23	123.69	128.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	a	1207	2MG	C6-C5-N7	3.20	136.12	130.29
25	A	2251	OMG	C6-C5-N7	3.19	136.10	130.29
1	a	1518	MA6	N1-C2-N3	-3.13	123.85	128.58
25	A	1835	2MG	C6-C5-N7	3.09	135.91	130.29
25	A	2445	2MG	C6-C5-N7	3.07	135.87	130.29
25	A	2030	6MZ	N1-C2-N3	-3.05	123.96	128.58
25	A	1939	5MU	O2-C2-N1	-3.00	118.89	122.80
25	A	2552	OMU	O4-C4-C5	-2.99	120.01	125.16
1	a	1519	MA6	N1-C6-N6	2.99	120.49	116.86
1	a	1518	MA6	C5-C6-N6	-2.96	120.65	125.33
25	A	1962	5MC	C5-C6-N1	-2.94	120.12	123.31
1	a	1519	MA6	C10-N6-C9	-2.91	106.82	116.18
1	a	1516	2MG	C4-C5-N7	-2.88	106.10	110.67
1	a	966	2MG	C6-C5-N7	2.87	135.52	130.29
22	x	8	4SU	O2-C2-N1	-2.80	119.16	122.80
25	A	2503	2MA	C2-N1-C6	2.75	122.33	118.10
22	x	37	MIA	C4-N9-C8	2.74	108.61	105.74
25	A	1962	5MC	C5-C4-N3	-2.71	118.98	121.75
25	A	2552	OMU	O2-C2-N1	-2.69	119.29	122.80
1	a	966	2MG	CM2-N2-C2	-2.67	117.90	123.65
25	A	2503	2MA	C5-N7-C8	2.67	107.64	103.45
1	a	1518	MA6	C4-N9-C8	2.65	108.53	105.74
25	A	2503	2MA	N6-C6-N1	2.64	120.59	117.03
1	a	1519	MA6	C5-N7-C8	2.62	107.56	103.45
1	a	967	5MC	C5-C4-N3	-2.62	119.07	121.75
1	a	1407	5MC	C5-C4-N3	-2.60	119.09	121.75
25	A	2030	6MZ	C5-N7-C8	2.59	107.53	103.45
25	A	2503	2MA	C6-C5-N7	2.58	137.07	132.09
1	a	1207	2MG	C4-C5-N7	-2.58	106.59	110.67
22	x	37	MIA	C5-N7-C8	2.58	107.50	103.45
25	A	2251	OMG	C4-C5-N7	-2.55	106.62	110.67
25	A	1618	6MZ	C5-N7-C8	2.53	107.43	103.45
25	A	745	1MG	C4-C5-N7	-2.53	106.67	110.67
25	A	2503	2MA	C4-N9-C8	2.52	108.38	105.74
25	A	1618	6MZ	C4-N9-C8	2.50	108.36	105.74
25	A	745	1MG	C6-C5-N7	2.49	134.90	129.36
1	a	1519	MA6	C4-N9-C8	2.48	108.34	105.74
22	x	46	G7M	O6-C6-C5	-2.44	122.57	128.01
25	A	2445	2MG	C4-C5-N7	-2.43	106.81	110.67
22	x	37	MIA	N3-C2-N1	-2.42	122.59	127.00
1	a	966	2MG	C4-C5-N7	-2.40	106.86	110.67
1	a	1518	MA6	C10-N6-C9	-2.40	108.47	116.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	a	527	G7M	O6-C6-C5	-2.40	122.66	128.01
11	k	119	IAS	OD1-CG-CB	-2.40	118.40	125.38
1	a	1518	MA6	C5-N7-C8	2.39	107.21	103.45
25	A	2580	PSU	O4'-C1'-C2'	2.39	108.45	105.15
25	A	2069	G7M	O6-C6-C5	-2.37	122.72	128.01
25	A	2030	6MZ	C4-N9-C8	2.36	108.22	105.74
25	A	1915	3TD	C6-C5-C4	2.36	119.77	118.19
1	a	966	2MG	C2-N1-C6	-2.33	121.74	124.55
25	A	1835	2MG	C4-C5-N7	-2.32	107.00	110.67
25	A	2498	OMC	C1'-N1-C2	2.28	123.49	118.44
22	x	39	PSU	C5-C6-N1	-2.28	118.97	122.14
1	a	527	G7M	C5-C6-N1	2.28	116.55	111.84
25	A	2445	2MG	CM2-N2-C2	-2.27	118.78	123.65
25	A	2449	H2U	O2-C2-N1	-2.25	120.40	123.10
25	A	2069	G7M	C5-C6-N1	2.25	116.48	111.84
25	A	2445	2MG	N1-C2-N2	2.24	118.85	116.56
25	A	747	5MU	C1'-N1-C2	2.23	121.61	117.59
1	a	1402	4OC	C6-C5-C4	2.20	119.66	117.00
22	x	37	MIA	C2-N1-C6	2.20	121.71	117.54
25	A	2445	2MG	O6-C6-C5	-2.17	120.80	126.53
1	a	967	5MC	O2-C2-N3	-2.17	118.91	122.33
25	A	2457	PSU	O4'-C1'-C2'	2.16	108.15	105.15
25	A	2605	PSU	C5-C6-N1	-2.15	119.15	122.14
22	x	37	MIA	N9-C8-N7	-2.15	110.88	113.94
25	A	2503	2MA	N9-C8-N7	-2.13	110.91	113.94
1	a	516	PSU	O4'-C1'-C2'	2.12	108.09	105.15
25	A	2449	H2U	N3-C2-N1	2.12	118.78	116.65
1	a	1407	5MC	O2-C2-N3	-2.11	119.00	122.33
25	A	955	PSU	C5-C6-N1	-2.11	119.22	122.14
1	a	1207	2MG	C2-N1-C6	-2.09	122.02	124.55
25	A	1835	2MG	O6-C6-C5	-2.09	121.02	126.53
25	A	2552	OMU	C5-C6-N1	-2.08	118.46	121.84
1	a	1519	MA6	C5-C6-N6	-2.07	122.05	125.33
25	A	2449	H2U	C5-C6-N1	-2.06	105.28	111.52
37	O	81	4D4	O-C-CA	-2.06	119.47	124.77
1	a	1519	MA6	N9-C8-N7	-2.06	111.01	113.94
1	a	1498	UR3	C3U-N3-C4	2.06	120.72	117.87
1	a	966	2MG	O6-C6-C5	-2.06	121.10	126.53
1	a	1519	MA6	C6-C5-N7	2.05	136.71	133.43
25	A	1835	2MG	CM2-N2-C2	-2.05	119.24	123.65
22	x	55	PSU	C5-C6-N1	-2.05	119.30	122.14
1	a	1207	2MG	O6-C6-C5	-2.04	121.14	126.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
25	A	1911	PSU	C5-C6-N1	-2.04	119.31	122.14
1	a	1516	2MG	C2-N1-C6	-2.04	122.09	124.55
1	a	1516	2MG	N1-C2-N2	2.04	118.64	116.56
25	A	2445	2MG	C2-N1-C6	-2.03	122.10	124.55
1	a	1498	UR3	C1'-N1-C2	2.02	120.34	117.04
22	x	32	PSU	C5-C6-N1	-2.01	119.34	122.14
25	A	1917	PSU	C6-C5-C4	-2.01	116.82	118.17
25	A	747	5MU	O2-C2-N1	-2.01	120.18	122.80
25	A	2445	2MG	N2-C2-N3	-2.01	117.95	120.51
25	A	2604	PSU	C5-C6-N1	-2.01	119.36	122.14

There are no chirality outliers.

All (44) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	a	966	2MG	N3-C2-N2-CM2
1	a	1518	MA6	C5-C6-N6-C9
1	a	1519	MA6	C5-C6-N6-C9
22	x	37	MIA	N6-C12-C13-C14
22	x	37	MIA	C12-C13-C14-C15
22	x	37	MIA	C12-C13-C14-C16
37	O	81	4D4	C-CA-CB-OB
37	O	81	4D4	C-CA-CB-CG
37	O	81	4D4	N-CA-CB-OB
37	O	81	4D4	N-CA-CB-CG
37	O	81	4D4	OB-CB-CG-CD
25	A	746	PSU	C2'-C1'-C5-C4
25	A	1835	2MG	N1-C2-N2-CM2
25	A	1835	2MG	N3-C2-N2-CM2
1	a	1402	4OC	O4'-C4'-C5'-O5'
1	a	1519	MA6	O4'-C4'-C5'-O5'
22	x	46	G7M	O4'-C1'-N9-C4
1	a	1519	MA6	C3'-C4'-C5'-O5'
25	A	2552	OMU	O4'-C4'-C5'-O5'
28	D	150	MEQ	NE2-CD-CG-CB
1	a	1518	MA6	N1-C6-N6-C9
28	D	150	MEQ	OE1-CD-CG-CB
1	a	1402	4OC	C3'-C4'-C5'-O5'
22	x	37	MIA	O4'-C4'-C5'-O5'
1	a	1519	MA6	N1-C6-N6-C9
25	A	1917	PSU	C3'-C4'-C5'-O5'
12	l	89	D2T	CG-CB-SB-CB1

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Mol	Chain	Res	Type	Atoms
22	x	46	G7M	O4'-C1'-N9-C8
37	O	81	4D4	CA-CB-CG-CD
1	a	1402	4OC	C3'-C2'-O2'-CM2
1	a	527	G7M	C4'-C5'-O5'-P
25	A	2030	6MZ	O4'-C4'-C5'-O5'
22	x	46	G7M	C4'-C5'-O5'-P
25	A	746	PSU	O4'-C1'-C5-C6
25	A	2498	OMC	C2'-C1'-N1-C6
1	a	966	2MG	C3'-C4'-C5'-O5'
25	A	1917	PSU	O4'-C4'-C5'-O5'
25	A	2445	2MG	C3'-C4'-C5'-O5'
25	A	2498	OMC	O4'-C4'-C5'-O5'
25	A	746	PSU	C2'-C1'-C5-C6
22	x	39	PSU	O4'-C4'-C5'-O5'
25	A	2503	2MA	O4'-C4'-C5'-O5'
11	k	119	IAS	O-C-CA-CB
25	A	2498	OMC	C2'-C1'-N1-C2

There are no ring outliers.

16 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
25	A	1911	PSU	1	0
1	a	1519	MA6	1	0
25	A	2030	6MZ	3	0
25	A	2498	OMC	2	0
1	a	1407	5MC	1	0
1	a	1402	4OC	2	0
25	A	2503	2MA	1	0
1	a	1207	2MG	2	0
22	x	55	PSU	2	0
1	a	527	G7M	1	0
22	x	54	5MU	1	0
12	l	89	D2T	1	0
25	A	2504	PSU	1	0
25	A	1915	3TD	1	0
1	a	1518	MA6	1	0
22	x	46	G7M	3	0

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 740 ligands modelled in this entry, 739 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	$\# Z > 2$
58	GCP	v	601	-	32,34,34	0.77	1 (3%)	49,54,54	0.56	1 (2%)

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
58	GCP	v	601	-	-	1/19/38/38	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
58	v	601	GCP	PB-O2B	-2.29	1.50	1.56

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
58	v	601	GCP	O1G-PG-C3B	-2.11	106.77	111.37

There are no chirality outliers.

All (1) torsion outliers are listed below:

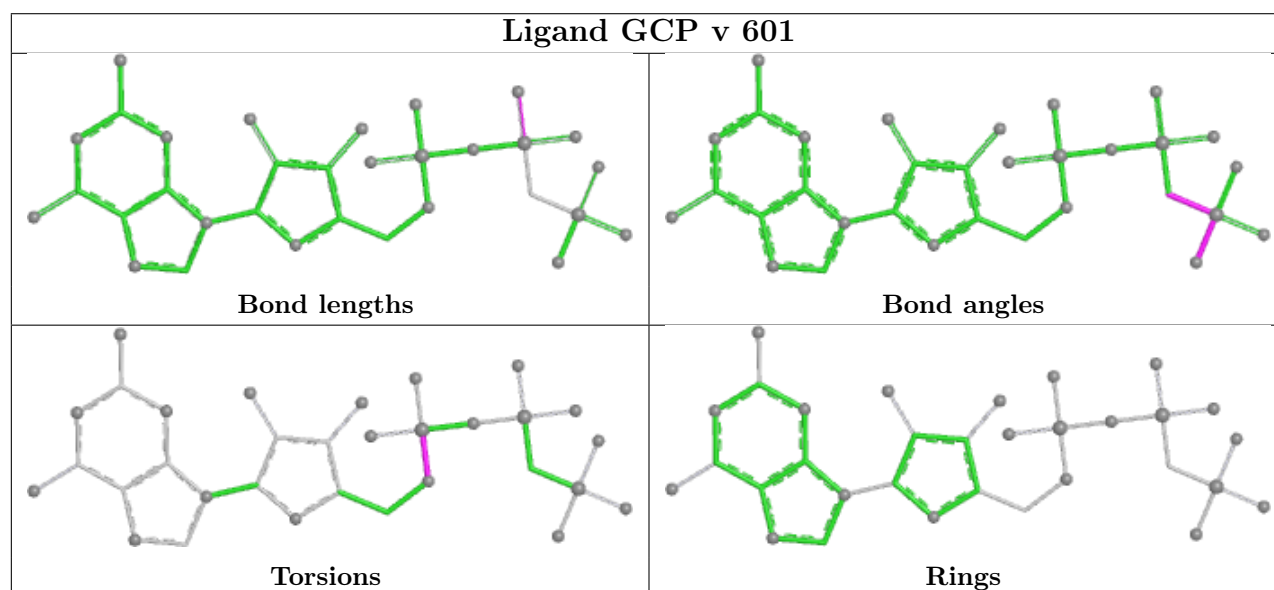
Mol	Chain	Res	Type	Atoms
58	v	601	GCP	C5'-O5'-PA-O1A

There are no ring outliers.

1 monomer is involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
58	v	601	GCP	9	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.

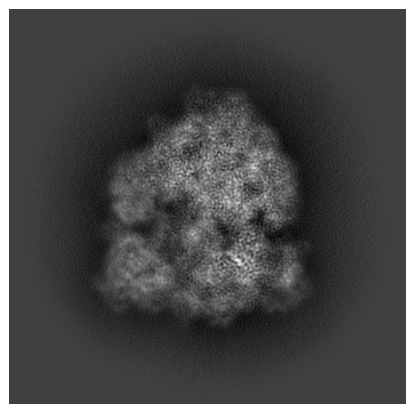
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-29631. These allow visual inspection of the internal detail of the map and identification of artifacts.

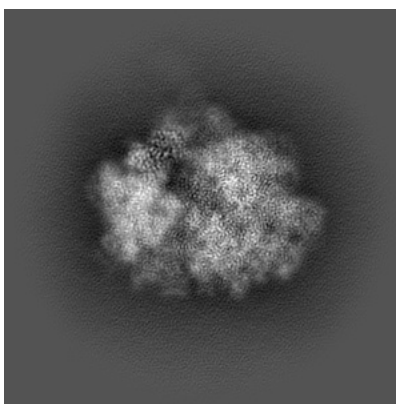
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

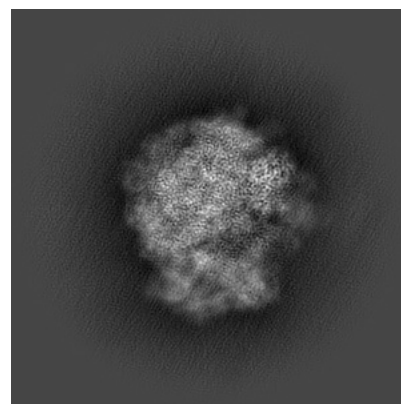
6.1.1 Primary map



X

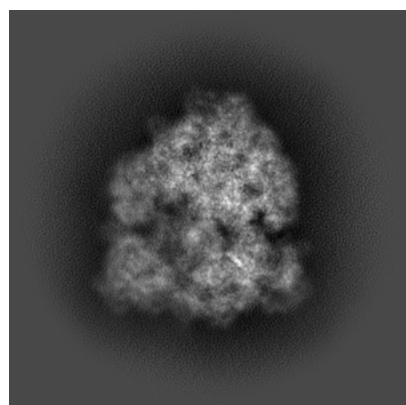


Y

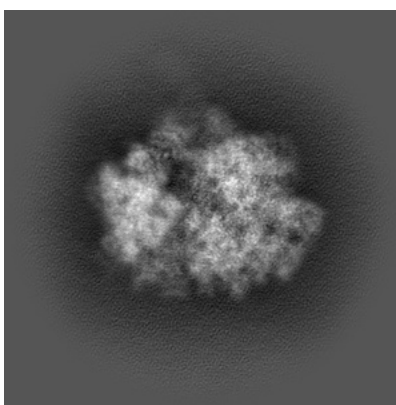


Z

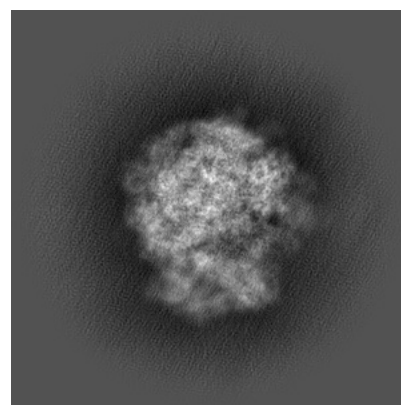
6.1.2 Raw map



X



Y

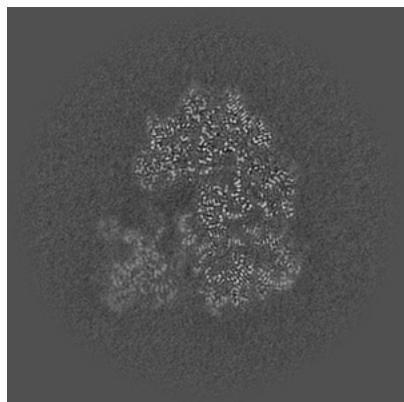


Z

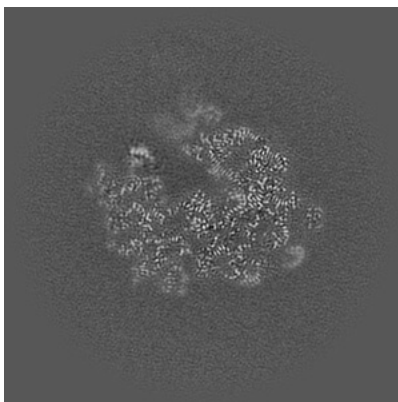
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

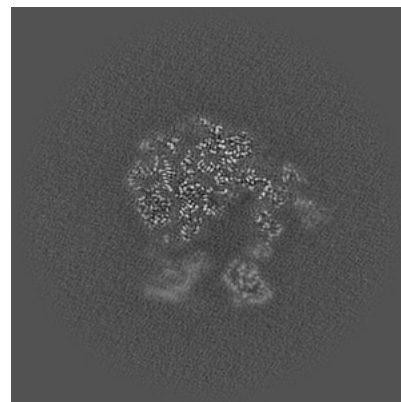
6.2.1 Primary map



X Index: 256

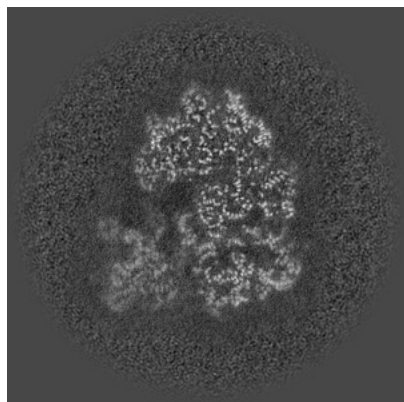


Y Index: 256

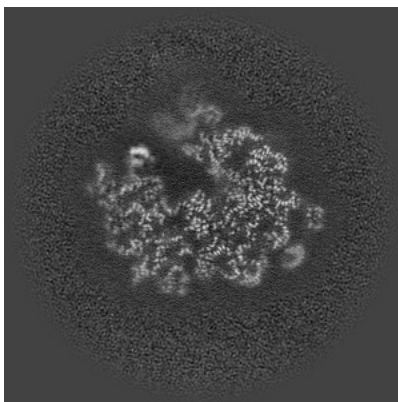


Z Index: 256

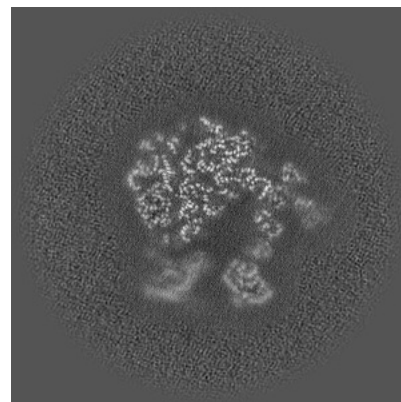
6.2.2 Raw map



X Index: 256



Y Index: 256

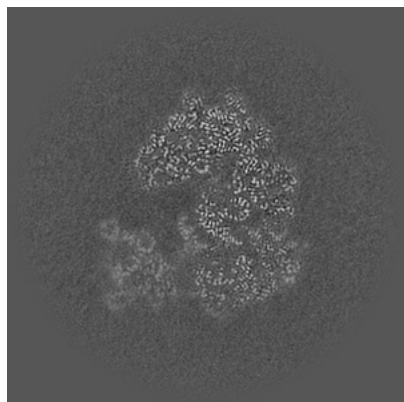


Z Index: 256

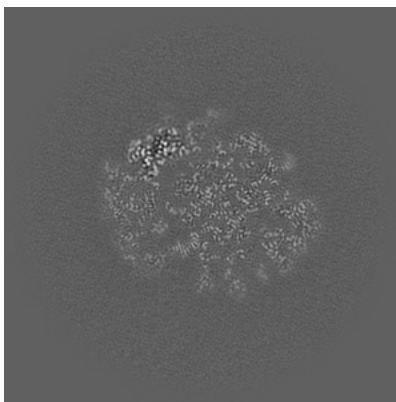
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

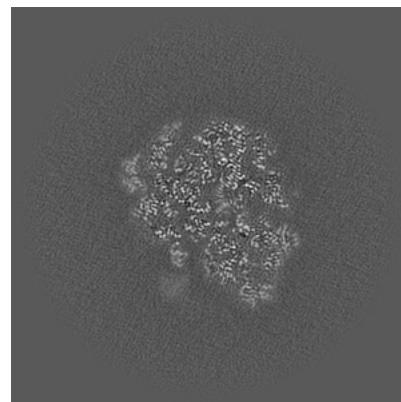
6.3.1 Primary map



X Index: 260

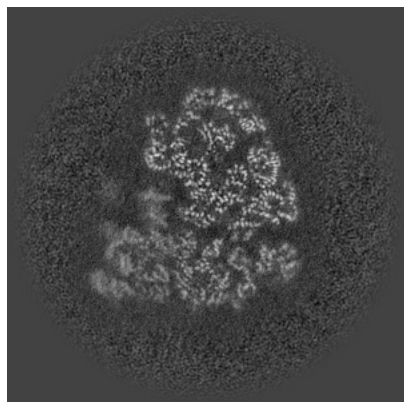


Y Index: 299

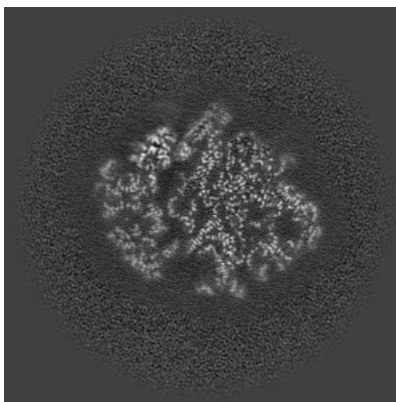


Z Index: 296

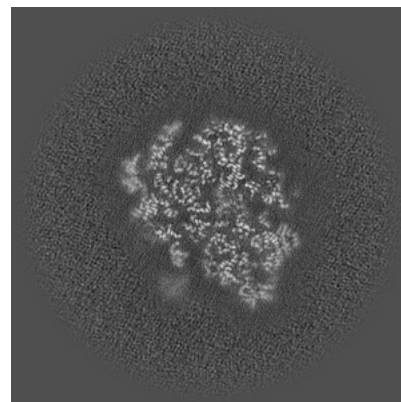
6.3.2 Raw map



X Index: 243



Y Index: 291

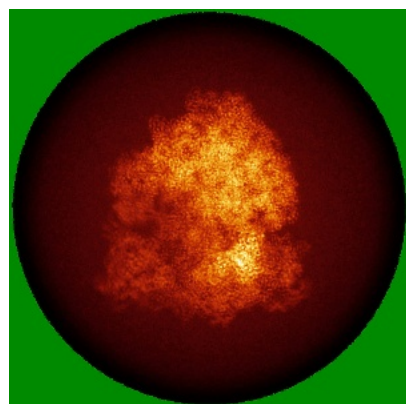


Z Index: 296

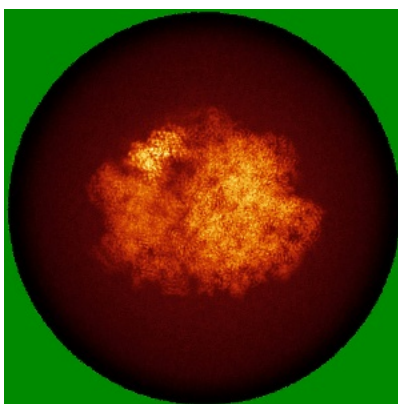
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

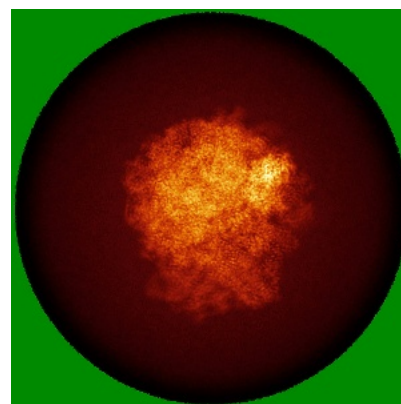
6.4.1 Primary map



X

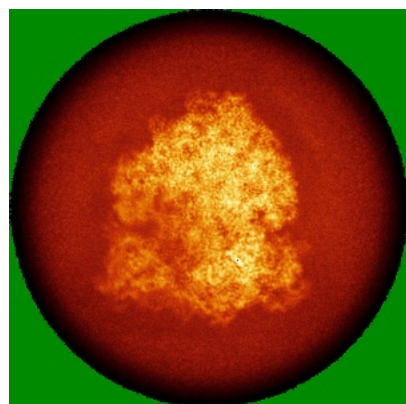


Y

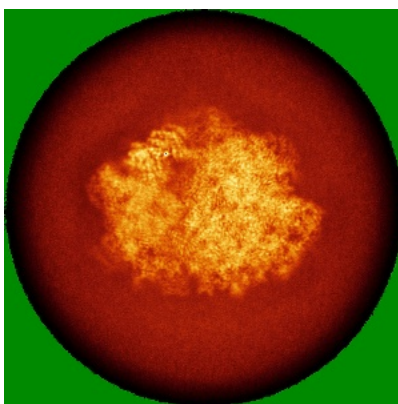


Z

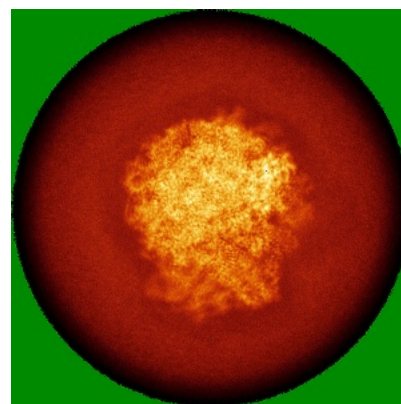
6.4.2 Raw map



X



Y

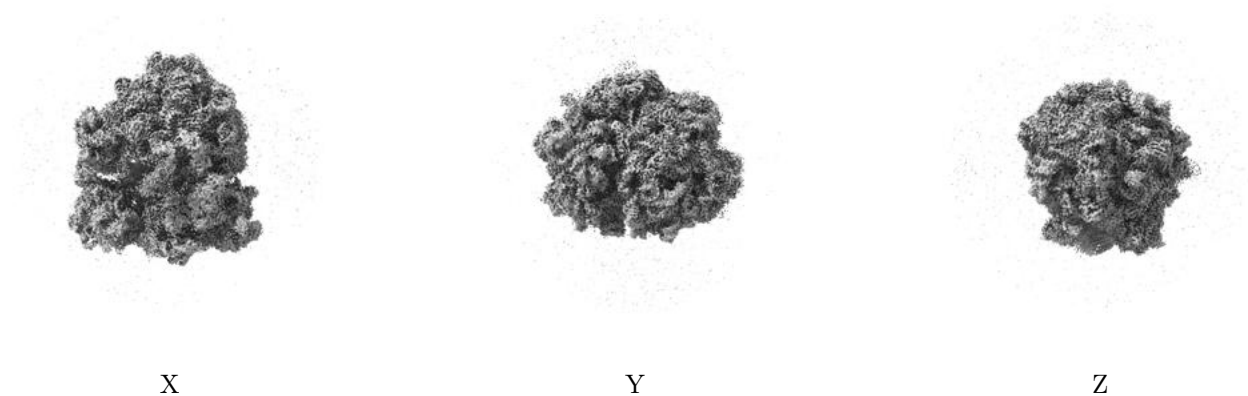


Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

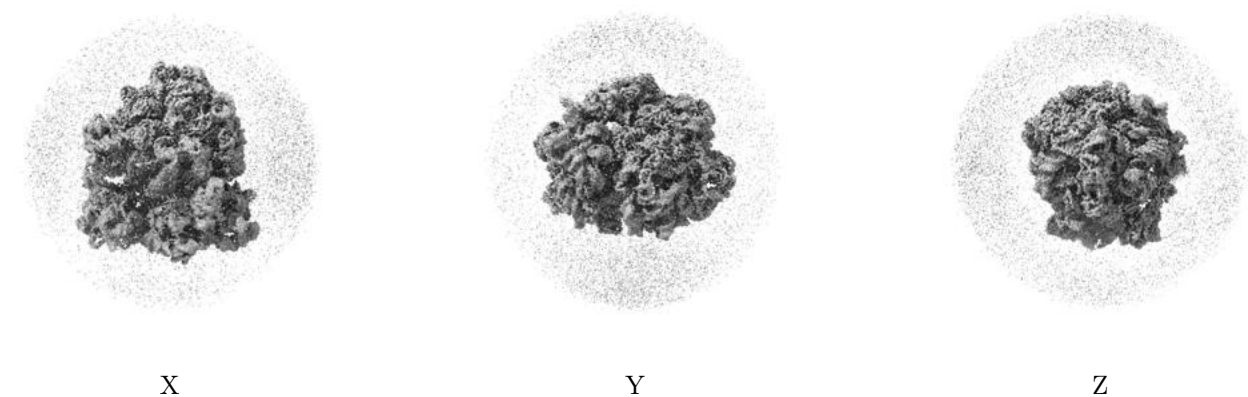
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 3.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

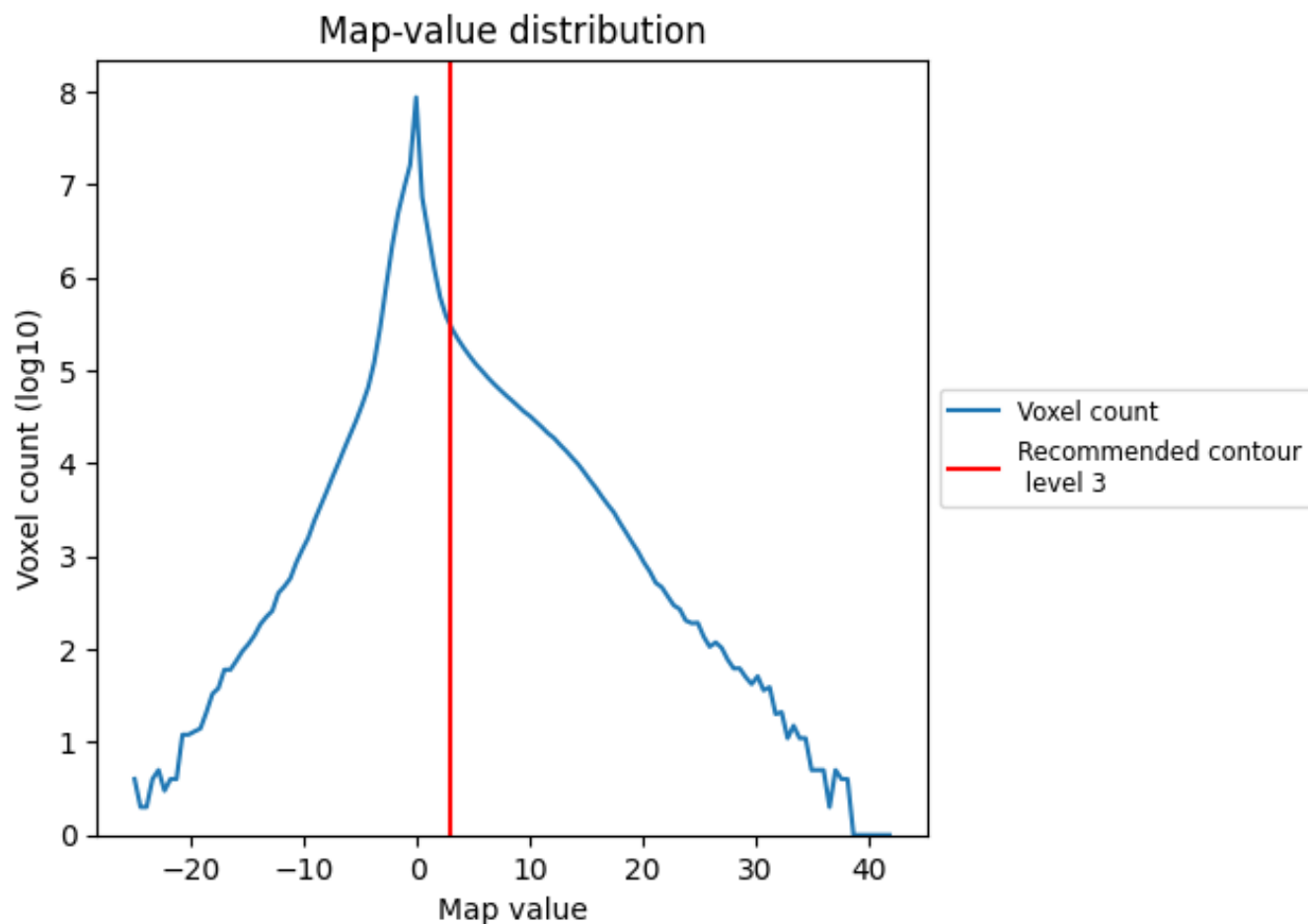
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

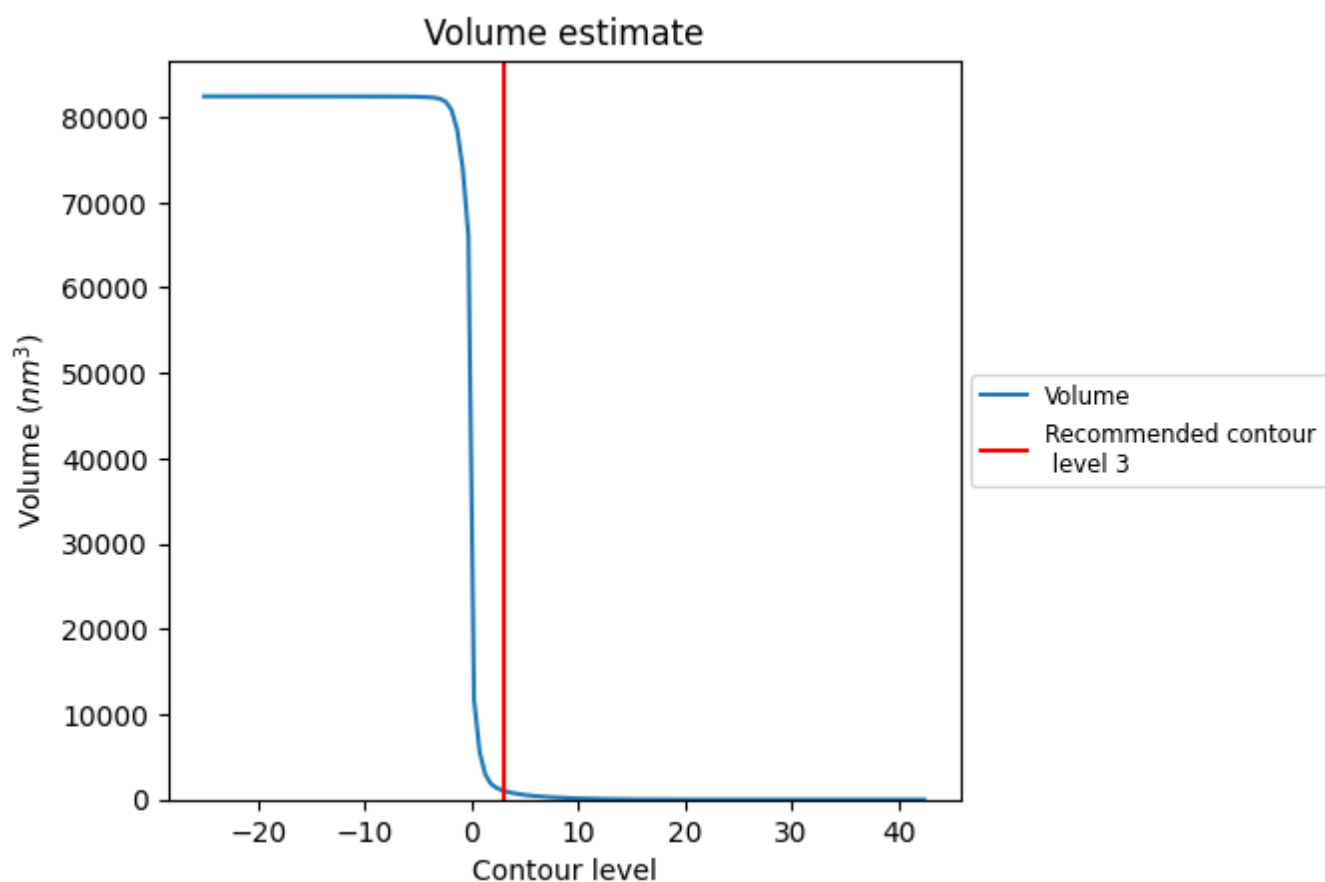
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

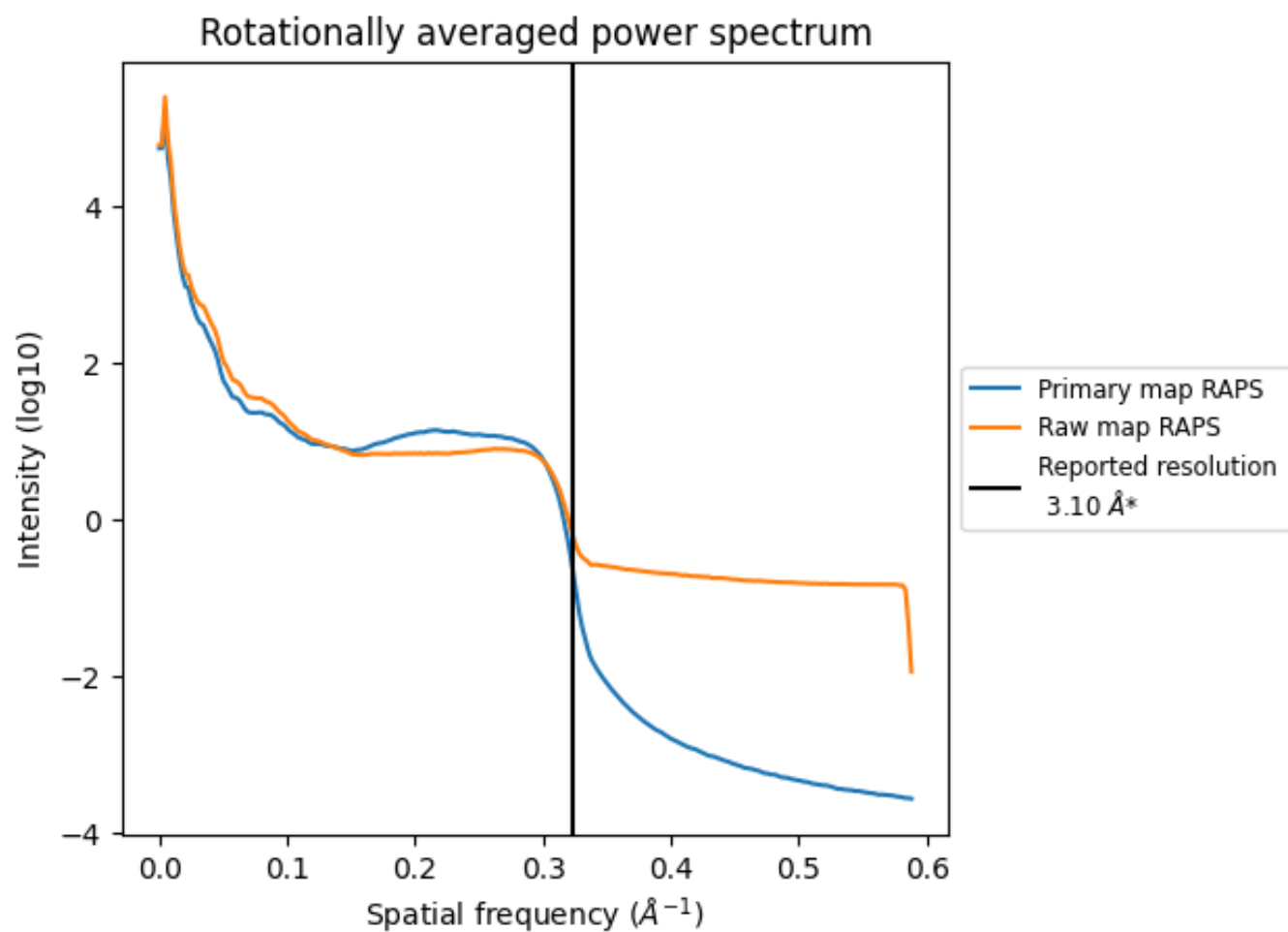
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1050 nm³; this corresponds to an approximate mass of 949 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ

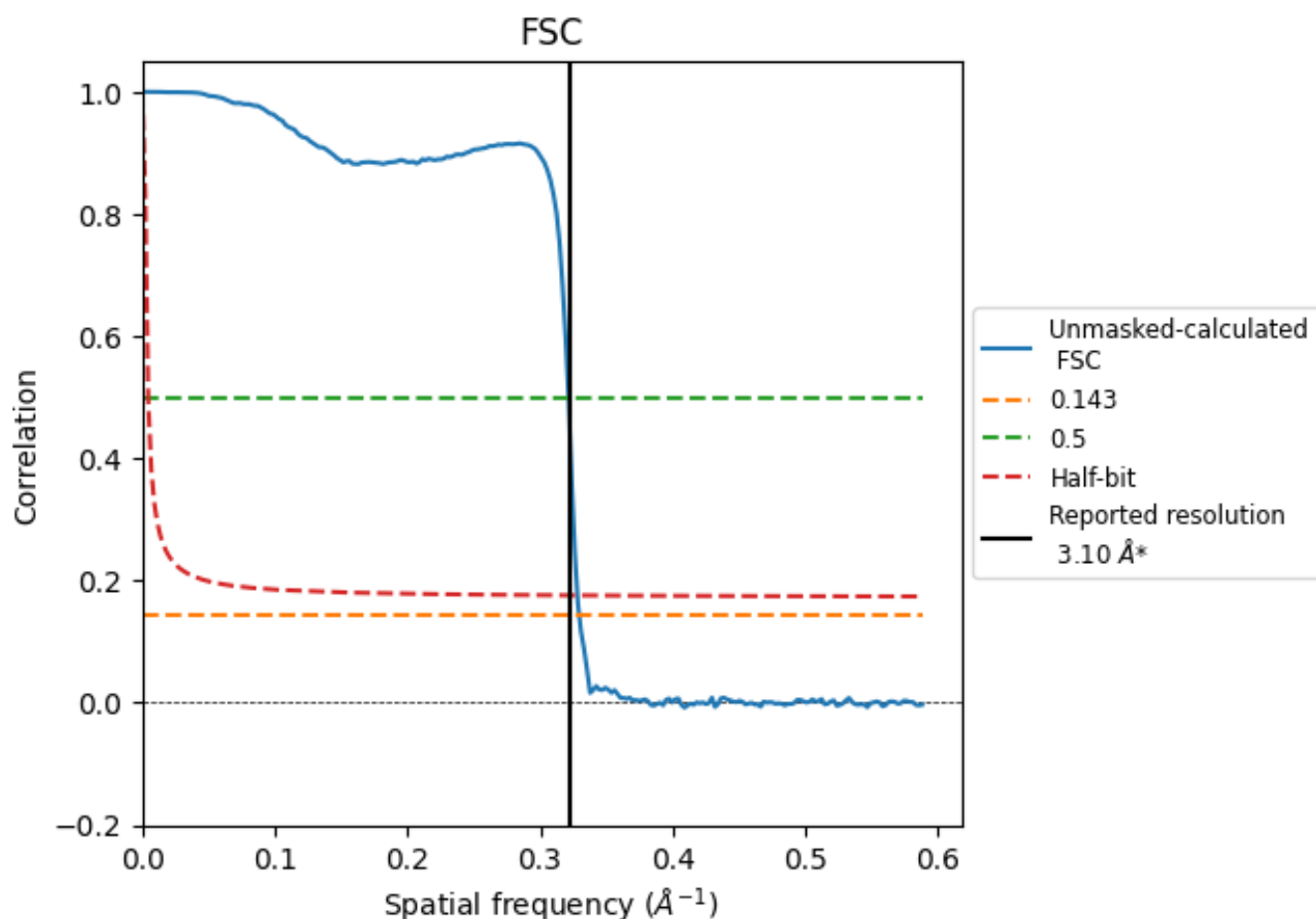


*Reported resolution corresponds to spatial frequency of $0.323 \text{ \AA}^{-1}$

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of $0.323 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

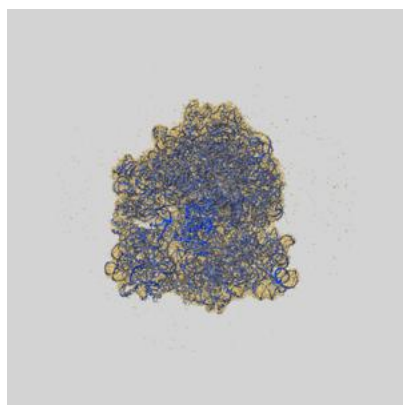
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.10	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.03	3.11	3.05

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

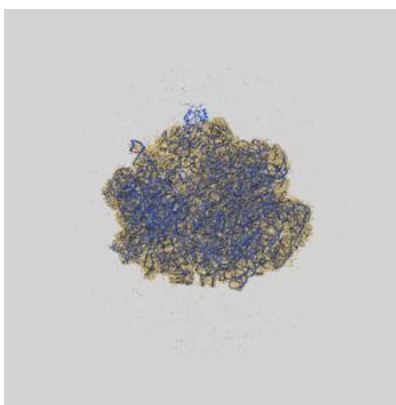
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-29631 and PDB model 8FZI. Per-residue inclusion information can be found in section 3 on page 17.

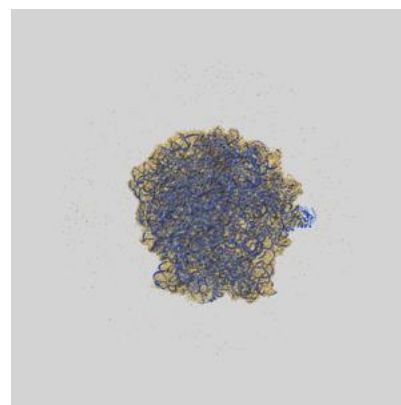
9.1 Map-model overlay [i](#)



X



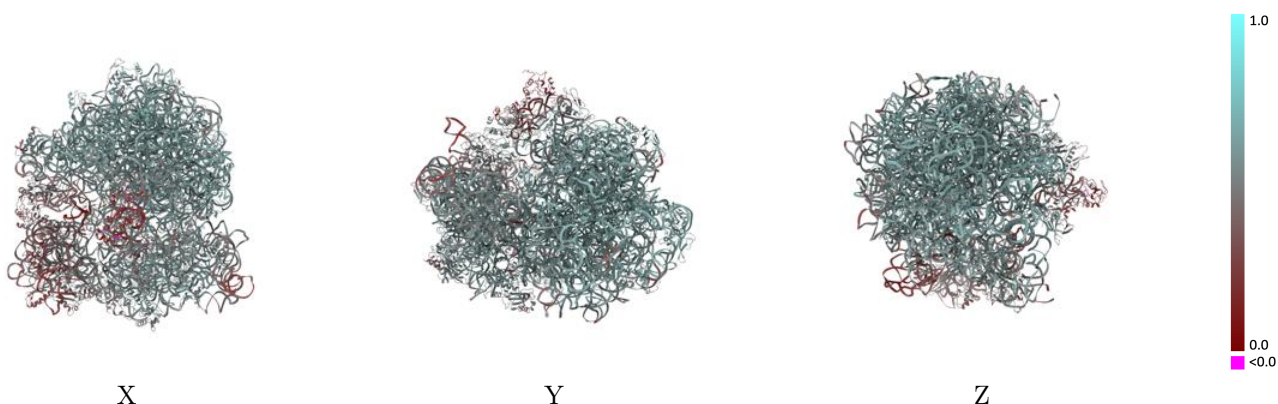
Y



Z

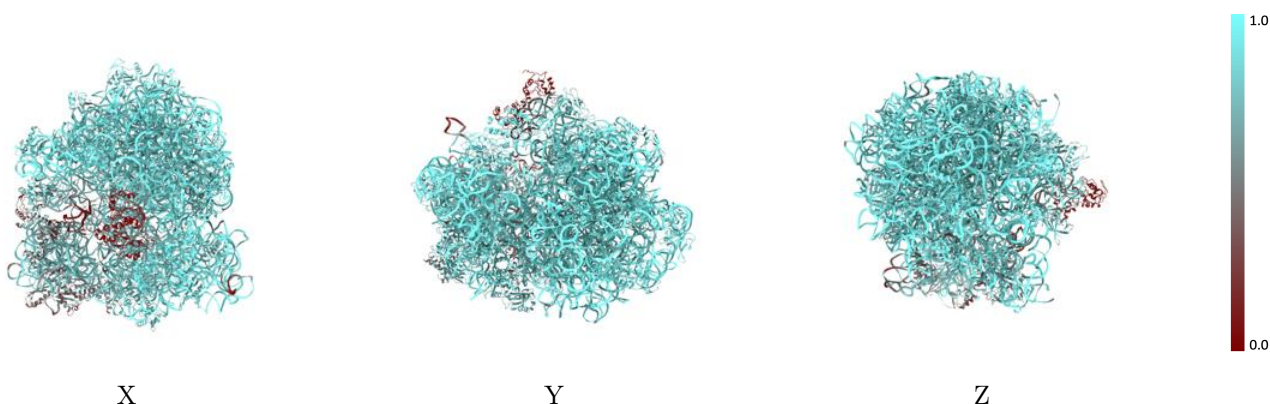
The images above show the 3D surface view of the map at the recommended contour level 3.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



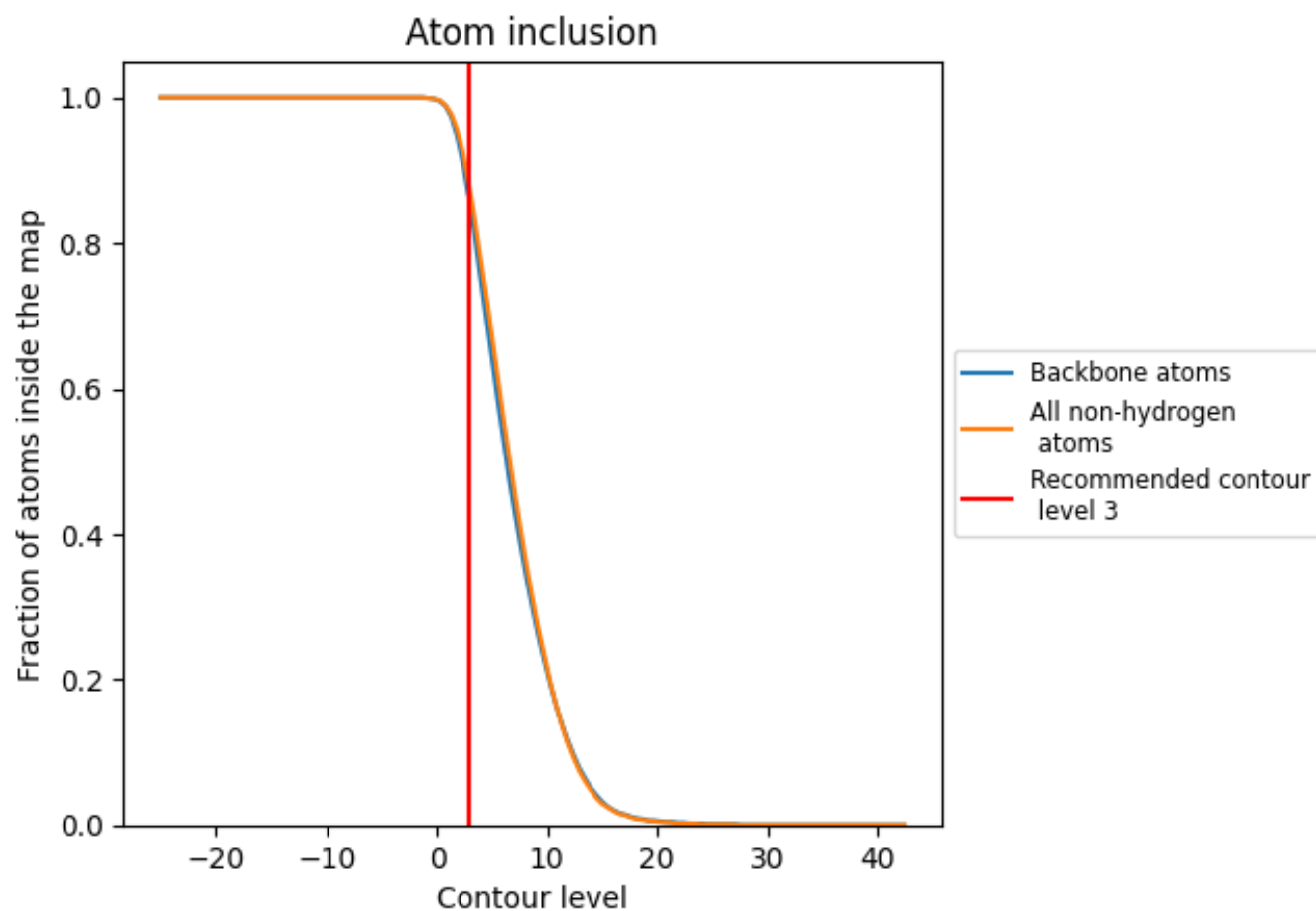
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (3).

9.4 Atom inclusion [i](#)



At the recommended contour level, 86% of all backbone atoms, 88% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (3) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8770	 0.5340
1	 0.8700	 0.5780
2	 0.8970	 0.5900
3	 0.4290	 0.3540
4	 0.9280	 0.6090
5	 0.8580	 0.5760
6	 0.9490	 0.6190
7	 0.9330	 0.6120
8	 0.8740	 0.5720
A	 0.9360	 0.5630
B	 0.9300	 0.5230
C	 0.9340	 0.6080
D	 0.9190	 0.5980
E	 0.8880	 0.5820
F	 0.7250	 0.4430
G	 0.8250	 0.5100
I	 0.0530	 0.2250
J	 0.0990	 0.2380
L	 0.9240	 0.6010
M	 0.9010	 0.5990
N	 0.9210	 0.5920
O	 0.8390	 0.5670
P	 0.9480	 0.6070
Q	 0.8670	 0.5390
R	 0.8710	 0.5840
S	 0.9530	 0.6120
T	 0.9140	 0.5930
U	 0.9170	 0.5950
V	 0.8920	 0.5810
W	 0.8720	 0.5630
X	 0.8370	 0.5570
Y	 0.8990	 0.5960
Z	 0.9220	 0.5930
a	 0.9080	 0.5090
b	 0.6360	 0.4590



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Chain	Atom inclusion	Q-score
c	 0.5640	 0.4450
d	 0.8220	 0.5240
e	 0.8510	 0.5520
f	 0.7840	 0.4920
g	 0.3570	 0.3450
h	 0.8640	 0.5560
i	 0.6990	 0.4380
j	 0.5930	 0.4050
k	 0.8230	 0.5290
l	 0.8470	 0.5650
m	 0.4920	 0.3830
n	 0.5920	 0.4030
o	 0.8300	 0.5430
p	 0.8590	 0.5620
q	 0.8670	 0.5520
r	 0.8010	 0.5360
s	 0.4740	 0.3330
t	 0.8430	 0.5280
u	 0.5850	 0.4930
v	 0.9560	 0.5050
x	 0.7330	 0.3880
z	 0.6570	 0.5030