



Full wwPDB EM Validation Report ⓘ

Jul 8, 2026 – 12:36 PM JST

PDB ID : 9KMT / pdb_00009kmt
EMDB ID : EMD-62444
Title : Bat SARSr-CoV RaTG15 Nsp1 bound to the Human 40S Ribosomal subunit-State1
Authors : Yuan, S.; Yan, R.; Wu, M.
Deposited on : 2024-11-18
Resolution : 2.81 Å(reported)

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

EMDB validation analysis : 0.0.1.dev133
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMD 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

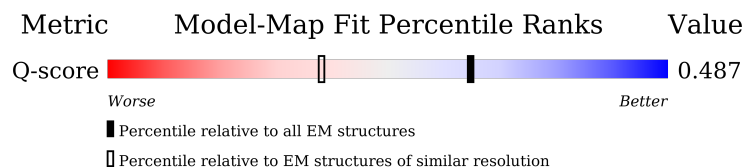
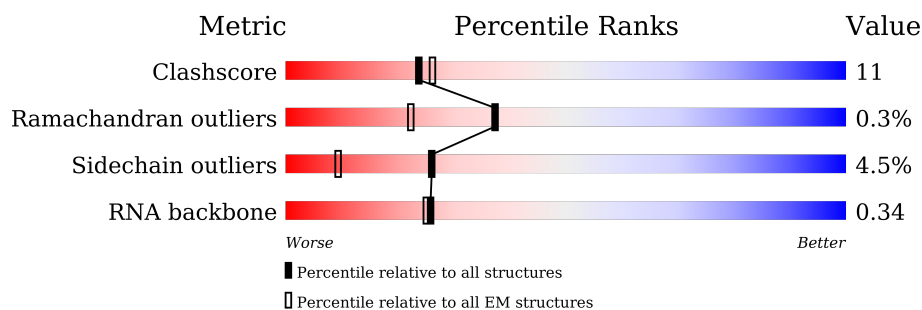
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.81 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.



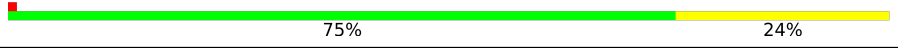
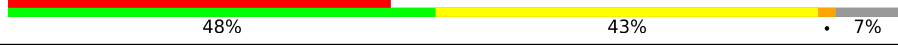
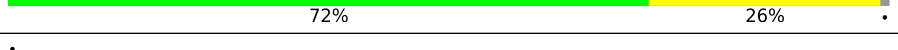
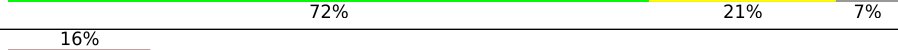
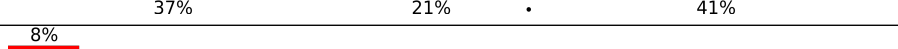
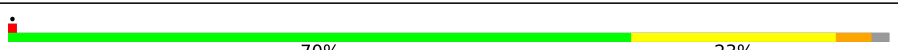
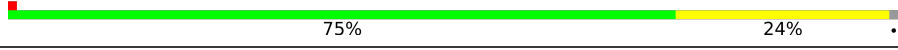
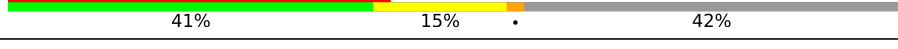
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	11740 (2.31 - 3.31)

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	2	1869	9% 35% 41% 13% 11%
2	A	295	56% 15% 28%
3	B	264	59% 20% 19%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
4	C	293	
5	D	243	
6	E	263	
7	F	204	
8	G	249	
9	H	194	
10	I	208	
11	J	194	
12	K	165	
13	L	158	
14	M	132	
15	N	151	
16	O	151	
17	P	145	
18	Q	146	
19	R	135	
20	S	152	
21	T	145	
22	U	119	
23	V	83	
24	W	130	
25	X	143	
26	Y	130	
27	Z	125	
28	a	101	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
29	b	82	5% 79% 21%
30	c	62	55% 44% 52% 5%
31	d	55	15% 69% 27%
32	e	56	12% 84% 16%
33	f	74	93% 81% 18%
34	g	315	52% 66% 32%
35	h	25	76% 12% 12%
36	n	180	16% 81%

2 Entry composition

There are 36 unique types of molecules in this entry. The entry contains 74832 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 18S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	2	1671	Total	C	N	O	P	0	0
			35647	15925	6406	11655	1661		

- Molecule 2 is a protein called Small ribosomal subunit protein uS2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	A	213	Total	C	N	O	S	0	0
			1686	1072	295	311	8		

- Molecule 3 is a protein called Small ribosomal subunit protein eS1.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	B	213	Total	C	N	O	S	0	0
			1729	1098	309	308	14		

- Molecule 4 is a protein called Small ribosomal subunit protein uS5.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	C	218	Total	C	N	O	S	0	0
			1690	1094	289	297	10		

- Molecule 5 is a protein called Small ribosomal subunit protein uS3.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	D	225	Total	C	N	O	S	0	0
			1752	1117	315	313	7		

- Molecule 6 is a protein called Small ribosomal subunit protein eS4, X isoform.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	E	262	Total	C	N	O	S	0	0
			2076	1324	386	358	8		

- Molecule 7 is a protein called Small ribosomal subunit protein uS7.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	F	189	Total	C	N	O	S	0	0
			1495	934	284	270	7		

- Molecule 8 is a protein called Small ribosomal subunit protein eS6.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	G	230	Total	C	N	O	S	0	0
			1864	1164	373	320	7		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	221	ARG	LYS	variant	UNP P62753

- Molecule 9 is a protein called Small ribosomal subunit protein eS7.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	H	186	Total	C	N	O	S	0	0
			1501	957	276	267	1		

- Molecule 10 is a protein called Small ribosomal subunit protein eS8.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	I	205	Total	C	N	O	S	0	0
			1682	1056	331	290	5		

- Molecule 11 is a protein called Small ribosomal subunit protein uS4.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	J	180	Total	C	N	O	S	0	0
			1499	955	300	242	2		

- Molecule 12 is a protein called Small ribosomal subunit protein eS10.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	K	97	Total	C	N	O	S	0	0
			816	533	144	133	6		

- Molecule 13 is a protein called Small ribosomal subunit protein uS17.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	L	151	Total	C	N	O	S	0	0
			1229	782	230	211	6		

- Molecule 14 is a protein called Small ribosomal subunit protein eS12.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	M	121	Total	C	N	O	S	0	0
			935	586	165	175	9		

- Molecule 15 is a protein called Small ribosomal subunit protein uS15.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	N	149	Total	C	N	O	S	0	0
			1202	770	228	203	1		

- Molecule 16 is a protein called Small ribosomal subunit protein uS11.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	O	135	Total	C	N	O	S	0	0
			1010	618	198	188	6		

- Molecule 17 is a protein called Small ribosomal subunit protein uS19.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	P	126	Total	C	N	O	S	0	0
			1037	659	196	175	7		

- Molecule 18 is a protein called Small ribosomal subunit protein uS9.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	Q	138	Total	C	N	O	S	0	0
			1097	698	206	190	3		

- Molecule 19 is a protein called Small ribosomal subunit protein eS17.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	R	132	Total	C	N	O	S	0	0
			1068	670	199	195	4		

- Molecule 20 is a protein called Small ribosomal subunit protein uS13.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	S	143	Total	C	N	O	S	0	0
			1184	743	240	200	1		

- Molecule 21 is a protein called Small ribosomal subunit protein eS19.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	T	144	Total	C	N	O	S	0	0
			1123	703	217	200	3		

- Molecule 22 is a protein called Small ribosomal subunit protein uS10.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	U	101	Total	C	N	O	S	0	0
			803	504	153	142	4		

- Molecule 23 is a protein called Small ribosomal subunit protein eS21.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	V	82	Total	C	N	O	S	0	0
			625	384	116	120	5		

- Molecule 24 is a protein called Small ribosomal subunit protein uS8.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	W	129	Total	C	N	O	S	0	0
			1034	659	193	176	6		

- Molecule 25 is a protein called Small ribosomal subunit protein uS12.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	X	141	Total	C	N	O	S	0	0
			1098	693	219	183	3		

- Molecule 26 is a protein called Small ribosomal subunit protein eS24.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	Y	124	Total	C	N	O	S	0	0
			1014	641	198	170	5		

- Molecule 27 is a protein called Small ribosomal subunit protein eS25.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	Z	72	Total	C	N	O	S	0	0
			574	368	104	101	1		

- Molecule 28 is a protein called Small ribosomal subunit protein eS26.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	a	99	Total	C	N	O	S	0	0
			794	494	165	130	5		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
a	78	VAL	ALA	conflict	UNP P62854

- Molecule 29 is a protein called Small ribosomal subunit protein eS27.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	b	82	Total	C	N	O	S	0	0
			641	402	118	114	7		

- Molecule 30 is a protein called Small ribosomal subunit protein eS28.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	c	62	Total	C	N	O	S	0	0
			489	297	97	93	2		

- Molecule 31 is a protein called Small ribosomal subunit protein uS14.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	d	55	Total	C	N	O	S	0	0
			459	286	94	74	5		

- Molecule 32 is a protein called Small ribosomal subunit protein eS30.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	e	56	Total	C	N	O	S	0	0
			442	273	96	72	1		

- Molecule 33 is a protein called Small ribosomal subunit protein eS31.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	f	74	Total	C	N	O	S	0	0
			611	385	117	102	7		

- Molecule 34 is a protein called Small ribosomal subunit protein RACK1.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	g	314	Total	C	N	O	S	0	0
			2441	1537	425	467	12		

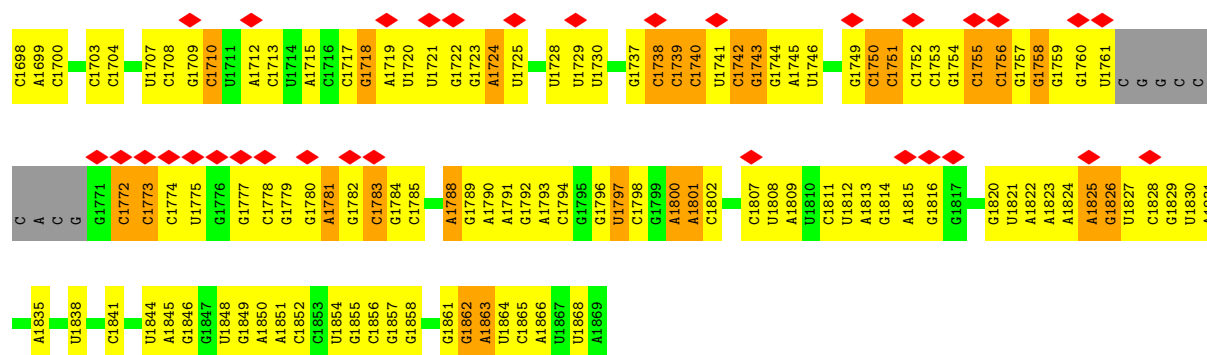
- Molecule 35 is a protein called Small ribosomal subunit protein eS32.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	h	22	Total	C	N	O	S	0	0
			213	130	57	23	3		

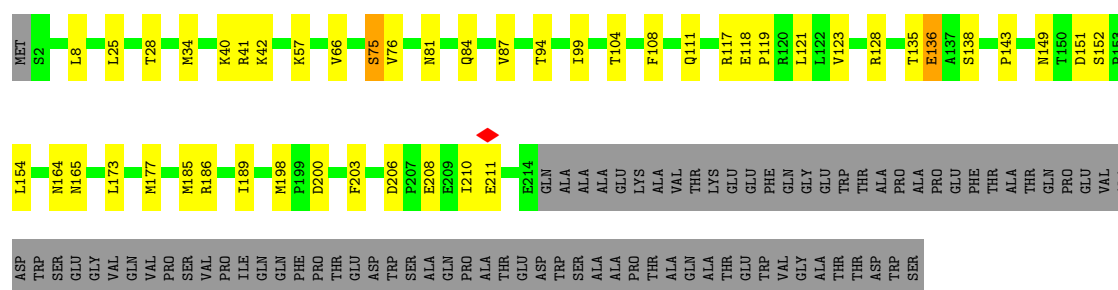
- Molecule 36 is a protein called ORF1ab polyprotein.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	n	35	Total	C	N	O	S	0	0
			272	167	46	57	2		

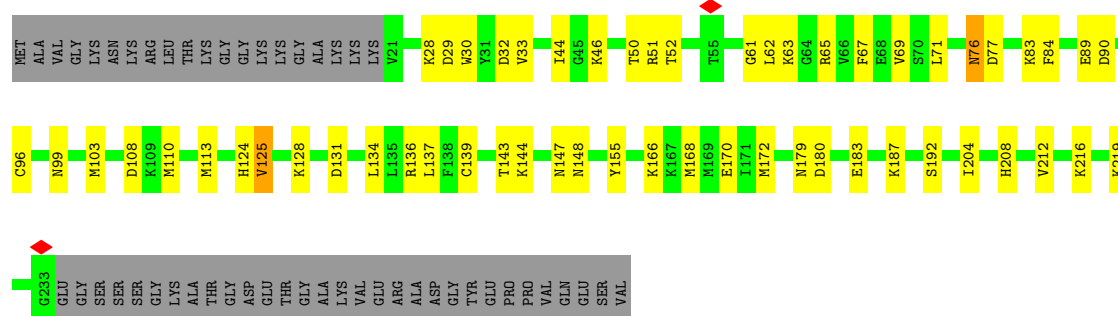




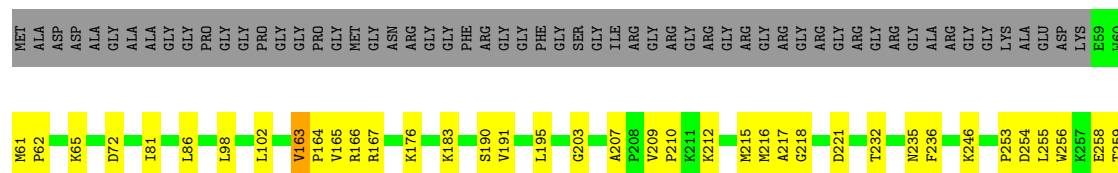
- Molecule 2: Small ribosomal subunit protein uS2

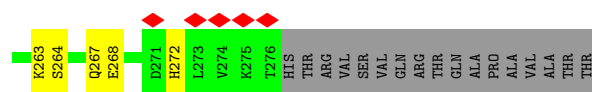


- Molecule 3: Small ribosomal subunit protein eS1

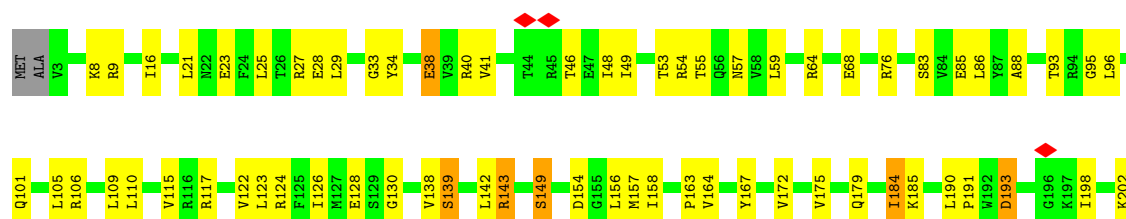


- Molecule 4: Small ribosomal subunit protein uS5

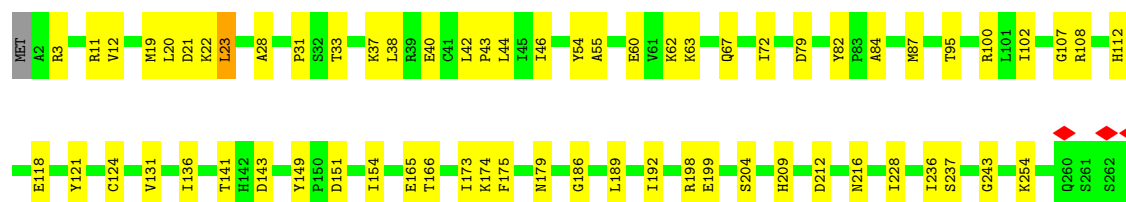
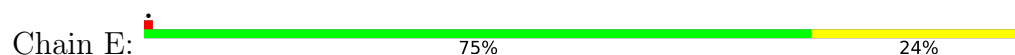




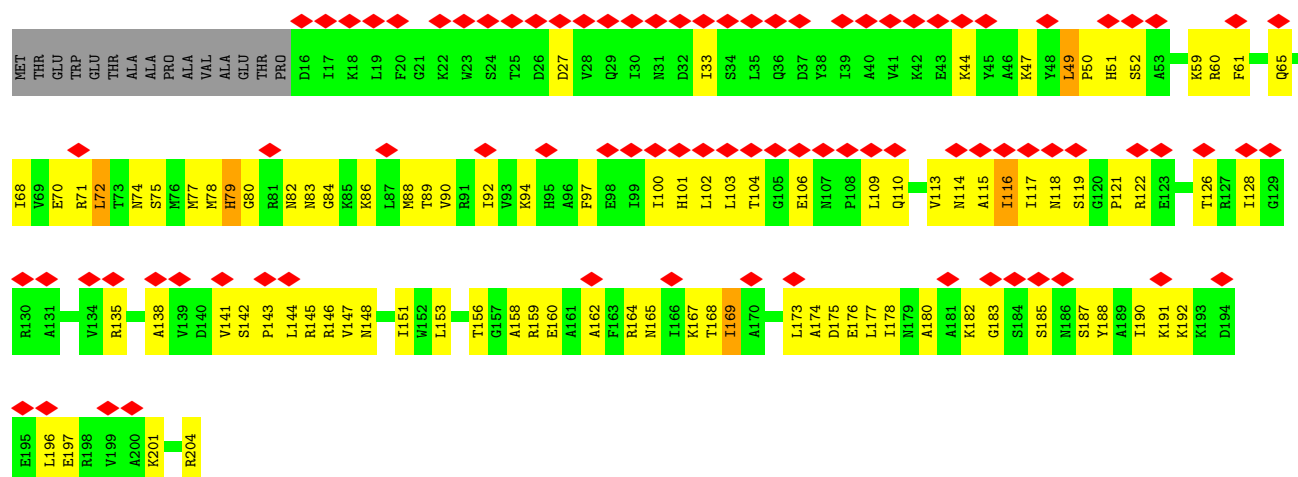
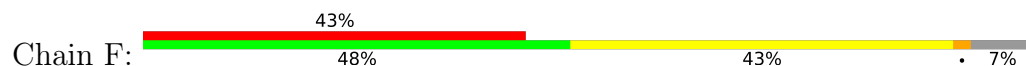
- Molecule 5: Small ribosomal subunit protein uS3



- Molecule 6: Small ribosomal subunit protein eS4, X isoform

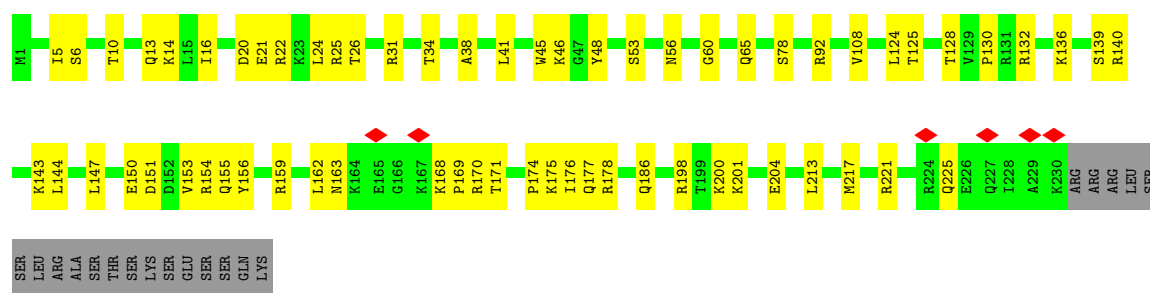


- Molecule 7: Small ribosomal subunit protein uS7



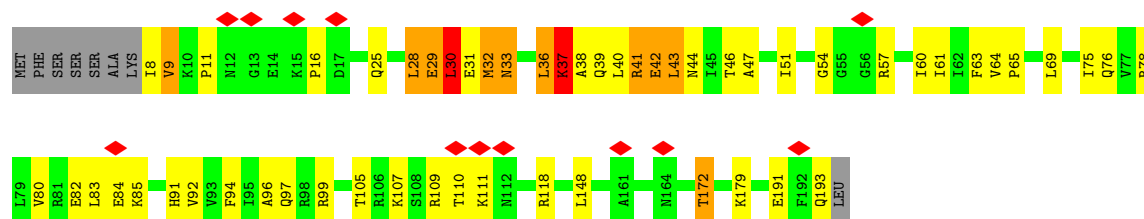
- Molecule 8: Small ribosomal subunit protein eS6

Chain G: 



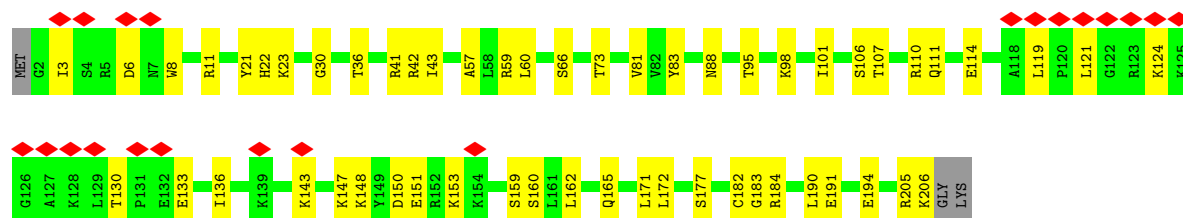
- Molecule 9: Small ribosomal subunit protein eS7

Chain H: 



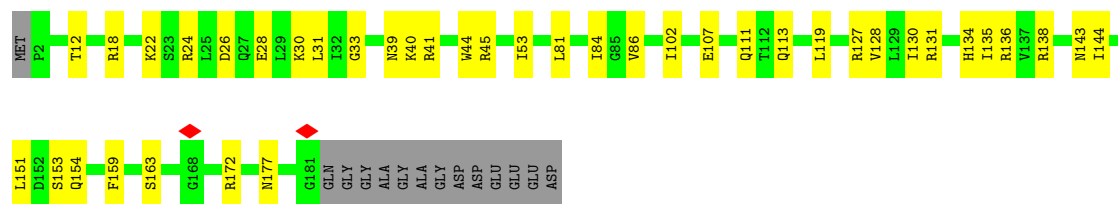
- Molecule 10: Small ribosomal subunit protein eS8

Chain I: 

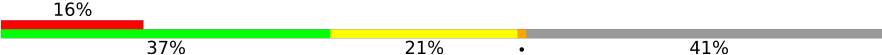


- Molecule 11: Small ribosomal subunit protein uS4

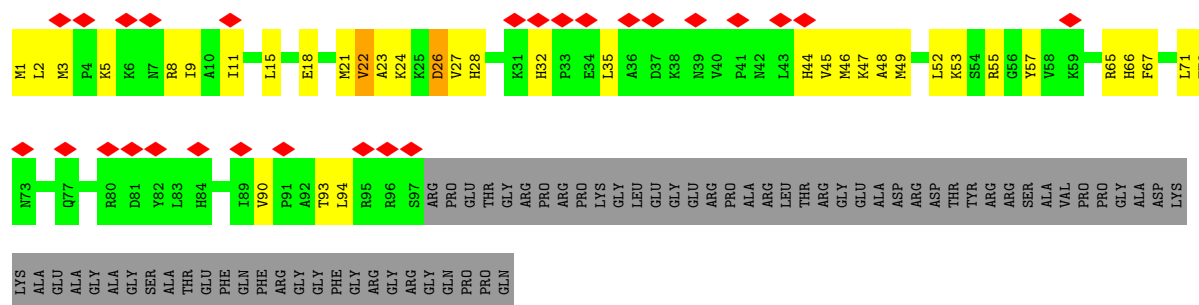
Chain J: 



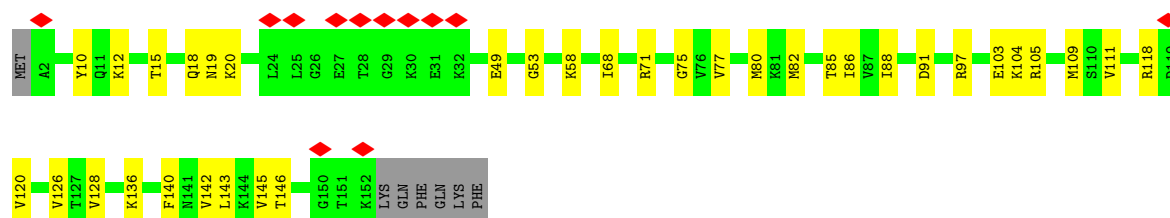
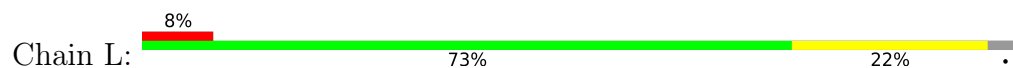
- Molecule 12: Small ribosomal subunit protein eS10

Chain K: 

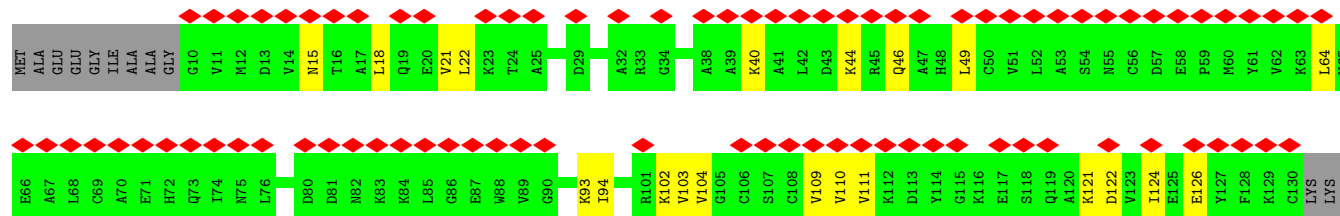
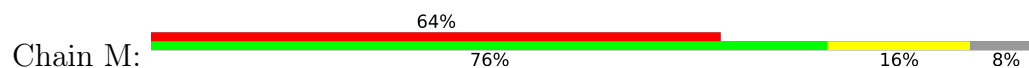




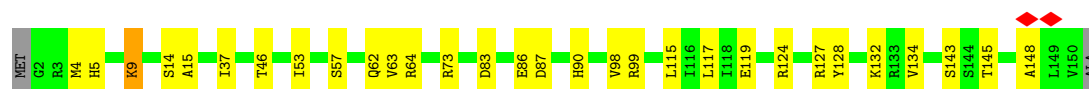
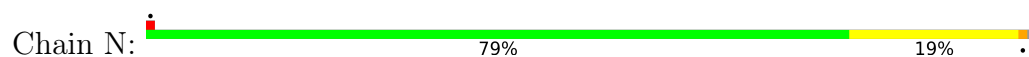
- Molecule 13: Small ribosomal subunit protein uS17



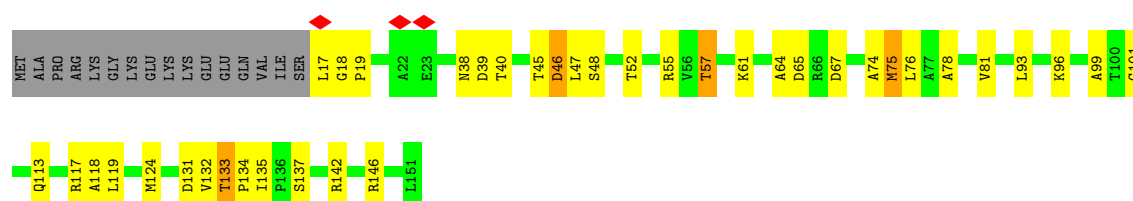
- Molecule 14: Small ribosomal subunit protein eS12



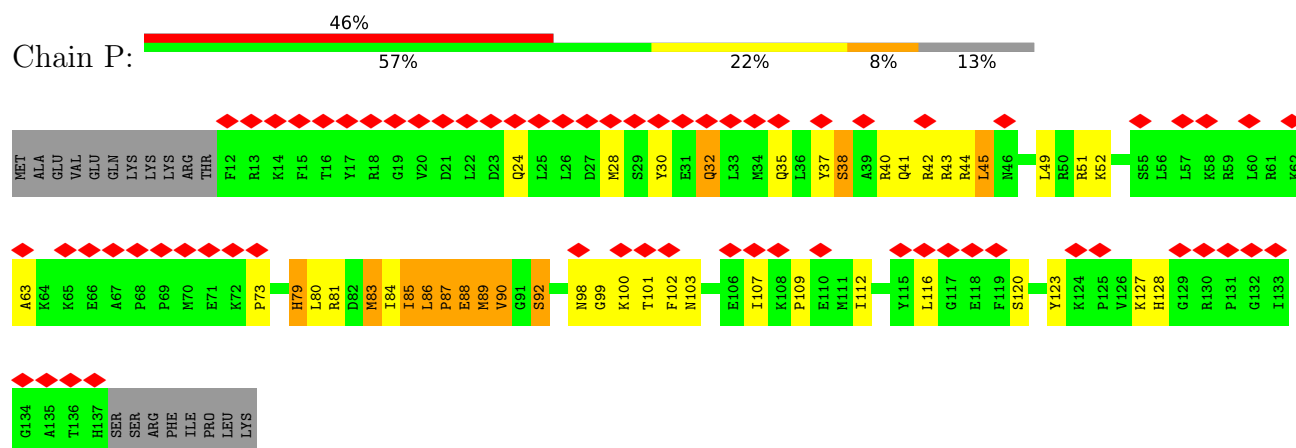
- Molecule 15: Small ribosomal subunit protein uS15



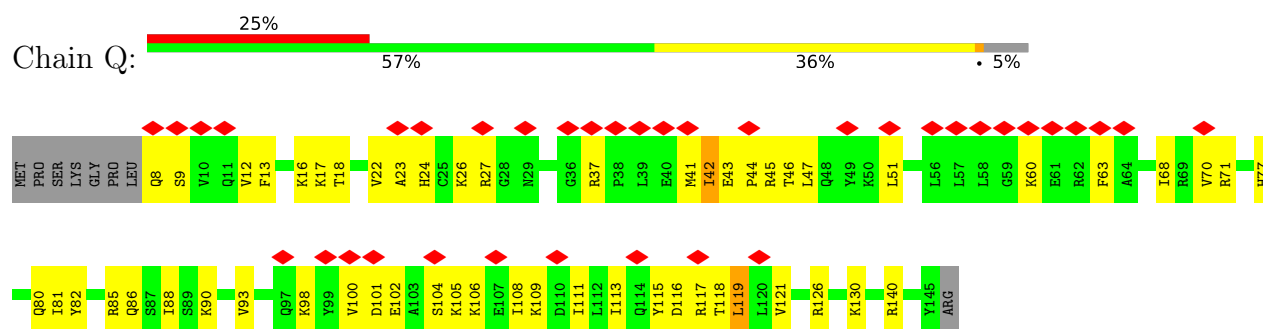
- Molecule 16: Small ribosomal subunit protein uS11



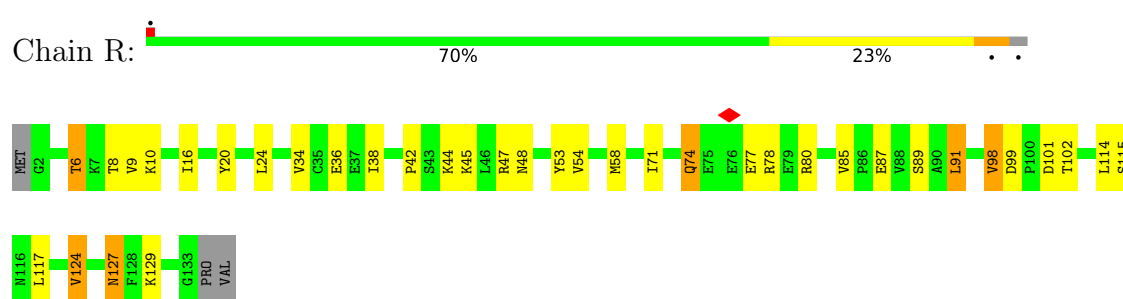
- Molecule 17: Small ribosomal subunit protein uS19



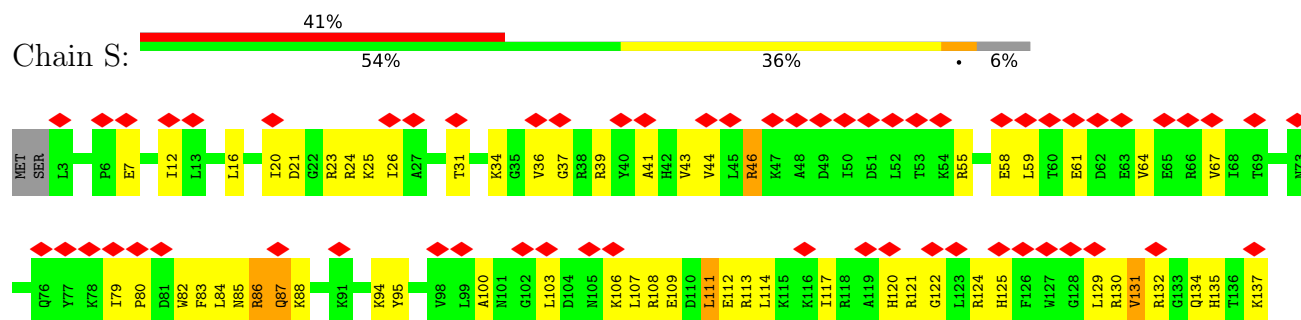
- Molecule 18: Small ribosomal subunit protein uS9

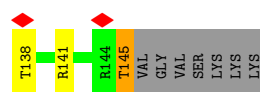


- Molecule 19: Small ribosomal subunit protein eS17

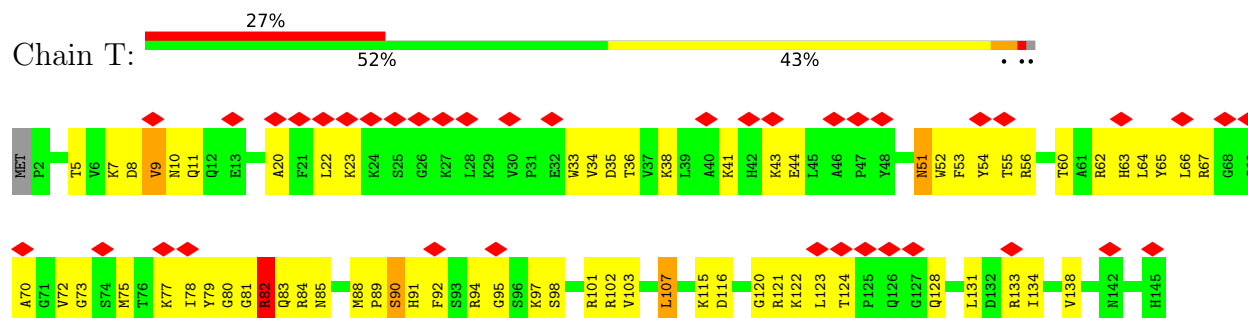


- Molecule 20: Small ribosomal subunit protein uS13

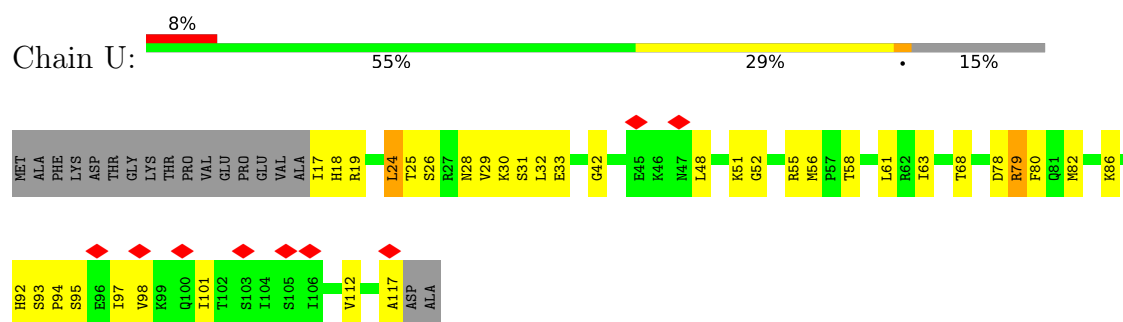




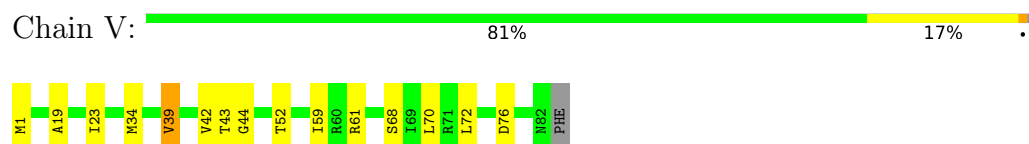
- Molecule 21: Small ribosomal subunit protein eS19



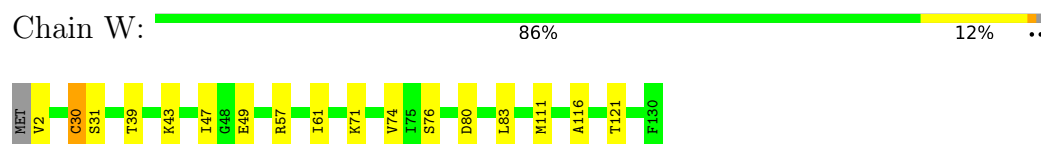
- Molecule 22: Small ribosomal subunit protein uS10



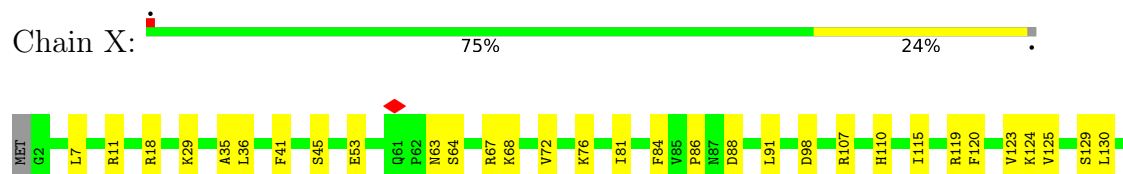
- Molecule 23: Small ribosomal subunit protein eS21

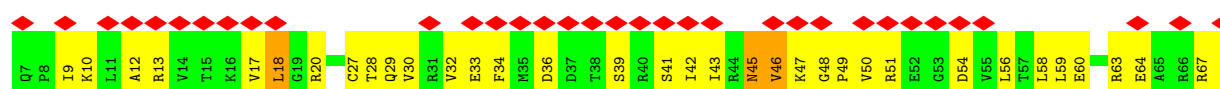


- Molecule 24: Small ribosomal subunit protein uS8

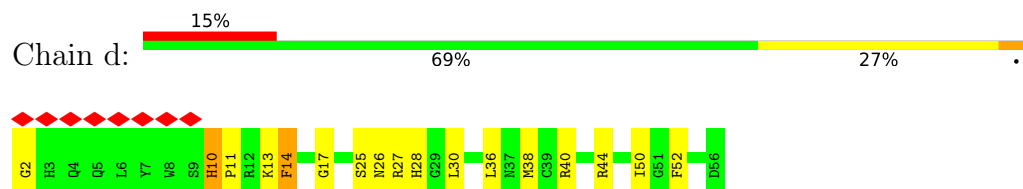


- Molecule 25: Small ribosomal subunit protein uS12

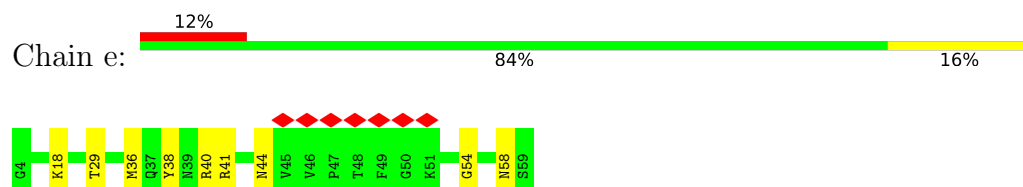




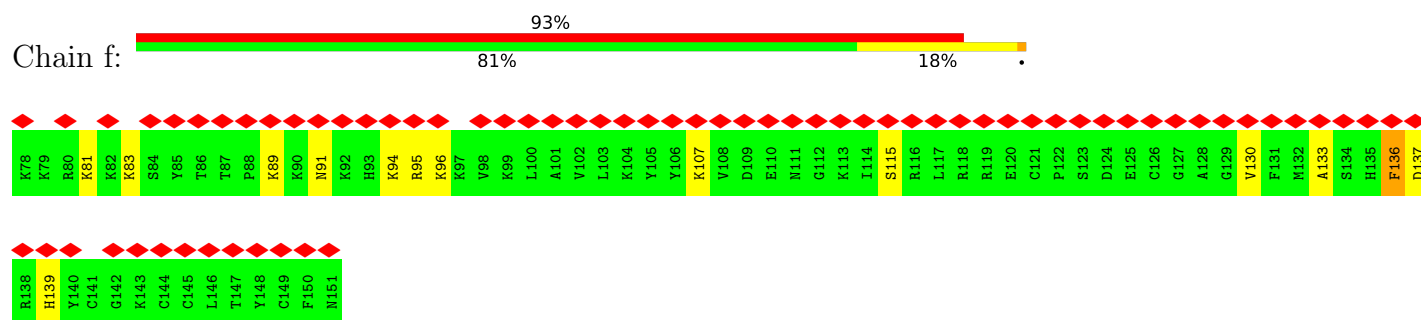
- Molecule 31: Small ribosomal subunit protein uS14



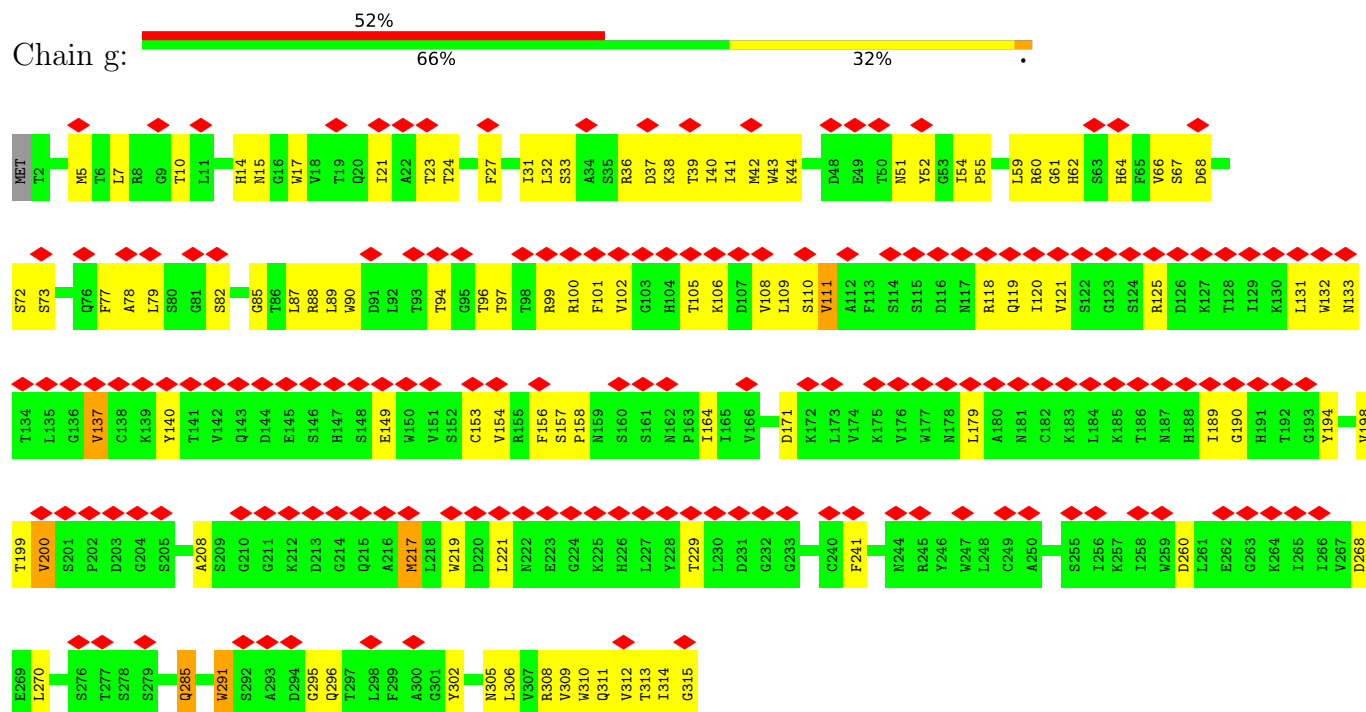
- Molecule 32: Small ribosomal subunit protein eS30



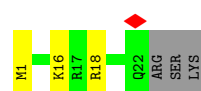
- Molecule 33: Small ribosomal subunit protein eS31



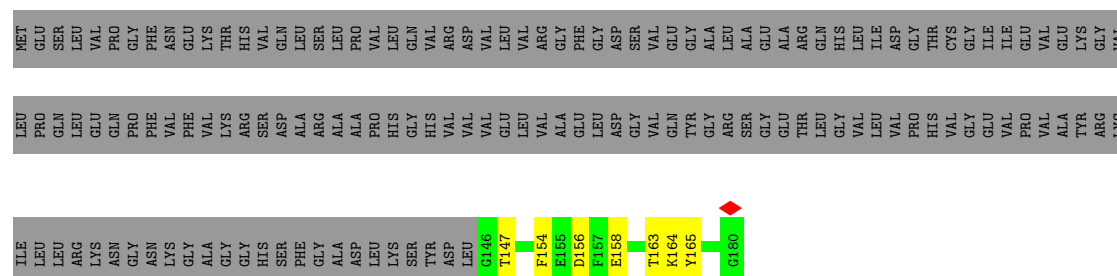
- Molecule 34: Small ribosomal subunit protein RACK1



- Molecule 35: Small ribosomal subunit protein eS32



Chain n: 16% 81%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	80699	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	JEOL CRYO ARM 300	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	2.042	Depositor
Minimum map value	-0.961	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.057	Depositor
Recommended contour level	0.177	Depositor
Map size ($\text{\AA}$)	486.4, 486.4, 486.4	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.95, 0.95, 0.95	Depositor

5 Model quality

5.1 Standard geometry

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	2	0.18	0/39865	0.33	10/62134 (0.0%)
2	A	0.15	0/1723	0.28	0/2341
3	B	0.16	0/1756	0.30	0/2350
4	C	0.16	0/1726	0.34	0/2332
5	D	0.16	0/1780	0.39	0/2397
6	E	0.16	0/2118	0.30	0/2849
7	F	0.21	0/1516	0.57	0/2037
8	G	0.13	0/1887	0.30	0/2513
9	H	0.29	0/1524	0.48	1/2042 (0.0%)
10	I	0.14	0/1711	0.33	0/2282
11	J	0.15	0/1524	0.30	0/2035
12	K	0.18	0/840	0.47	0/1133
13	L	0.16	0/1250	0.31	0/1673
14	M	0.12	0/945	0.33	0/1269
15	N	0.14	0/1226	0.27	0/1649
16	O	0.16	0/1023	0.37	0/1372
17	P	0.30	0/1058	0.61	0/1414
18	Q	0.19	0/1114	0.51	0/1492
19	R	0.15	0/1082	0.36	0/1452
20	S	0.16	0/1202	0.44	0/1610
21	T	0.21	0/1143	0.52	0/1530
22	U	0.15	0/813	0.34	0/1092
23	V	0.14	0/631	0.28	0/844
24	W	0.18	0/1051	0.33	0/1406
25	X	0.16	0/1116	0.31	0/1490
26	Y	0.15	0/1031	0.28	0/1370
27	Z	0.13	0/580	0.37	0/780
28	a	0.16	0/807	0.31	0/1082
29	b	0.15	0/654	0.32	0/876
30	c	0.17	0/491	0.47	0/656
31	d	0.16	0/470	0.33	0/623
32	e	0.16	0/447	0.32	0/587
33	f	0.11	0/623	0.34	0/822
34	g	0.16	0/2498	0.41	1/3399 (0.0%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
35	h	0.13	0/214	0.22	0/272
36	n	0.20	0/276	0.38	0/371
All	All	0.17	0/79715	0.35	12/115576 (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
20	S	0	1
21	T	0	1
All	All	0	2

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	2	579	C	O3'-P-O5'	-12.19	85.71	104.00
1	2	550	C	O3'-P-O5'	-8.81	90.79	104.00
1	2	804	U	O3'-P-O5'	-8.62	91.06	104.00
1	2	500	A	P-O3'-C3'	8.25	132.58	120.20
1	2	1022	U	C4'-C3'-C2'	-7.58	95.02	102.60
1	2	1231	C	C2'-C3'-O3'	-7.47	102.49	113.70
1	2	804	U	OP2-P-O3'	6.23	126.68	108.00
1	2	1230	C	O3'-P-O5'	6.10	113.14	104.00
1	2	579	C	C4'-C3'-C2'	-5.87	96.73	102.60
34	g	55	PRO	CA-N-CD	-5.84	103.83	112.00
1	2	500	A	OP1-P-O3'	5.48	124.45	108.00
9	H	37	LYS	N-CA-C	-5.08	105.87	111.71

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
20	S	135	HIS	Peptide
21	T	82	ARG	Peptide

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	2	35647	0	18022	593	0
2	A	1686	0	1688	28	0
3	B	1729	0	1803	31	0
4	C	1690	0	1777	27	0
5	D	1752	0	1848	48	0
6	E	2076	0	2177	46	0
7	F	1495	0	1549	98	0
8	G	1864	0	2018	52	0
9	H	1501	0	1593	46	0
10	I	1682	0	1769	39	0
11	J	1499	0	1618	31	0
12	K	816	0	841	35	0
13	L	1229	0	1302	24	0
14	M	935	0	964	12	0
15	N	1202	0	1289	20	0
16	O	1010	0	1034	34	0
17	P	1037	0	1082	37	0
18	Q	1097	0	1161	46	0
19	R	1068	0	1121	33	0
20	S	1184	0	1244	56	0
21	T	1123	0	1153	89	0
22	U	803	0	873	27	0
23	V	625	0	628	10	0
24	W	1034	0	1080	9	0
25	X	1098	0	1167	24	0
26	Y	1014	0	1082	22	0
27	Z	574	0	627	16	0
28	a	794	0	849	19	0
29	b	641	0	665	15	0
30	c	489	0	514	35	0
31	d	459	0	452	16	0
32	e	442	0	487	8	0
33	f	611	0	638	15	0
34	g	2441	0	2396	68	0
35	h	213	0	258	3	0
36	n	272	0	248	5	0
All	All	74832	0	59017	1439	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

All (1439) 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)
1:2:1607:A:N6	1:2:1632:G:N2	1.99	1.09
1:2:1607:A:N6	1:2:1632:G:H21	1.48	1.07
18:Q:22:VAL:O	18:Q:70:VAL:HA	1.56	1.04
1:2:1743:G:H21	1:2:1791:A:N6	1.58	1.01
20:S:130:ARG:HE	20:S:134:GLN:HB2	1.28	0.99
1:2:1743:G:N2	1:2:1791:A:C6	2.35	0.94
34:g:217:MET:HB3	34:g:229:THR:HG22	1.51	0.93
30:c:28:THR:HB	30:c:46:VAL:O	1.71	0.90
1:2:13:C:H5	1:2:1198:G:H1	1.18	0.90
27:Z:98:LYS:HB3	27:Z:110:THR:O	1.74	0.87
7:F:158:ALA:O	7:F:162:ALA:HB3	1.76	0.86
7:F:90:VAL:O	7:F:94:LYS:HB3	1.76	0.85
1:2:1327:G:H1	1:2:1502:C:H5	1.23	0.85
1:2:1541:G:H22	1:2:1591:C:H42	1.18	0.84
1:2:1607:A:H62	1:2:1632:G:N2	1.73	0.84
7:F:117:ILE:H	7:F:183:GLY:H	1.25	0.83
2:A:81:ASN:HA	2:A:84:GLN:HG3	1.60	0.83
9:H:69:LEU:HD22	9:H:96:ALA:HB2	1.61	0.81
1:2:1788:A:H62	1:2:1789:G:H21	1.29	0.81
1:2:1608:U:H4'	20:S:130:ARG:HH22	1.46	0.80
1:2:1589:A:OP1	21:T:83:GLN:NE2	2.14	0.80
1:2:125:C:O2	1:2:339:A:N6	2.15	0.79
9:H:29:GLU:HG2	9:H:33:ASN:HD21	1.47	0.79
21:T:116:ASP:HB3	21:T:122:LYS:HB3	1.63	0.78
1:2:1532:C:N4	7:F:79:HIS:O	2.17	0.78
1:2:1540:G:N2	1:2:1589:A:OP2	2.16	0.78
1:2:1513:C:H42	31:d:10:HIS:HB2	1.50	0.77
7:F:119:SER:HB3	7:F:190:ILE:HG21	1.65	0.77
1:2:1628:C:O2'	20:S:39:ARG:NH2	2.17	0.77
1:2:928:G:H1	1:2:1013:U:H3	1.31	0.77
1:2:1610:G:N2	20:S:85:ASN:O	2.18	0.77
1:2:925:G:H1	1:2:1017:U:H3	1.33	0.76
1:2:1569:A:O4'	21:T:94:ARG:NH1	2.18	0.76
3:B:128:LYS:HE3	3:B:134:LEU:HB2	1.66	0.76
1:2:1243:U:H3	1:2:1264:C:N4	1.82	0.76
7:F:138:ALA:O	30:c:63:ARG:NH2	2.18	0.76
1:2:1243:U:N3	1:2:1264:C:N4	2.34	0.75

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1611:G:H1	1:2:1627:C:H42	1.34	0.75
9:H:31:GLU:HB3	9:H:32:MET:HE3	1.68	0.75
1:2:1223:A:H61	1:2:1644:C:H42	1.31	0.75
21:T:72:VAL:HG13	21:T:95:GLY:HA3	1.67	0.75
1:2:1587:G:H1	21:T:97:LYS:HE2	1.51	0.75
1:2:1260:A:OP2	1:2:1623:A:N6	2.17	0.74
1:2:1143:A:H5'	4:C:190:SER:HB3	1.70	0.74
1:2:1743:G:N2	1:2:1791:A:N6	2.33	0.74
1:2:588:G:H4'	1:2:589:G:H5'	1.70	0.74
1:2:1862:G:O2'	28:a:5:ARG:NH2	2.21	0.74
7:F:121:PRO:O	7:F:146:ARG:NH1	2.21	0.74
22:U:33:GLU:OE2	22:U:55:ARG:NH1	2.21	0.73
1:2:145:G:N2	1:2:146:G:O6	2.19	0.73
1:2:1590:C:H1'	21:T:81:GLY:HA3	1.68	0.73
7:F:147:VAL:HG23	7:F:151:ILE:HD13	1.68	0.73
1:2:1274:G:OP2	12:K:1:MET:N	2.20	0.73
6:E:100:ARG:NH2	6:E:121:TYR:O	2.21	0.73
1:2:1534:C:H41	1:2:1599:U:H3	1.37	0.73
34:g:313:THR:O	34:g:315:GLY:N	2.21	0.73
1:2:1507:G:O2'	1:2:1511:U:O4	2.06	0.73
5:D:138:VAL:HG22	5:D:184:ILE:HG22	1.71	0.72
1:2:1276:A:O2'	12:K:57:TYR:OH	2.06	0.72
25:X:67:ARG:HG3	25:X:115:ILE:HG12	1.70	0.71
30:c:28:THR:O	30:c:45:ASN:ND2	2.22	0.71
34:g:66:VAL:HA	34:g:82:SER:HA	1.71	0.71
29:b:75:GLU:OE1	29:b:75:GLU:N	2.21	0.71
1:2:913:A:H2	9:H:99:ARG:H	1.39	0.71
19:R:127:ASN:ND2	19:R:127:ASN:O	2.23	0.71
1:2:145:G:O6	8:G:178:ARG:NH2	2.24	0.71
1:2:640:A:H2'	1:2:641:A:C8	2.26	0.71
1:2:1488:C:H3'	1:2:1489:A:H4'	1.73	0.71
12:K:65:ARG:HD3	31:d:25:SER:HB2	1.74	0.70
18:Q:46:THR:HG22	18:Q:47:LEU:HG	1.73	0.70
1:2:1618:C:H5''	1:2:1619:A:H2	1.56	0.70
1:2:952:G:H21	16:O:52:THR:HG21	1.57	0.70
1:2:1863:A:OP2	28:a:4:LYS:NZ	2.24	0.70
1:2:77:A:C8	8:G:154:ARG:HG3	2.27	0.70
1:2:1538:C:H5''	1:2:1593:C:H42	1.57	0.70
5:D:124:ARG:O	5:D:128:GLU:HG2	1.92	0.70
21:T:131:LEU:HA	21:T:134:ILE:HD12	1.74	0.70
1:2:686:U:O2	9:H:118:ARG:NH2	2.25	0.69

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:H:38:ALA:HA	9:H:41:ARG:HH21	1.57	0.69
12:K:46:MET:HA	12:K:49:MET:HE3	1.73	0.69
1:2:551:U:O2'	32:e:44:ASN:OD1	2.11	0.69
1:2:499:G:H2'	1:2:501:C:C5	2.28	0.69
10:I:3:ILE:HB	10:I:30:GLY:HA3	1.73	0.69
1:2:1588:A:N1	1:2:1626:C:O2'	2.25	0.69
10:I:106:SER:HB3	10:I:171:LEU:HG	1.75	0.69
21:T:64:LEU:HD11	21:T:70:ALA:H	1.58	0.69
7:F:176:GLU:O	7:F:180:ALA:HB2	1.93	0.68
5:D:143:ARG:NH1	36:n:156:ASP:OD2	2.26	0.68
3:B:63:LYS:HE2	3:B:90:ASP:HA	1.74	0.68
34:g:15:ASN:ND2	34:g:37:ASP:OD2	2.26	0.68
1:2:1024:A:OP2	15:N:124:ARG:NH2	2.26	0.68
10:I:6:ASP:OD2	10:I:8:TRP:NE1	2.26	0.68
1:2:1400:U:O4	1:2:1401:A:N6	2.25	0.68
17:P:38:SER:HB3	20:S:88:LYS:HD3	1.76	0.68
3:B:32:ASP:OD2	3:B:46:LYS:NZ	2.27	0.68
7:F:72:LEU:HD11	7:F:176:GLU:HB3	1.74	0.68
34:g:120:ILE:HB	34:g:132:TRP:HB2	1.75	0.68
30:c:67:ARG:HG3	30:c:68:LEU:HG	1.75	0.67
1:2:973:C:N3	16:O:55:ARG:NH1	2.42	0.67
1:2:643:A:OP1	11:J:39:ASN:ND2	2.27	0.67
5:D:106:ARG:HG3	5:D:175:VAL:HG22	1.75	0.67
10:I:41:ARG:HE	10:I:43:ILE:HD12	1.59	0.67
1:2:305:U:H5'	10:I:41:ARG:HH12	1.59	0.67
1:2:1541:G:H22	1:2:1591:C:N4	1.92	0.67
1:2:1858:G:OP2	16:O:146:ARG:NH2	2.28	0.67
1:2:1724:A:N6	1:2:1812:U:O4	2.27	0.67
1:2:1758:G:N2	1:2:1773:C:N3	2.42	0.67
21:T:98:SER:O	21:T:102:ARG:HG3	1.94	0.67
22:U:78:ASP:OD2	31:d:44:ARG:NH1	2.26	0.67
5:D:172:VAL:HG22	5:D:185:LYS:HG3	1.76	0.67
5:D:29:LEU:HD12	5:D:34:TYR:HB2	1.76	0.67
1:2:1431:G:H5''	1:2:1432:U:H5'	1.76	0.67
21:T:20:ALA:HA	21:T:23:LYS:HE3	1.76	0.67
1:2:560:A:OP2	11:J:177:ASN:ND2	2.28	0.66
1:2:1453:C:H5	1:2:1476:A:H2	1.44	0.66
7:F:59:LYS:HD2	7:F:60:ARG:N	2.10	0.66
6:E:54:TYR:O	26:Y:15:ASN:ND2	2.28	0.66
3:B:125:VAL:HG22	3:B:172:MET:HE3	1.76	0.66
6:E:100:ARG:NH1	6:E:118:GLU:OE2	2.27	0.66

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:F:204:ARG:O	16:O:117:ARG:NH2	2.28	0.66
21:T:103:VAL:O	21:T:107:LEU:HB2	1.95	0.66
7:F:128:ILE:HD13	7:F:135:ARG:HG3	1.76	0.66
1:2:77:A:H2	8:G:175:LYS:HG3	1.60	0.66
1:2:567:C:N4	1:2:578:C:OP1	2.28	0.66
12:K:11:ILE:HG21	12:K:48:ALA:HB1	1.78	0.66
30:c:12:ALA:HB1	30:c:32:VAL:HB	1.76	0.66
1:2:332:G:O6	8:G:186:GLN:NE2	2.29	0.66
1:2:629:A:O2'	1:2:631:U:OP1	2.14	0.66
9:H:29:GLU:O	9:H:30:LEU:C	2.38	0.66
34:g:89:LEU:HD12	34:g:99:ARG:HB2	1.77	0.66
1:2:1534:C:H5''	1:2:1535:U:H2'	1.78	0.65
1:2:1587:G:H4'	21:T:92:PHE:HB2	1.78	0.65
1:2:1587:G:H21	21:T:94:ARG:HD2	1.61	0.65
7:F:74:ASN:O	7:F:74:ASN:ND2	2.22	0.65
30:c:46:VAL:HG23	30:c:56:LEU:HD11	1.77	0.65
1:2:1754:G:N2	1:2:1781:A:N3	2.41	0.65
17:P:63:ALA:HB1	17:P:73:PRO:HG3	1.78	0.65
1:2:1129:G:OP1	29:b:22:LYS:NZ	2.30	0.65
1:2:1322:G:H5''	1:2:1324:G:N7	2.12	0.65
7:F:141:VAL:HG13	30:c:46:VAL:HG22	1.77	0.65
1:2:499:G:H2'	1:2:501:C:H5	1.62	0.65
1:2:928:G:H2'	1:2:929:G:C8	2.32	0.65
1:2:373:G:H4'	13:L:85:THR:HG21	1.78	0.65
1:2:1588:A:N6	1:2:1627:C:OP2	2.29	0.65
5:D:16:ILE:HD11	31:d:36:LEU:HD23	1.79	0.65
9:H:40:LEU:HA	9:H:43:LEU:HB2	1.79	0.65
34:g:37:ASP:O	34:g:39:THR:OG1	2.15	0.65
22:U:80:PHE:HB3	31:d:52:PHE:HB3	1.79	0.65
7:F:92:ILE:HD11	7:F:169:ILE:HD11	1.80	0.64
33:f:136:PHE:HD1	33:f:137:ASP:H	1.45	0.64
1:2:1243:U:O2	1:2:1264:C:N3	2.30	0.64
27:Z:71:ALA:O	27:Z:75:GLU:HG3	1.97	0.64
10:I:119:LEU:HG	10:I:153:LYS:HG2	1.79	0.64
13:L:82:MET:HE3	13:L:85:THR:HG23	1.78	0.64
1:2:1445:U:O4	1:2:1446:A:N6	2.31	0.64
9:H:29:GLU:O	9:H:32:MET:HG2	1.98	0.64
18:Q:37:ARG:NH1	21:T:10:ASN:OD1	2.31	0.64
20:S:82:TRP:HH2	21:T:38:LYS:HE2	1.62	0.64
1:2:617:G:N7	25:X:67:ARG:NH1	2.43	0.64
1:2:64:A:H2	1:2:83:A:H62	1.44	0.64

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:G:150:GLU:OE1	8:G:150:GLU:N	2.22	0.64
10:I:150:ASP:OD1	10:I:153:LYS:NZ	2.31	0.64
11:J:131:ARG:NH1	11:J:143:ASN:OD1	2.31	0.64
20:S:26:ILE:HG21	20:S:59:LEU:H	1.62	0.64
7:F:59:LYS:HD2	7:F:60:ARG:H	1.63	0.63
7:F:116:ILE:HG22	7:F:182:LYS:HA	1.79	0.63
1:2:1568:C:H2'	21:T:94:ARG:CZ	2.28	0.63
2:A:108:PHE:HB2	2:A:136:GLU:HG2	1.81	0.63
17:P:41:GLN:HA	17:P:44:ARG:HB3	1.81	0.63
1:2:919:A:OP2	15:N:64:ARG:NH2	2.30	0.63
1:2:989:C:OP2	3:B:155:TYR:OH	2.12	0.63
5:D:38:GLU:HG3	5:D:49:ILE:HB	1.80	0.63
5:D:163:PRO:O	5:D:167:TYR:HB2	1.99	0.63
7:F:156:THR:O	7:F:160:GLU:HG3	1.98	0.63
21:T:77:LYS:HG3	21:T:78:ILE:HG23	1.81	0.63
1:2:541:U:H3	1:2:543:C:H41	1.46	0.63
1:2:617:G:H4'	25:X:88:ASP:HB3	1.79	0.63
20:S:23:ARG:HA	27:Z:43:LYS:HE2	1.80	0.63
1:2:77:A:H8	8:G:154:ARG:HG3	1.63	0.63
1:2:1428:G:H4'	1:2:1430:C:H41	1.64	0.63
1:2:1745:A:H1'	8:G:31:ARG:HH22	1.64	0.63
8:G:159:ARG:NH2	8:G:171:THR:O	2.32	0.63
19:R:36:GLU:HG3	19:R:47:ARG:HD2	1.79	0.63
20:S:141:ARG:O	20:S:145:THR:OG1	2.17	0.63
1:2:587:A:O2'	1:2:588:G:N2	2.31	0.62
12:K:15:LEU:HG	12:K:52:LEU:HD12	1.80	0.62
18:Q:13:PHE:CB	18:Q:23:ALA:H	2.12	0.62
1:2:1569:A:H61	1:2:1613:G:N2	1.97	0.62
20:S:46:ARG:HE	21:T:35:ASP:HB2	1.64	0.62
1:2:94:G:HO2'	1:2:508:A:HO2'	1.45	0.62
1:2:1661:A:C8	31:d:14:PHE:HB3	2.35	0.62
1:2:1749:G:OP2	1:2:1750:C:N4	2.27	0.62
13:L:77:VAL:HG11	13:L:80:MET:HE3	1.81	0.62
18:Q:105:LYS:HA	18:Q:108:ILE:HD11	1.80	0.62
5:D:8:LYS:HB3	22:U:61:LEU:HD21	1.79	0.62
11:J:18:ARG:O	11:J:24:ARG:NH1	2.33	0.62
1:2:296:U:O2'	6:E:131:VAL:O	2.12	0.62
1:2:537:C:H5'	1:2:539:C:H5''	1.82	0.61
10:I:11:ARG:O	13:L:136:LYS:NZ	2.28	0.61
7:F:110:GLN:O	7:F:114:ASN:HB2	2.00	0.61
21:T:101:ARG:HH12	21:T:115:LYS:HE3	1.65	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:64:A:OP1	8:G:136:LYS:NZ	2.32	0.61
7:F:116:ILE:H	7:F:183:GLY:N	1.97	0.61
12:K:11:ILE:HG23	12:K:52:LEU:HD11	1.82	0.61
1:2:455:A:H2'	1:2:456:C:C6	2.36	0.61
1:2:1230:C:O2'	1:2:1231:C:OP2	2.17	0.61
1:2:1772:C:H5'	1:2:1773:C:H5	1.64	0.61
14:M:93:LYS:HG3	14:M:94:ILE:H	1.65	0.61
34:g:61:GLY:O	34:g:100:ARG:NH1	2.30	0.60
1:2:1587:G:C4	21:T:73:GLY:HA3	2.35	0.60
34:g:42:MET:HE3	34:g:59:LEU:HD21	1.83	0.60
34:g:133:ASN:HD21	34:g:137:VAL:H	1.49	0.60
1:2:1153:C:OP2	24:W:71:LYS:NZ	2.34	0.60
7:F:144:LEU:HD23	30:c:49:PRO:HB3	1.83	0.60
1:2:92:A:H1'	6:E:3:ARG:HB3	1.82	0.60
1:2:959:G:OP2	16:O:38:ASN:ND2	2.34	0.60
1:2:1474:A:H5''	18:Q:121:VAL:HG12	1.83	0.60
3:B:33:VAL:HG12	3:B:96:CYS:HB2	1.83	0.60
7:F:72:LEU:HD21	7:F:176:GLU:HG3	1.84	0.60
10:I:110:ARG:O	10:I:114:GLU:HG2	2.02	0.60
26:Y:55:ILE:HG12	26:Y:75:ILE:HG12	1.82	0.60
34:g:296:GLN:O	34:g:311:GLN:NE2	2.35	0.60
1:2:107:A:H2'	1:2:108:G:C8	2.36	0.60
1:2:84:A:N3	1:2:150:A:O2'	2.34	0.60
1:2:1241:A:O2'	1:2:1266:C:N4	2.31	0.60
8:G:221:ARG:HH11	8:G:225:GLN:HG3	1.66	0.60
1:2:125:C:OP1	1:2:127:C:N4	2.34	0.60
1:2:1174:U:H2'	1:2:1175:G:C8	2.36	0.60
1:2:1679:A:H2'	7:F:60:ARG:HD3	1.83	0.60
9:H:46:THR:HG23	9:H:65:PRO:HD3	1.83	0.60
10:I:101:ILE:HD12	10:I:190:LEU:HD11	1.83	0.60
25:X:68:LYS:HB3	25:X:91:LEU:HD22	1.84	0.60
15:N:99:ARG:NH2	15:N:119:GLU:OE2	2.31	0.60
1:2:94:G:O2'	1:2:508:A:O2'	2.18	0.60
1:2:1261:C:O2	1:2:1513:C:O2'	2.19	0.60
1:2:1587:G:N3	21:T:94:ARG:HG2	2.16	0.60
1:2:1601:A:N6	1:2:1636:G:OP2	2.35	0.60
20:S:124:ARG:HB2	20:S:131:VAL:HG23	1.84	0.60
1:2:16:G:H2'	1:2:17:C:C6	2.37	0.60
7:F:52:SER:H	7:F:86:LYS:HZ3	1.50	0.60
13:L:103:GLU:OE2	13:L:105:ARG:NH1	2.30	0.60
1:2:1457:U:H2'	1:2:1458:G:H8	1.67	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1522:A:N3	17:P:128:HIS:ND1	2.50	0.59
3:B:166:LYS:O	3:B:170:GLU:HG3	2.01	0.59
1:2:904:A:C8	1:2:905:C:H5	2.20	0.59
1:2:1203:G:H2'	1:2:1204:A:C8	2.37	0.59
30:c:17:VAL:HA	30:c:30:VAL:HG12	1.84	0.59
1:2:1446:A:H5''	22:U:58:THR:HG22	1.84	0.59
1:2:1854:U:H2'	1:2:1855:G:H8	1.68	0.59
6:E:11:ARG:HA	6:E:28:ALA:HB2	1.85	0.59
7:F:117:ILE:HG23	7:F:183:GLY:N	2.18	0.59
9:H:109:ARG:HG3	9:H:110:THR:H	1.67	0.59
20:S:87:GLN:O	20:S:88:LYS:HG3	2.02	0.59
30:c:67:ARG:HE	30:c:68:LEU:H	1.49	0.59
1:2:17:C:O2'	1:2:1194:A:N1	2.32	0.59
22:U:42:GLY:HA3	22:U:101:ILE:HD11	1.84	0.59
8:G:21:GLU:O	8:G:25:ARG:HD3	2.02	0.59
12:K:71:LEU:HD23	12:K:93:THR:HG21	1.84	0.59
34:g:67:SER:OG	34:g:110:SER:OG	2.17	0.59
6:E:192:ILE:HB	6:E:243:GLY:HA3	1.84	0.59
7:F:77:MET:HB2	7:F:84:GLY:H	1.68	0.59
1:2:188:C:OP2	1:2:208:G:N2	2.35	0.59
1:2:1543:U:OP2	21:T:79:TYR:OH	2.20	0.59
1:2:1602:U:H1'	20:S:24:ARG:HD2	1.85	0.59
1:2:1739:C:O3'	1:2:1740:C:H4'	2.02	0.59
1:2:855:G:O2'	13:L:71:ARG:NH1	2.35	0.59
1:2:952:G:H2'	1:2:953:C:C6	2.38	0.59
29:b:19:HIS:HB3	29:b:22:LYS:HG3	1.84	0.59
34:g:51:ASN:HD21	34:g:54:ILE:HG13	1.67	0.59
1:2:455:A:H2'	1:2:456:C:H6	1.67	0.58
1:2:1217:A:H2'	1:2:1218:C:C6	2.38	0.58
7:F:145:ARG:HG3	30:c:48:GLY:HA3	1.85	0.58
9:H:9:VAL:HG13	9:H:11:PRO:HD3	1.84	0.58
9:H:40:LEU:HD12	9:H:40:LEU:H	1.67	0.58
19:R:6:THR:HG22	19:R:8:THR:H	1.67	0.58
1:2:534:G:H2'	1:2:535:G:C8	2.37	0.58
17:P:79:HIS:CE1	17:P:98:ASN:HA	2.38	0.58
1:2:963:A:H1'	1:2:964:A:C8	2.38	0.58
1:2:1290:G:N7	31:d:2:GLY:N	2.51	0.58
3:B:103:MET:HE1	3:B:212:VAL:O	2.04	0.58
9:H:29:GLU:HG2	9:H:33:ASN:ND2	2.17	0.58
22:U:48:LEU:HD11	22:U:93:SER:HB3	1.85	0.58
1:2:1613:G:N7	17:P:42:ARG:NH2	2.50	0.58

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1636:G:C8	7:F:164:ARG:HD2	2.39	0.58
7:F:78:MET:HG3	7:F:159:ARG:HG3	1.85	0.58
20:S:130:ARG:NE	20:S:134:GLN:HB2	2.10	0.58
25:X:41:PHE:O	25:X:76:LYS:NZ	2.35	0.58
30:c:9:ILE:HA	30:c:59:LEU:HA	1.85	0.58
1:2:664:A:O2'	1:2:670:A:N1	2.32	0.58
10:I:150:ASP:HA	10:I:153:LYS:HD3	1.85	0.58
22:U:17:ILE:HG23	22:U:19:ARG:H	1.68	0.58
8:G:22:ARG:H	8:G:22:ARG:HE	1.52	0.58
14:M:44:LYS:HB3	14:M:46:GLN:HG3	1.85	0.58
22:U:68:THR:O	31:d:40:ARG:NH1	2.37	0.58
1:2:942:G:H2'	1:2:943:U:C6	2.39	0.58
16:O:46:ASP:OD1	16:O:47:LEU:N	2.37	0.58
21:T:123:LEU:HD21	21:T:131:LEU:HD21	1.86	0.58
1:2:1615:U:OP2	17:P:43:ARG:NH1	2.37	0.58
17:P:30:TYR:OH	17:P:85:ILE:O	2.22	0.58
20:S:44:VAL:HG11	20:S:67:VAL:HG22	1.86	0.58
33:f:94:LYS:O	33:f:96:LYS:NZ	2.37	0.58
1:2:678:U:H2'	1:2:679:A:H8	1.69	0.57
1:2:957:A:OP1	16:O:57:THR:OG1	2.14	0.57
4:C:263:LYS:HZ3	4:C:268:GLU:HA	1.69	0.57
18:Q:109:LYS:O	18:Q:113:ILE:HG12	2.04	0.57
24:W:57:ARG:NH1	29:b:26:GLN:OE1	2.37	0.57
27:Z:99:LEU:HA	27:Z:109:TYR:HD1	1.67	0.57
30:c:33:GLU:HG2	30:c:41:SER:HA	1.86	0.57
7:F:204:ARG:NH1	30:c:60:GLU:OE1	2.34	0.57
17:P:92:SER:O	17:P:107:ILE:HG12	2.04	0.57
27:Z:73:VAL:HG21	27:Z:84:ALA:HB2	1.85	0.57
1:2:1354:G:N2	1:2:1357:A:OP2	2.32	0.57
15:N:37:ILE:HD11	15:N:63:VAL:HG21	1.87	0.57
20:S:130:ARG:HD3	20:S:134:GLN:O	2.04	0.57
7:F:115:ALA:H	7:F:182:LYS:C	2.12	0.57
18:Q:42:ILE:HG23	18:Q:44:PRO:HD2	1.84	0.57
18:Q:118:THR:HA	18:Q:121:VAL:HG23	1.85	0.57
6:E:165:GLU:HG3	6:E:166:THR:HG23	1.87	0.57
9:H:36:LEU:HB3	9:H:39:GLN:HB2	1.85	0.57
34:g:87:LEU:HD22	34:g:111:VAL:HG21	1.87	0.57
34:g:302:TYR:HE2	34:g:308:ARG:HE	1.53	0.57
1:2:63:U:O2'	1:2:170:A:N3	2.37	0.57
34:g:106:LYS:HB3	34:g:125:ARG:HB2	1.87	0.57
2:A:8:LEU:HD11	23:V:39:VAL:HG11	1.87	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:c:28:THR:CB	30:c:46:VAL:O	2.47	0.57
1:2:15:U:H2'	1:2:16:G:O4'	2.04	0.57
4:C:72:ASP:OD2	4:C:272:HIS:NE2	2.37	0.57
6:E:175:PHE:HE2	6:E:198:ARG:HD2	1.70	0.57
18:Q:113:ILE:HD12	18:Q:119:LEU:HD13	1.87	0.57
1:2:17:C:H2'	1:2:18:C:C6	2.40	0.57
1:2:106:C:OP1	1:2:431:G:O2'	2.21	0.57
1:2:678:U:OP1	15:N:127:ARG:NH2	2.38	0.57
34:g:121:VAL:HB	34:g:154:VAL:HG21	1.87	0.57
1:2:104:A:N6	1:2:356:C:O2	2.38	0.57
1:2:153:G:H21	8:G:13:GLN:NE2	2.03	0.57
7:F:80:GLY:N	7:F:83:ASN:OD1	2.38	0.57
7:F:116:ILE:H	7:F:182:LYS:C	2.13	0.57
7:F:143:PRO:HA	7:F:146:ARG:HG3	1.85	0.57
22:U:28:ASN:OD1	22:U:31:SER:OG	2.14	0.57
1:2:51:U:H2'	1:2:52:G:H8	1.70	0.56
1:2:1598:G:H5'	27:Z:80:ARG:HD2	1.86	0.56
1:2:1642:U:O2'	1:2:1643:U:O5'	2.21	0.56
19:R:77:GLU:HG2	19:R:80:ARG:HH21	1.68	0.56
1:2:118:C:H1'	1:2:445:A:C5	2.40	0.56
1:2:1263:U:O2	31:d:28:HIS:NE2	2.38	0.56
1:2:1401:A:H61	1:2:1446:A:H61	1.54	0.56
1:2:1453:C:C5	1:2:1476:A:H2	2.22	0.56
1:2:1571:G:O2'	1:2:1572:C:OP1	2.22	0.56
1:2:1605:G:H1	1:2:1635:C:H1'	1.70	0.56
11:J:28:GLU:HB2	11:J:40:LYS:HE3	1.87	0.56
20:S:124:ARG:NE	20:S:130:ARG:O	2.38	0.56
1:2:301:A:H4'	1:2:302:A:H5''	1.88	0.56
1:2:1253:A:N6	1:2:1665:G:O2'	2.38	0.56
1:2:1275:G:OP1	12:K:1:MET:N	2.38	0.56
1:2:1845:A:H2'	1:2:1846:G:C8	2.41	0.56
5:D:40:ARG:HE	5:D:49:ILE:HD11	1.69	0.56
9:H:31:GLU:O	9:H:37:LYS:HG3	2.04	0.56
18:Q:111:ILE:O	18:Q:115:TYR:HB2	2.05	0.56
1:2:301:A:O2'	1:2:302:A:OP2	2.17	0.56
1:2:1231:C:H4'	1:2:1665:G:C8	2.39	0.56
5:D:48:ILE:HB	5:D:86:LEU:HB3	1.87	0.56
7:F:78:MET:O	7:F:79:HIS:HB2	2.04	0.56
12:K:1:MET:HA	12:K:1:MET:HE2	1.86	0.56
17:P:87:PRO:HG2	17:P:112:ILE:HD11	1.86	0.56
34:g:21:ILE:HD12	34:g:21:ILE:O	2.06	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1098:C:H2'	1:2:1099:G:C8	2.40	0.56
1:2:1144:A:H5'	1:2:1355:C:H41	1.71	0.56
1:2:1273:C:O2'	1:2:1274:G:O5'	2.24	0.56
10:I:143:LYS:O	10:I:147:LYS:HG2	2.06	0.56
18:Q:108:ILE:HD12	18:Q:109:LYS:H	1.71	0.56
19:R:71:ILE:HD12	19:R:74:GLN:H	1.71	0.56
11:J:53:ILE:HD13	11:J:81:LEU:HD21	1.88	0.56
20:S:79:ILE:HG23	21:T:33:TRP:CH2	2.41	0.56
21:T:51:ASN:HB2	21:T:54:TYR:HB2	1.87	0.56
23:V:39:VAL:HG13	23:V:44:GLY:HA2	1.87	0.56
1:2:29:G:H2'	1:2:30:C:C6	2.41	0.56
1:2:1611:G:H1	1:2:1627:C:N4	2.03	0.56
7:F:201:LYS:O	7:F:204:ARG:HG2	2.06	0.56
18:Q:47:LEU:HB2	18:Q:81:ILE:HD13	1.88	0.56
34:g:44:LYS:HB2	34:g:54:ILE:HD13	1.88	0.56
1:2:1596:U:O4	27:Z:104:ARG:NE	2.32	0.56
9:H:91:HIS:HD2	9:H:172:THR:HG21	1.69	0.56
12:K:55:ARG:NE	12:K:55:ARG:O	2.39	0.56
21:T:124:THR:O	21:T:128:GLN:N	2.35	0.56
34:g:5:MET:HB2	34:g:312:VAL:HG12	1.88	0.56
1:2:1856:C:H2'	1:2:1857:G:C8	2.40	0.56
6:E:151:ASP:HB3	6:E:154:ILE:HG13	1.87	0.56
17:P:87:PRO:HG2	17:P:112:ILE:CD1	2.35	0.56
7:F:165:ASN:O	7:F:168:THR:OG1	2.19	0.55
13:L:128:VAL:HG12	13:L:142:VAL:HA	1.87	0.55
18:Q:13:PHE:HB3	18:Q:23:ALA:H	1.71	0.55
34:g:85:GLY:HA2	34:g:108:VAL:HB	1.88	0.55
1:2:165:G:H4'	8:G:53:SER:HB2	1.88	0.55
1:2:1174:U:H2'	1:2:1175:G:H8	1.71	0.55
8:G:45:TRP:HA	8:G:48:TYR:HD2	1.71	0.55
9:H:63:PHE:HB3	9:H:97:GLN:OE1	2.06	0.55
20:S:103:LEU:HG	20:S:106:LYS:HB2	1.88	0.55
1:2:385:G:H3'	13:L:136:LYS:HB2	1.87	0.55
1:2:1033:G:N1	1:2:1080:A:O2'	2.32	0.55
1:2:1590:C:H4'	21:T:82:ARG:HG3	1.87	0.55
1:2:1644:C:H4'	18:Q:140:ARG:HB2	1.88	0.55
6:E:55:ALA:HB1	6:E:60:GLU:HB2	1.88	0.55
8:G:10:THR:HG22	8:G:128:THR:HG23	1.89	0.55
1:2:549:C:H2'	1:2:550:C:H6	1.72	0.55
1:2:1611:G:H5'	20:S:88:LYS:HE2	1.88	0.55
10:I:42:ARG:HB3	10:I:59:ARG:HB2	1.88	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:805:U:O2	1:2:860:G:N2	2.34	0.55
7:F:116:ILE:HG13	7:F:119:SER:HB2	1.88	0.55
8:G:168:LYS:HD3	8:G:168:LYS:H	1.71	0.55
1:2:1091:C:HO2'	24:W:2:VAL:N	2.04	0.55
3:B:76:ASN:OD1	3:B:76:ASN:N	2.38	0.55
13:L:111:VAL:HG12	13:L:140:PHE:HB2	1.88	0.55
25:X:123:VAL:HG12	25:X:124:LYS:HG3	1.88	0.55
1:2:1224:G:H22	1:2:1643:U:H5	1.54	0.55
16:O:40:THR:HG21	16:O:74:ALA:HB2	1.89	0.55
20:S:134:GLN:N	20:S:134:GLN:OE1	2.38	0.55
1:2:185:G:H3'	1:2:186:C:H5''	1.87	0.55
1:2:1139:C:H2'	1:2:1140:G:O4'	2.07	0.55
1:2:1589:A:O2'	21:T:79:TYR:O	2.16	0.55
16:O:96:LYS:HG2	16:O:132:VAL:HG11	1.89	0.55
16:O:131:ASP:OD1	16:O:133:THR:HG22	2.07	0.55
1:2:489:A:N1	1:2:511:U:H5	2.05	0.55
33:f:81:LYS:HB3	33:f:83:LYS:HG3	1.89	0.55
1:2:51:U:H2'	1:2:52:G:C8	2.42	0.54
1:2:549:C:H2'	1:2:550:C:C6	2.41	0.54
30:c:20:ARG:HA	30:c:27:CYS:O	2.07	0.54
1:2:92:A:O2'	6:E:3:ARG:O	2.19	0.54
10:I:162:LEU:HD11	10:I:191:GLU:HG2	1.89	0.54
17:P:79:HIS:HE1	17:P:98:ASN:HA	1.71	0.54
5:D:210:ILE:HD11	19:R:16:ILE:HG13	1.88	0.54
21:T:52:TRP:O	21:T:55:THR:OG1	2.13	0.54
1:2:1643:U:H2'	1:2:1644:C:C6	2.42	0.54
2:A:135:THR:O	2:A:138:SER:OG	2.25	0.54
10:I:177:SER:OG	10:I:182:CYS:SG	2.66	0.54
11:J:84:ILE:HG13	11:J:86:VAL:HG13	1.90	0.54
24:W:111:MET:HE3	24:W:116:ALA:HA	1.89	0.54
1:2:1296:U:O2'	1:2:1298:G:OP1	2.26	0.54
1:2:1374:C:H2'	1:2:1375:G:O4'	2.07	0.54
7:F:82:ASN:HD21	7:F:88:MET:HG3	1.72	0.54
21:T:101:ARG:HH22	21:T:121:ARG:NH2	2.06	0.54
25:X:84:PHE:CE2	25:X:86:PRO:HA	2.42	0.54
34:g:217:MET:HE1	34:g:219:TRP:CE2	2.43	0.54
1:2:1096:G:H1	1:2:1136:U:H5	1.55	0.54
1:2:1825:A:OP2	1:2:1826:G:N2	2.40	0.54
7:F:122:ARG:N	7:F:197:GLU:OE1	2.41	0.54
13:L:80:MET:HE1	13:L:120:VAL:HB	1.89	0.54
34:g:33:SER:HB3	34:g:43:TRP:HE1	1.73	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1738:C:N4	1:2:1790:A:OP1	2.41	0.54
2:A:76:VAL:HG12	2:A:123:VAL:HB	1.90	0.54
34:g:64:HIS:HB3	34:g:82:SER:OG	2.08	0.54
36:n:163:THR:HG22	36:n:164:LYS:H	1.73	0.54
1:2:3:C:O2	11:J:18:ARG:NH2	2.41	0.54
1:2:808:A:O2'	1:2:809:A:OP1	2.26	0.54
1:2:1590:C:H3'	1:2:1591:C:C6	2.43	0.54
22:U:25:THR:HG22	22:U:86:LYS:HB2	1.89	0.54
33:f:136:PHE:HD1	33:f:137:ASP:N	2.04	0.54
26:Y:18:LEU:O	26:Y:85:ASN:ND2	2.39	0.54
1:2:5:U:H2'	1:2:6:G:H8	1.73	0.53
1:2:1224:G:H1	1:2:1643:U:H5	1.56	0.53
1:2:1457:U:H2'	1:2:1458:G:C8	2.42	0.53
2:A:173:LEU:HD11	2:A:177:MET:HE3	1.89	0.53
1:2:1013:U:OP1	1:2:1129:G:O2'	2.25	0.53
1:2:1562:C:H2'	1:2:1563:G:H8	1.73	0.53
1:2:1566:G:O2'	1:2:1567:G:OP1	2.23	0.53
1:2:1863:A:H1'	28:a:79:ILE:HD13	1.90	0.53
12:K:35:LEU:HD21	12:K:49:MET:HE1	1.90	0.53
26:Y:7:ILE:HD13	26:Y:43:LYS:HD2	1.89	0.53
1:2:1402:A:OP1	22:U:52:GLY:N	2.32	0.53
1:2:1528:G:O2'	1:2:1666:C:OP1	2.23	0.53
6:E:174:LYS:O	6:E:179:ASN:ND2	2.41	0.53
7:F:71:ARG:HH22	7:F:148:ASN:HA	1.72	0.53
20:S:80:PRO:HD2	21:T:33:TRP:CZ2	2.43	0.53
34:g:7:LEU:HD11	34:g:308:ARG:HB3	1.89	0.53
1:2:579:C:OP2	1:2:579:C:H6	1.91	0.53
17:P:90:VAL:CG2	17:P:109:PRO:HG3	2.38	0.53
1:2:678:U:H2'	1:2:679:A:C8	2.43	0.53
1:2:1544:C:N4	21:T:80:GLY:O	2.22	0.53
1:2:1607:A:N7	1:2:1632:G:N3	2.57	0.53
7:F:118:ASN:HB3	7:F:185:SER:HB3	1.89	0.53
12:K:26:ASP:H	12:K:46:MET:HE1	1.74	0.53
19:R:24:LEU:HB2	19:R:58:MET:SD	2.49	0.53
19:R:87:GLU:H	19:R:87:GLU:CD	2.16	0.53
20:S:86:ARG:HG3	20:S:108:ARG:HH21	1.73	0.53
16:O:113:GLN:OE1	28:a:46:GLU:N	2.35	0.53
1:2:678:U:O2	1:2:1027:A:N7	2.41	0.53
1:2:1264:C:O2'	1:2:1517:G:O6	2.16	0.53
4:C:263:LYS:NZ	4:C:268:GLU:HA	2.24	0.53
8:G:21:GLU:O	8:G:24:LEU:N	2.37	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1201:U:H2'	1:2:1202:U:C6	2.42	0.53
1:2:1210:G:OP1	28:a:82:LYS:NZ	2.33	0.53
1:2:1270:G:H22	1:2:1515:G:N2	2.07	0.53
1:2:1569:A:H61	1:2:1613:G:H21	1.57	0.53
17:P:89:MET:HE3	17:P:116:LEU:HD11	1.91	0.53
18:Q:26:LYS:HG2	18:Q:27:ARG:H	1.73	0.53
21:T:8:ASP:OD1	21:T:9:VAL:N	2.41	0.53
1:2:562:U:H2'	1:2:563:G:C8	2.43	0.53
1:2:903:A:H2'	1:2:904:A:C8	2.44	0.53
7:F:78:MET:CG	7:F:159:ARG:HG3	2.39	0.53
1:2:594:A:H61	1:2:643:A:H5''	1.74	0.53
1:2:1564:C:OP1	21:T:101:ARG:HD2	2.08	0.53
3:B:134:LEU:HD23	3:B:219:LYS:HG3	1.90	0.53
1:2:25:A:H2'	1:2:26:U:C6	2.44	0.52
1:2:1249:C:H2'	1:2:1250:A:C2	2.44	0.52
5:D:41:VAL:HG22	5:D:46:THR:HG22	1.90	0.52
9:H:109:ARG:HB3	9:H:111:LYS:HE3	1.90	0.52
20:S:131:VAL:O	20:S:132:ARG:NH1	2.42	0.52
1:2:962:A:O2'	16:O:65:ASP:OD2	2.20	0.52
4:C:255:LEU:HD13	23:V:23:ILE:HD11	1.91	0.52
7:F:100:ILE:HA	7:F:103:LEU:HB3	1.92	0.52
17:P:32:GLN:HB3	17:P:35:GLN:HB2	1.91	0.52
20:S:24:ARG:HH21	20:S:25:LYS:HB3	1.74	0.52
21:T:53:PHE:HA	21:T:56:ARG:HE	1.74	0.52
29:b:67:THR:HG22	29:b:68:GLY:H	1.75	0.52
1:2:508:A:H3'	1:2:509:G:H8	1.74	0.52
1:2:996:A:H2'	1:2:997:A:C8	2.44	0.52
1:2:1759:G:N2	1:2:1773:C:OP1	2.37	0.52
6:E:44:LEU:HD13	6:E:72:ILE:HD11	1.91	0.52
16:O:45:THR:HA	16:O:52:THR:HA	1.90	0.52
1:2:1788:A:H62	1:2:1789:G:N2	2.01	0.52
12:K:15:LEU:HD21	12:K:49:MET:HB2	1.91	0.52
16:O:99:ALA:H	16:O:133:THR:HB	1.75	0.52
28:a:60:ASP:OD1	28:a:60:ASP:N	2.42	0.52
19:R:98:VAL:HG13	19:R:102:THR:HG23	1.90	0.52
30:c:58:LEU:C	30:c:60:GLU:H	2.17	0.52
34:g:78:ALA:O	34:g:89:LEU:HA	2.10	0.52
1:2:1025:U:OP1	1:2:1090:C:O2'	2.27	0.52
1:2:1653:U:H2'	1:2:1654:G:C8	2.45	0.52
6:E:60:GLU:HB3	26:Y:18:LEU:HD11	1.92	0.52
1:2:1088:U:H4'	1:2:1089:G:OP2	2.09	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:I:3:ILE:O	10:I:30:GLY:N	2.42	0.52
17:P:24:GLN:H	17:P:28:MET:HB2	1.75	0.52
18:Q:108:ILE:HD12	18:Q:109:LYS:N	2.23	0.52
30:c:36:ASP:N	30:c:36:ASP:OD1	2.43	0.52
3:B:124:HIS:HA	3:B:137:LEU:O	2.09	0.52
7:F:119:SER:OG	7:F:185:SER:HA	2.10	0.52
12:K:90:VAL:HG23	12:K:94:LEU:HD12	1.91	0.52
21:T:33:TRP:H	21:T:102:ARG:NH2	2.07	0.52
1:2:1587:G:H1	21:T:97:LYS:CE	2.20	0.52
5:D:25:LEU:HA	5:D:28:GLU:HB3	1.92	0.52
5:D:123:LEU:HD11	5:D:154:ASP:HB3	1.91	0.52
6:E:20:LEU:HD21	6:E:46:ILE:HD12	1.91	0.52
18:Q:86:GLN:O	18:Q:90:LYS:HG2	2.10	0.52
19:R:10:LYS:HD3	19:R:53:TYR:CZ	2.45	0.52
1:2:221:A:H4'	1:2:222:U:C5	2.45	0.52
1:2:301:A:H4'	1:2:302:A:C5'	2.40	0.52
1:2:359:U:OP2	25:X:18:ARG:HD3	2.10	0.52
1:2:945:U:H2'	1:2:946:U:C6	2.46	0.52
1:2:1231:C:H4'	1:2:1665:G:N7	2.24	0.52
1:2:1274:G:H8	1:2:1275:G:C5	2.28	0.52
1:2:1587:G:C5	21:T:73:GLY:HA3	2.45	0.52
36:n:154:PHE:O	36:n:158:GLU:HG3	2.11	0.52
9:H:42:GLU:O	9:H:44:ASN:N	2.44	0.51
21:T:35:ASP:HA	21:T:52:TRP:CH2	2.45	0.51
22:U:61:LEU:HB2	22:U:82:MET:HB3	1.92	0.51
1:2:96:C:H2'	1:2:97:U:C6	2.45	0.51
1:2:536:A:H2'	1:2:548:C:H42	1.75	0.51
1:2:1669:G:OP1	22:U:79:ARG:NH1	2.44	0.51
7:F:173:LEU:O	7:F:177:LEU:HB2	2.10	0.51
14:M:64:LEU:HD13	33:f:107:LYS:HE3	1.93	0.51
19:R:80:ARG:HG3	19:R:80:ARG:HH11	1.75	0.51
1:2:1084:A:OP1	1:2:1858:G:O2'	2.27	0.51
3:B:134:LEU:HB3	3:B:219:LYS:HB2	1.93	0.51
3:B:144:LYS:HD3	3:B:208:HIS:CG	2.45	0.51
12:K:48:ALA:O	12:K:52:LEU:HG	2.09	0.51
4:C:165:VAL:HG21	4:C:217:ALA:HB1	1.93	0.51
26:Y:86:GLU:OE1	26:Y:90:ARG:NE	2.43	0.51
1:2:953:C:H2'	1:2:954:U:O4'	2.10	0.51
1:2:1596:U:O2'	1:2:1597:C:O4'	2.21	0.51
4:C:254:ASP:HB2	23:V:1:MET:HG2	1.92	0.51
6:E:228:ILE:HD12	6:E:236:ILE:HD11	1.91	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:U:56:MET:HG3	22:U:86:LYS:HG2	1.93	0.51
1:2:388:U:H2'	1:2:389:A:C8	2.45	0.51
1:2:1241:A:N3	1:2:1518:C:H5''	2.26	0.51
1:2:1569:A:N3	21:T:94:ARG:NH2	2.58	0.51
1:2:377:G:H5''	10:I:98:LYS:HB3	1.92	0.51
1:2:1448:A:H2'	1:2:1449:G:O4'	2.11	0.51
12:K:8:ARG:O	12:K:11:ILE:HG22	2.11	0.51
20:S:86:ARG:HG3	20:S:108:ARG:NH2	2.25	0.51
28:a:36:ILE:HD13	28:a:78:VAL:HG11	1.92	0.51
1:2:1274:G:H1'	1:2:1275:G:C8	2.46	0.51
1:2:1538:C:H4'	21:T:44:GLU:H	1.75	0.51
7:F:158:ALA:O	7:F:162:ALA:CB	2.55	0.51
10:I:57:ALA:HB2	10:I:183:GLY:HA2	1.93	0.51
20:S:112:GLU:HG2	20:S:113:ARG:N	2.25	0.51
1:2:174:C:H42	1:2:175:A:H62	1.59	0.51
19:R:34:VAL:O	19:R:38:ILE:HG13	2.11	0.51
1:2:107:A:H2'	1:2:108:G:H8	1.76	0.51
1:2:809:A:OP2	6:E:186:GLY:HA3	2.11	0.51
1:2:823:U:H5	11:J:143:ASN:HB3	1.76	0.51
1:2:1264:C:O2'	1:2:1265:A:OP2	2.29	0.51
14:M:18:LEU:HA	14:M:21:VAL:HG12	1.93	0.51
18:Q:111:ILE:O	18:Q:115:TYR:CB	2.59	0.51
1:2:637:U:H2'	1:2:638:C:H6	1.75	0.50
1:2:747:U:O2'	1:2:795:A:N1	2.39	0.50
1:2:1519:U:H5''	1:2:1520:G:O4'	2.11	0.50
7:F:51:HIS:CD2	18:Q:47:LEU:HD21	2.46	0.50
10:I:107:THR:HG22	10:I:111:GLN:HE21	1.75	0.50
12:K:2:LEU:O	12:K:47:LYS:NZ	2.44	0.50
26:Y:20:ARG:HE	26:Y:22:GLN:HE21	1.59	0.50
1:2:318:A:N6	8:G:186:GLN:OE1	2.45	0.50
1:2:628:A:O2'	1:2:629:A:OP1	2.24	0.50
1:2:1121:G:O2'	3:B:204:ILE:O	2.25	0.50
1:2:1495:G:H4'	31:d:26:ASN:HD21	1.75	0.50
1:2:1667:U:H2'	1:2:1668:U:C6	2.47	0.50
10:I:121:LEU:HD11	10:I:124:LYS:HB2	1.93	0.50
17:P:90:VAL:HG23	17:P:109:PRO:HG3	1.94	0.50
21:T:56:ARG:H	21:T:56:ARG:HD3	1.76	0.50
27:Z:92:LEU:HD23	27:Z:97:ILE:HG13	1.93	0.50
1:2:1021:U:H4'	1:2:1022:U:H5'	1.92	0.50
1:2:1126:G:OP1	2:A:41:ARG:NH1	2.44	0.50
1:2:1611:G:H4'	20:S:88:LYS:HG2	1.93	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1710:C:N3	1:2:1824:A:N6	2.58	0.50
4:C:203:GLY:H	4:C:221:ASP:HB3	1.76	0.50
5:D:64:ARG:O	5:D:68:GLU:HG3	2.12	0.50
1:2:495:U:H2'	1:2:496:C:O4'	2.11	0.50
1:2:1569:A:O5'	21:T:94:ARG:NH1	2.36	0.50
1:2:546:G:H5''	1:2:547:G:N7	2.27	0.50
1:2:1627:C:C5'	21:T:41:LYS:HD3	2.41	0.50
1:2:1661:A:N7	31:d:14:PHE:HB3	2.26	0.50
6:E:95:THR:HA	26:Y:16:ARG:HH21	1.77	0.50
7:F:115:ALA:N	7:F:182:LYS:O	2.35	0.50
7:F:187:SER:HB3	7:F:190:ILE:HB	1.94	0.50
25:X:29:LYS:HE2	25:X:35:ALA:HB2	1.94	0.50
1:2:546:G:H5''	1:2:547:G:C8	2.46	0.50
1:2:579:C:H2'	1:2:580:U:O5'	2.10	0.50
1:2:581:U:H4'	26:Y:66:GLY:HA2	1.92	0.50
1:2:1125:C:OP1	19:R:129:LYS:HE2	2.10	0.50
1:2:1324:G:O2'	1:2:1325:G:N2	2.41	0.50
1:2:1389:C:H5	1:2:1481:G:H1	1.59	0.50
1:2:1590:C:OP2	21:T:82:ARG:NH2	2.45	0.50
1:2:1598:G:N7	27:Z:82:SER:N	2.60	0.50
3:B:168:MET:O	3:B:172:MET:HG3	2.11	0.50
4:C:264:SER:O	4:C:268:GLU:N	2.39	0.50
5:D:76:ARG:HB2	12:K:22:VAL:HG21	1.94	0.50
14:M:15:ASN:HA	14:M:18:LEU:HB2	1.94	0.50
34:g:72:SER:OG	34:g:77:PHE:O	2.24	0.50
1:2:12:U:H2'	1:2:13:C:O2	2.12	0.50
1:2:501:C:O2	1:2:501:C:H2'	2.12	0.50
1:2:625:G:O6	25:X:63:ASN:HB2	2.10	0.50
1:2:1083:A:N7	1:2:1841:C:O2'	2.37	0.50
1:2:1618:C:H4'	31:d:10:HIS:CE1	2.47	0.50
7:F:146:ARG:HD2	7:F:147:VAL:N	2.26	0.50
12:K:21:MET:SD	12:K:53:LYS:NZ	2.69	0.50
12:K:24:LYS:HB3	12:K:26:ASP:OD1	2.12	0.50
34:g:32:LEU:HD22	34:g:40:ILE:HG22	1.93	0.50
1:2:672:A:N6	1:2:1027:A:OP1	2.40	0.50
1:2:916:A:C5	15:N:73:ARG:HD3	2.47	0.50
1:2:1016:U:H5''	15:N:14:SER:HB3	1.94	0.50
3:B:33:VAL:HG23	3:B:44:ILE:HG13	1.93	0.50
9:H:30:LEU:O	9:H:31:GLU:C	2.55	0.50
1:2:547:G:H4'	1:2:549:C:H5	1.76	0.49
1:2:649:U:H1'	25:X:45:SER:HB3	1.94	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:F:74:ASN:HD22	7:F:74:ASN:C	2.09	0.49
12:K:23:ALA:O	12:K:67:PHE:N	2.31	0.49
34:g:52:TYR:OH	34:g:311:GLN:OE1	2.21	0.49
34:g:153:CYS:SG	34:g:154:VAL:N	2.85	0.49
34:g:268:ASP:OD1	34:g:268:ASP:N	2.45	0.49
1:2:605:A:H4'	1:2:639:C:H5''	1.94	0.49
1:2:1521:C:C2	20:S:137:LYS:HB3	2.47	0.49
1:2:1560:U:H2'	1:2:1561:A:H8	1.78	0.49
17:P:86:LEU:HG	17:P:88:GLU:H	1.76	0.49
18:Q:104:SER:O	18:Q:108:ILE:HG13	2.12	0.49
21:T:83:GLN:HB3	21:T:90:SER:HB2	1.94	0.49
1:2:823:U:C5	11:J:143:ASN:HB3	2.47	0.49
1:2:943:U:C2	1:2:944:A:C8	3.01	0.49
1:2:1232:U:H2'	1:2:1233:G:C8	2.47	0.49
1:2:1617:G:O2'	1:2:1619:A:N1	2.36	0.49
1:2:29:G:H2'	1:2:30:C:H6	1.77	0.49
1:2:1192:U:P	25:X:119:ARG:HH22	2.35	0.49
1:2:1300:U:O2'	17:P:51:ARG:NH2	2.46	0.49
4:C:98:LEU:HD13	4:C:102:LEU:HD21	1.92	0.49
4:C:183:LYS:HA	4:C:195:LEU:O	2.12	0.49
5:D:46:THR:OG1	5:D:83:SER:O	2.30	0.49
7:F:74:ASN:O	7:F:77:MET:HG2	2.12	0.49
7:F:192:LYS:HE3	7:F:196:LEU:HD11	1.93	0.49
14:M:22:LEU:HD21	14:M:102:LYS:HB3	1.94	0.49
20:S:112:GLU:HG2	20:S:113:ARG:H	1.77	0.49
1:2:5:U:H2'	1:2:6:G:C8	2.47	0.49
1:2:54:A:OP1	26:Y:111:LYS:NZ	2.38	0.49
1:2:643:A:OP1	11:J:41:ARG:NH2	2.45	0.49
7:F:115:ALA:HB3	7:F:182:LYS:N	2.27	0.49
8:G:56:ASN:ND2	8:G:60:GLY:O	2.45	0.49
30:c:34:PHE:HD2	30:c:39:SER:HA	1.78	0.49
1:2:1110:G:H4'	19:R:124:VAL:HG12	1.93	0.49
1:2:1217:A:H2'	1:2:1218:C:H6	1.77	0.49
1:2:1692:U:H2'	1:2:1693:G:C8	2.48	0.49
13:L:105:ARG:HG3	25:X:11:ARG:HH22	1.78	0.49
21:T:63:HIS:ND1	21:T:75:MET:HB2	2.28	0.49
1:2:125:C:H5''	8:G:198:ARG:HD2	1.93	0.49
1:2:1144:A:H2'	1:2:1145:A:C8	2.48	0.49
1:2:1447:G:HO2'	1:2:1448:A:H8	1.59	0.49
2:A:66:VAL:HG21	2:A:185:MET:HB3	1.95	0.49
7:F:117:ILE:HG23	7:F:183:GLY:H	1.76	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:F:145:ARG:HE	30:c:47:LYS:HG3	1.77	0.49
15:N:145:THR:HA	15:N:148:ALA:HB3	1.95	0.49
20:S:80:PRO:HD2	21:T:33:TRP:HZ2	1.76	0.49
1:2:76:U:O2'	8:G:154:ARG:NH2	2.45	0.49
1:2:375:U:H2'	1:2:376:A:C8	2.47	0.49
1:2:77:A:C2	8:G:175:LYS:HG3	2.45	0.49
1:2:106:C:H2'	1:2:107:A:H8	1.78	0.49
1:2:521:A:OP1	11:J:45:ARG:NH1	2.42	0.49
1:2:1450:G:OP2	19:R:44:LYS:NZ	2.40	0.49
36:n:163:THR:HG22	36:n:164:LYS:N	2.27	0.49
1:2:571:U:O2'	26:Y:60:PHE:O	2.31	0.49
1:2:1037:G:H4'	1:2:1845:A:H4'	1.94	0.49
1:2:1387:G:H22	5:D:206:ASP:HB2	1.78	0.49
1:2:1618:C:H5''	1:2:1619:A:C2	2.41	0.49
1:2:1856:C:H2'	1:2:1857:G:H8	1.77	0.49
5:D:139:SER:HB3	5:D:149:SER:HB2	1.94	0.49
6:E:31:PRO:HB3	6:E:43:PRO:HG3	1.95	0.49
9:H:30:LEU:HD11	9:H:82:GLU:OE1	2.12	0.49
18:Q:13:PHE:HB2	18:Q:23:ALA:H	1.75	0.49
1:2:221:A:N6	1:2:297:A:H62	2.11	0.48
1:2:372:U:O2'	13:L:82:MET:SD	2.68	0.48
1:2:443:U:H2'	1:2:444:G:O4'	2.13	0.48
6:E:112:HIS:NE2	6:E:237:SER:O	2.46	0.48
13:L:68:ILE:HG13	13:L:143:LEU:HD11	1.94	0.48
15:N:46:THR:HB	15:N:86:GLU:HG3	1.95	0.48
1:2:549:C:H2'	1:2:550:C:O4'	2.13	0.48
1:2:639:C:HO2'	1:2:640:A:H8	1.59	0.48
6:E:22:LYS:HG3	6:E:23:LEU:HD23	1.94	0.48
13:L:91:ASP:HB3	13:L:104:LYS:HE3	1.95	0.48
15:N:5:HIS:HB3	15:N:117:LEU:HD13	1.94	0.48
18:Q:16:LYS:O	18:Q:18:THR:N	2.46	0.48
23:V:72:LEU:O	23:V:76:ASP:HB2	2.12	0.48
1:2:558:G:H2'	1:2:559:G:C8	2.48	0.48
1:2:1743:G:H5''	1:2:1744:G:C8	2.48	0.48
3:B:33:VAL:HG11	3:B:67:PHE:CZ	2.48	0.48
8:G:213:LEU:O	8:G:217:MET:HG3	2.13	0.48
1:2:70:G:H21	1:2:79:A:H62	1.61	0.48
1:2:375:U:H2'	1:2:376:A:H8	1.79	0.48
13:L:49:GLU:O	13:L:53:GLY:N	2.46	0.48
20:S:114:LEU:HD13	20:S:122:GLY:HA2	1.96	0.48
30:c:29:GLN:HE22	30:c:43:ILE:HB	1.78	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:367:U:H4'	1:2:371:A:C8	2.49	0.48
1:2:453:C:O2'	8:G:92:ARG:O	2.23	0.48
4:C:209:VAL:HB	4:C:210:PRO:HD3	1.94	0.48
19:R:127:ASN:HD22	19:R:127:ASN:C	2.21	0.48
20:S:94:LYS:HG2	20:S:95:TYR:H	1.78	0.48
25:X:81:ILE:HD12	25:X:120:PHE:CE2	2.49	0.48
26:Y:6:THR:HB	26:Y:28:LEU:HB2	1.95	0.48
1:2:1665:G:H21	1:2:1665:G:P	2.37	0.48
9:H:148:LEU:N	24:W:49:GLU:OE2	2.45	0.48
1:2:433:A:H2'	1:2:434:G:C8	2.48	0.48
1:2:441:C:H2'	1:2:442:C:C6	2.49	0.48
1:2:963:A:H8	1:2:963:A:OP1	1.96	0.48
1:2:1740:C:H41	1:2:1742:C:H5	1.62	0.48
9:H:32:MET:HG2	9:H:32:MET:H	1.49	0.48
9:H:105:THR:OG1	9:H:107:LYS:O	2.30	0.48
20:S:36:VAL:HG11	20:S:64:VAL:HG12	1.95	0.48
21:T:134:ILE:O	21:T:138:VAL:HG13	2.13	0.48
2:A:40:LYS:NZ	19:R:101:ASP:OD1	2.44	0.48
7:F:97:PHE:O	7:F:100:ILE:HG13	2.14	0.48
9:H:191:GLU:O	9:H:193:GLN:NE2	2.47	0.48
1:2:119:U:O2	6:E:33:THR:OG1	2.27	0.48
1:2:1537:A:O2'	21:T:44:GLU:OE1	2.32	0.48
9:H:11:PRO:HG3	9:H:16:PRO:HG3	1.95	0.48
22:U:98:VAL:HA	22:U:101:ILE:HG22	1.96	0.48
34:g:164:ILE:HD11	34:g:200:VAL:HG21	1.95	0.48
1:2:1653:U:O2'	21:T:82:ARG:HB3	2.14	0.48
5:D:21:LEU:O	5:D:25:LEU:HG	2.14	0.48
6:E:79:ASP:HB3	6:E:82:TYR:HB2	1.96	0.48
6:E:124:CYS:HB3	6:E:141:THR:HB	1.96	0.48
9:H:84:GLU:OE2	9:H:91:HIS:ND1	2.27	0.48
10:I:107:THR:O	10:I:111:GLN:HG3	2.13	0.48
11:J:113:GLN:OE1	11:J:154:GLN:NE2	2.47	0.48
18:Q:108:ILE:HA	18:Q:111:ILE:HG22	1.96	0.48
20:S:21:ASP:O	20:S:23:ARG:HD2	2.14	0.48
24:W:39:THR:O	24:W:43:LYS:HG2	2.14	0.48
1:2:582:U:H2'	1:2:583:A:C8	2.49	0.47
1:2:604:A:H2'	1:2:605:A:C8	2.48	0.47
1:2:1588:A:C5	21:T:41:LYS:HE2	2.49	0.47
7:F:117:ILE:HG12	7:F:183:GLY:HA2	1.96	0.47
20:S:124:ARG:HA	20:S:129:LEU:HB2	1.96	0.47
21:T:7:LYS:HD2	21:T:65:TYR:CE2	2.49	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:c:45:ASN:HB2	30:c:63:ARG:HH22	1.79	0.47
1:2:1588:A:O4'	21:T:94:ARG:NE	2.46	0.47
5:D:85:GLU:OE1	5:D:85:GLU:N	2.47	0.47
1:2:225:G:N2	1:2:294:U:OP2	2.47	0.47
1:2:578:C:H2'	1:2:579:C:C6	2.49	0.47
1:2:1495:G:H4'	31:d:26:ASN:ND2	2.29	0.47
4:C:61:MET:HE3	4:C:62:PRO:HD2	1.96	0.47
7:F:122:ARG:H	7:F:197:GLU:CD	2.22	0.47
10:I:153:LYS:HB2	10:I:153:LYS:HE2	1.56	0.47
12:K:27:VAL:HG12	12:K:28:HIS:ND1	2.29	0.47
21:T:5:THR:HG22	21:T:7:LYS:H	1.80	0.47
21:T:53:PHE:HA	21:T:56:ARG:NE	2.28	0.47
32:e:40:ARG:HG2	32:e:40:ARG:HH11	1.79	0.47
2:A:143:PRO:HB2	23:V:34:MET:HE3	1.96	0.47
3:B:28:LYS:O	16:O:17:LEU:N	2.47	0.47
12:K:1:MET:HG3	12:K:3:MET:HG3	1.97	0.47
17:P:87:PRO:CG	17:P:112:ILE:HD11	2.44	0.47
26:Y:105:LYS:O	26:Y:109:GLU:HG3	2.14	0.47
1:2:595:U:H2'	1:2:596:U:C6	2.49	0.47
1:2:815:U:C2	1:2:816:A:C8	3.02	0.47
1:2:929:G:H2'	1:2:930:C:O4'	2.15	0.47
1:2:964:A:H2'	1:2:965:U:O2	2.15	0.47
1:2:1189:A:H2'	1:2:1190:A:C8	2.50	0.47
8:G:38:ALA:HB1	8:G:41:LEU:HD13	1.96	0.47
8:G:56:ASN:HB2	8:G:108:VAL:HG12	1.97	0.47
20:S:37:GLY:H	20:S:41:ALA:HB2	1.78	0.47
20:S:125:HIS:CD2	20:S:131:VAL:HG21	2.49	0.47
1:2:290:U:O2'	1:2:292:A:N7	2.35	0.47
1:2:641:A:O2'	1:2:645:C:OP1	2.33	0.47
1:2:1079:C:O2'	1:2:1182:A:N1	2.47	0.47
6:E:42:LEU:N	6:E:84:ALA:O	2.47	0.47
7:F:68:ILE:HA	7:F:71:ARG:HB3	1.96	0.47
9:H:54:GLY:O	9:H:57:ARG:NH2	2.48	0.47
11:J:136:ARG:NH1	11:J:138:ARG:O	2.44	0.47
15:N:87:ASP:N	15:N:87:ASP:OD1	2.47	0.47
20:S:87:GLN:HA	20:S:95:TYR:HE1	1.79	0.47
28:a:93:LYS:HB3	28:a:93:LYS:HE2	1.58	0.47
1:2:17:C:H2'	1:2:18:C:H6	1.78	0.47
1:2:302:A:H4'	10:I:73:THR:O	2.14	0.47
1:2:639:C:OP1	32:e:41:ARG:NH1	2.45	0.47
1:2:980:A:H2'	1:2:981:A:C8	2.50	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1022:U:O2'	1:2:1023:A:H5'	2.15	0.47
1:2:1365:G:H2'	1:2:1366:G:C8	2.50	0.47
1:2:1590:C:H3'	1:2:1591:C:H6	1.80	0.47
25:X:124:LYS:HG2	25:X:129:SER:HA	1.97	0.47
28:a:40:VAL:HG23	28:a:69:VAL:HB	1.97	0.47
30:c:18:LEU:N	30:c:29:GLN:O	2.48	0.47
1:2:160:U:O2'	1:2:162:C:OP2	2.27	0.47
1:2:613:G:N1	1:2:629:A:OP2	2.30	0.47
1:2:1339:U:H2'	1:2:1340:U:C6	2.50	0.47
1:2:1669:G:H5''	18:Q:130:LYS:HB3	1.97	0.47
4:C:167:ARG:NH1	4:C:218:GLY:O	2.47	0.47
8:G:22:ARG:H	8:G:22:ARG:NE	2.13	0.47
1:2:684:G:H2'	1:2:685:A:H5''	1.96	0.47
1:2:1538:C:H4'	21:T:43:LYS:HA	1.97	0.47
20:S:61:GLU:HG2	20:S:61:GLU:O	2.15	0.47
33:f:83:LYS:HE2	33:f:83:LYS:HB3	1.64	0.47
1:2:210:U:H5''	1:2:211:G:N7	2.30	0.47
1:2:638:C:H2'	1:2:639:C:H5'	1.97	0.47
1:2:1199:A:H2'	1:2:1200:A:C8	2.49	0.47
1:2:1550:G:O2'	1:2:1558:C:O2	2.28	0.47
4:C:191:VAL:HG11	4:C:236:PHE:HA	1.96	0.47
8:G:5:ILE:HD12	8:G:16:ILE:HD12	1.97	0.47
1:2:563:G:N7	11:J:172:ARG:NH2	2.63	0.46
1:2:879:C:O2'	1:2:908:A:N1	2.46	0.46
1:2:1427:C:H1'	21:T:5:THR:HG23	1.97	0.46
1:2:1543:U:O2	18:Q:77:HIS:ND1	2.47	0.46
1:2:1703:C:H2'	1:2:1704:C:O4'	2.15	0.46
7:F:61:PHE:HE1	7:F:65:GLN:HG3	1.80	0.46
17:P:84:ILE:C	17:P:86:LEU:N	2.73	0.46
21:T:84:ARG:HH22	21:T:89:PRO:HA	1.80	0.46
1:2:482:G:OP1	25:X:76:LYS:HA	2.14	0.46
1:2:1535:U:C6	7:F:88:MET:HE3	2.50	0.46
1:2:1603:G:O5'	1:2:1630:A:N6	2.48	0.46
5:D:54:ARG:HG3	5:D:54:ARG:HH11	1.79	0.46
20:S:79:ILE:HG23	21:T:33:TRP:CZ2	2.50	0.46
21:T:35:ASP:HA	21:T:52:TRP:CZ3	2.50	0.46
21:T:101:ARG:HH22	21:T:121:ARG:HH22	1.63	0.46
34:g:285:GLN:N	34:g:285:GLN:NE2	2.63	0.46
1:2:954:U:O2'	1:2:955:A:H5'	2.16	0.46
1:2:1450:G:H4'	34:g:64:HIS:HE2	1.81	0.46
1:2:1866:A:N6	28:a:84:VAL:HB	2.31	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:30:TRP:CE2	16:O:19:PRO:HD3	2.51	0.46
7:F:143:PRO:O	7:F:147:VAL:HG12	2.15	0.46
26:Y:100:LYS:HD3	26:Y:102:THR:HG22	1.96	0.46
1:2:949:G:H2'	1:2:950:C:C6	2.51	0.46
1:2:1330:G:N7	1:2:1493:C:H5'	2.29	0.46
1:2:1665:G:OP1	1:2:1665:G:N2	2.37	0.46
2:A:34:MET:HE3	2:A:34:MET:HB3	1.72	0.46
2:A:57:LYS:HD3	23:V:70:LEU:HD13	1.98	0.46
7:F:49:LEU:HD13	7:F:50:PRO:HD2	1.97	0.46
8:G:217:MET:HE2	8:G:217:MET:HB3	1.79	0.46
1:2:90:G:H2'	1:2:91:A:O4'	2.16	0.46
1:2:875:A:H61	1:2:911:C:H42	1.63	0.46
1:2:1016:U:O2	29:b:30:SER:OG	2.34	0.46
1:2:1238:U:O4'	17:P:127:LYS:HD3	2.15	0.46
1:2:1845:A:H2'	1:2:1846:G:H8	1.78	0.46
5:D:105:LEU:HB2	5:D:122:VAL:HG21	1.96	0.46
6:E:212:ASP:HB2	6:E:216:ASN:HB2	1.97	0.46
11:J:119:LEU:HD23	11:J:135:ILE:HD12	1.97	0.46
19:R:58:MET:HB3	19:R:58:MET:HE2	1.69	0.46
26:Y:57:VAL:HB	26:Y:60:PHE:CE2	2.51	0.46
1:2:376:A:OP2	13:L:58:LYS:NZ	2.40	0.46
1:2:604:A:H4'	11:J:22:LYS:HD3	1.98	0.46
1:2:656:G:H5'	1:2:662:G:N2	2.30	0.46
1:2:802:A:N6	1:2:861:A:H2	2.13	0.46
1:2:1271:C:OP2	1:2:1507:G:N2	2.49	0.46
16:O:113:GLN:NE2	28:a:44:ILE:O	2.43	0.46
1:2:433:A:H5''	10:I:22:HIS:HB3	1.96	0.46
1:2:1630:A:H5''	20:S:34:LYS:HA	1.97	0.46
34:g:36:ARG:HH12	34:g:64:HIS:CE1	2.34	0.46
34:g:37:ASP:O	34:g:60:ARG:NH2	2.49	0.46
1:2:4:C:H4'	4:C:207:ALA:HB2	1.98	0.46
1:2:1049:A:OP2	1:2:1068:G:N1	2.27	0.46
1:2:1589:A:O2'	1:2:1590:C:O5'	2.34	0.46
7:F:176:GLU:O	7:F:180:ALA:CB	2.63	0.46
28:a:66:LYS:HB3	28:a:66:LYS:HE3	1.79	0.46
1:2:500:A:H1'	1:2:501:C:OP1	2.16	0.46
1:2:1018:U:OP2	15:N:62:GLN:NE2	2.49	0.46
1:2:1678:A:O2'	1:2:1679:A:H5'	2.14	0.46
6:E:62:LYS:HB3	6:E:62:LYS:HE3	1.53	0.46
7:F:60:ARG:O	7:F:61:PHE:CD2	2.69	0.46
9:H:30:LEU:HD21	9:H:82:GLU:OE1	2.15	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:I:21:TYR:CE2	10:I:22:HIS:HD2	2.34	0.46
24:W:30:CYS:SG	24:W:31:SER:N	2.89	0.46
34:g:118:ARG:HG2	34:g:119:GLN:HG2	1.98	0.46
1:2:1329:U:O2'	1:2:1332:A:OP2	2.22	0.46
1:2:1365:G:H2'	1:2:1366:G:H8	1.80	0.46
1:2:1534:C:H3'	1:2:1535:U:C6	2.50	0.46
5:D:193:ASP:OD2	5:D:198:ILE:HG22	2.16	0.46
7:F:113:VAL:O	7:F:183:GLY:HA3	2.16	0.46
9:H:91:HIS:CD2	9:H:172:THR:HG21	2.51	0.46
21:T:83:GLN:HG2	21:T:85:ASN:OD1	2.16	0.46
1:2:57:U:OP1	1:2:504:G:O2'	2.31	0.45
1:2:305:U:H5'	10:I:41:ARG:NH1	2.29	0.45
10:I:148:LYS:HA	10:I:151:GLU:OE1	2.16	0.45
12:K:18:GLU:HG2	12:K:21:MET:HE3	1.98	0.45
17:P:100:LYS:HG3	17:P:101:THR:HG23	1.97	0.45
18:Q:51:LEU:HD11	18:Q:88:ILE:HD12	1.96	0.45
27:Z:73:VAL:HA	27:Z:77:LEU:HB3	1.98	0.45
30:c:51:ARG:HH21	30:c:54:ASP:HA	1.81	0.45
1:2:124:U:OP1	8:G:201:LYS:NZ	2.42	0.45
1:2:1025:U:H2'	1:2:1026:C:O4'	2.16	0.45
1:2:1092:G:OP2	15:N:9:LYS:NZ	2.43	0.45
1:2:1164:G:O2'	1:2:1165:G:H5'	2.15	0.45
1:2:1517:G:P	1:2:1517:G:H21	2.38	0.45
1:2:1800:A:H3'	1:2:1801:A:H5''	1.98	0.45
1:2:1824:A:H2'	1:2:1826:G:C6	2.51	0.45
5:D:179:GLN:OE1	5:D:179:GLN:HA	2.15	0.45
8:G:14:LYS:HB2	8:G:124:LEU:HD11	1.98	0.45
17:P:44:ARG:HE	17:P:83:MET:HB2	1.82	0.45
1:2:674:C:H2'	1:2:675:U:C6	2.51	0.45
2:A:75:SER:HB2	2:A:119:PRO:HB3	1.99	0.45
13:L:18:GLN:C	13:L:19:ASN:HD22	2.25	0.45
18:Q:68:ILE:HG12	18:Q:70:VAL:HG13	1.97	0.45
23:V:42:VAL:HG13	23:V:43:THR:HG23	1.98	0.45
29:b:36:LYS:HD3	29:b:36:LYS:HA	1.66	0.45
30:c:34:PHE:HB2	30:c:39:SER:HA	1.96	0.45
34:g:105:THR:HG23	34:g:106:LYS:HG2	1.97	0.45
12:K:49:MET:HB3	12:K:49:MET:HE2	1.71	0.45
13:L:126:VAL:HG12	13:L:145:VAL:HG22	1.98	0.45
19:R:45:LYS:HE3	19:R:45:LYS:HB2	1.59	0.45
21:T:60:THR:O	21:T:64:LEU:HB2	2.16	0.45
1:2:500:A:H1'	1:2:501:C:P	2.57	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:500:A:O2'	1:2:501:C:OP1	2.31	0.45
1:2:1019:C:H2'	1:2:1020:A:O4'	2.15	0.45
1:2:1394:G:H5''	18:Q:126:ARG:NH1	2.31	0.45
1:2:1448:A:OP1	22:U:30:LYS:NZ	2.34	0.45
1:2:1587:G:H1'	21:T:94:ARG:HG2	1.99	0.45
1:2:1602:U:H2'	1:2:1603:G:H5''	1.97	0.45
6:E:100:ARG:HD3	6:E:102:ILE:HD11	1.98	0.45
20:S:25:LYS:HG2	20:S:26:ILE:H	1.82	0.45
1:2:71:G:N7	8:G:170:ARG:NH2	2.65	0.45
1:2:943:U:O2'	16:O:135:ILE:O	2.34	0.45
1:2:1269:G:H3'	33:f:83:LYS:HZ3	1.82	0.45
6:E:199:GLU:OE2	6:E:209:HIS:NE2	2.42	0.45
11:J:131:ARG:HA	11:J:131:ARG:HD2	1.59	0.45
20:S:20:ILE:HD12	20:S:20:ILE:HA	1.85	0.45
1:2:81:U:H2'	1:2:82:G:O4'	2.17	0.45
1:2:388:U:H2'	1:2:389:A:H8	1.81	0.45
1:2:1101:U:H2'	1:2:1102:G:H8	1.82	0.45
1:2:1452:A:H5''	19:R:48:ASN:ND2	2.32	0.45
7:F:144:LEU:HB3	30:c:49:PRO:HB2	1.97	0.45
11:J:28:GLU:OE2	11:J:44:TRP:NE1	2.46	0.45
18:Q:98:LYS:HA	18:Q:98:LYS:HD2	1.85	0.45
24:W:30:CYS:HB2	24:W:61:ILE:HG13	1.99	0.45
34:g:149:GLU:HB3	34:g:171:ASP:HB2	1.98	0.45
1:2:419:G:N2	1:2:661:U:O2	2.50	0.45
1:2:1020:A:H1'	1:2:1022:U:O4	2.17	0.45
1:2:1388:A:C2	5:D:205:PRO:HG2	2.51	0.45
1:2:1820:G:H2'	1:2:1821:U:C6	2.52	0.45
2:A:198:MET:HE3	19:R:85:VAL:HB	1.99	0.45
11:J:31:LEU:HD13	11:J:102:ILE:HD13	1.99	0.45
1:2:65:C:C5	8:G:174:PRO:HB3	2.52	0.45
1:2:345:U:O2'	1:2:346:C:OP1	2.31	0.45
1:2:1447:G:O2'	1:2:1448:A:H8	2.00	0.45
8:G:130:PRO:HG2	8:G:132:ARG:HH11	1.82	0.45
8:G:201:LYS:HB3	8:G:201:LYS:HE3	1.78	0.45
21:T:120:GLY:C	21:T:121:ARG:HD2	2.42	0.45
35:h:16:LYS:HD3	35:h:16:LYS:HA	1.81	0.45
1:2:106:C:H2'	1:2:107:A:C8	2.52	0.45
1:2:1222:G:H2'	1:2:1223:A:C8	2.52	0.45
1:2:1397:U:O2'	18:Q:12:VAL:HB	2.17	0.45
4:C:215:MET:HE3	4:C:215:MET:HB3	1.75	0.45
5:D:106:ARG:HB2	5:D:184:ILE:HD11	1.98	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:T:22:LEU:HB3	21:T:54:TYR:OH	2.17	0.45
26:Y:3:ASP:HB3	26:Y:4:THR:H	1.59	0.45
30:c:42:ILE:HB	30:c:64:GLU:OE2	2.17	0.45
1:2:981:A:H2'	1:2:982:G:C8	2.52	0.44
3:B:89:GLU:HG3	3:B:99:ASN:HB2	1.99	0.44
5:D:158:ILE:HG13	5:D:164:VAL:HG22	1.98	0.44
9:H:40:LEU:O	9:H:41:ARG:C	2.59	0.44
18:Q:16:LYS:HD2	18:Q:82:TYR:O	2.16	0.44
18:Q:80:GLN:NE2	18:Q:81:ILE:HG13	2.32	0.44
21:T:85:ASN:HB2	21:T:88:MET:SD	2.57	0.44
32:e:58:ASN:OD1	32:e:58:ASN:N	2.49	0.44
1:2:1:U:H5'	1:2:417:C:N4	2.33	0.44
1:2:77:A:OP1	8:G:155:GLN:NE2	2.50	0.44
1:2:170:A:OP2	8:G:140:ARG:NH1	2.50	0.44
1:2:420:G:O2'	1:2:660:C:N3	2.37	0.44
1:2:606:G:OP1	32:e:54:GLY:HA3	2.17	0.44
1:2:1615:U:H5''	1:2:1616:U:O4'	2.16	0.44
1:2:1619:A:O2'	1:2:1620:A:OP1	2.36	0.44
3:B:136:ARG:HB3	3:B:216:LYS:HG3	1.98	0.44
5:D:185:LYS:HE3	5:D:185:LYS:HB2	1.59	0.44
9:H:51:ILE:HG21	9:H:179:LYS:HE2	1.99	0.44
9:H:85:LYS:HB2	9:H:85:LYS:HE3	1.82	0.44
11:J:151:LEU:HD23	11:J:151:LEU:HA	1.79	0.44
20:S:103:LEU:O	20:S:107:LEU:HG	2.17	0.44
21:T:11:GLN:CD	21:T:62:ARG:HH21	2.25	0.44
22:U:26:SER:HB3	22:U:32:LEU:HD13	1.99	0.44
30:c:10:LYS:HB2	30:c:58:LEU:HB2	1.99	0.44
34:g:131:LEU:HD12	34:g:140:TYR:HB3	1.98	0.44
1:2:289:G:N2	1:2:293:C:H5	2.14	0.44
1:2:1204:A:O2'	1:2:1700:C:OP2	2.29	0.44
1:2:1545:A:O2'	1:2:1546:G:H5'	2.17	0.44
1:2:1673:U:H2'	1:2:1674:G:O4'	2.17	0.44
3:B:139:CYS:HB2	3:B:172:MET:SD	2.56	0.44
5:D:33:GLY:O	5:D:53:THR:HG23	2.17	0.44
6:E:254:LYS:HE2	6:E:254:LYS:HB3	1.62	0.44
10:I:23:LYS:HB3	10:I:23:LYS:HE2	1.57	0.44
14:M:121:LYS:HB2	14:M:121:LYS:HE2	1.77	0.44
14:M:121:LYS:O	14:M:124:ILE:HG13	2.17	0.44
19:R:20:TYR:O	19:R:24:LEU:HG	2.17	0.44
20:S:43:VAL:HA	21:T:35:ASP:OD2	2.17	0.44
1:2:824:C:C2	11:J:144:ILE:HG13	2.52	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1182:A:C5	1:2:1183:A:H1'	2.53	0.44
1:2:1203:G:H2'	1:2:1204:A:H8	1.81	0.44
1:2:1627:C:H4'	21:T:41:LYS:HD3	1.98	0.44
2:A:42:LYS:HE3	19:R:99:ASP:OD2	2.16	0.44
5:D:59:LEU:HD21	5:D:88:ALA:HB3	1.98	0.44
8:G:78:SER:HB2	8:G:92:ARG:HG2	1.98	0.44
15:N:83:ASP:OD1	15:N:83:ASP:N	2.49	0.44
23:V:19:ALA:HB2	23:V:68:SER:OG	2.18	0.44
28:a:78:VAL:HG13	28:a:84:VAL:HG22	1.99	0.44
29:b:5:LYS:HG2	29:b:7:LEU:HD23	1.99	0.44
30:c:28:THR:HB	30:c:46:VAL:C	2.40	0.44
32:e:18:LYS:HE2	32:e:18:LYS:HB2	1.67	0.44
14:M:40:LYS:HE2	33:f:130:VAL:HA	1.99	0.44
1:2:434:G:H5''	10:I:23:LYS:HE3	1.98	0.44
1:2:1308:U:O4	33:f:95:ARG:HD3	2.18	0.44
1:2:1378:A:H4'	1:2:1379:A:O5'	2.18	0.44
1:2:1546:G:H8	1:2:1670:C:O2'	2.00	0.44
1:2:1609:C:O2'	20:S:109:GLU:OE2	2.20	0.44
5:D:205:PRO:HA	19:R:42:PRO:HG2	1.99	0.44
34:g:67:SER:OG	34:g:68:ASP:N	2.51	0.44
1:2:501:C:O2	1:2:501:C:C2'	2.65	0.44
1:2:1533:A:C8	1:2:1604:G:H1'	2.52	0.44
1:2:1824:A:H2'	1:2:1826:G:N1	2.33	0.44
8:G:65:GLN:O	8:G:65:GLN:NE2	2.50	0.44
12:K:11:ILE:HD13	12:K:45:VAL:HG13	2.00	0.44
16:O:75:MET:HG3	16:O:76:LEU:N	2.31	0.44
18:Q:41:MET:HE1	21:T:10:ASN:HB2	2.00	0.44
34:g:291:TRP:HB3	34:g:295:GLY:HA2	1.99	0.44
1:2:108:G:H1'	1:2:853:C:H42	1.82	0.44
1:2:1013:U:C2	1:2:1014:G:C8	3.06	0.44
1:2:1022:U:O2'	1:2:1022:U:O2	2.27	0.44
1:2:1516:G:H8	17:P:99:GLY:O	1.99	0.44
2:A:94:THR:O	2:A:186:ARG:NH2	2.50	0.44
20:S:83:PHE:CE1	21:T:36:THR:HB	2.52	0.44
34:g:40:ILE:O	34:g:59:LEU:HB2	2.18	0.44
1:2:92:A:H4'	1:2:93:U:OP2	2.17	0.44
3:B:61:GLY:O	3:B:65:ARG:NH2	2.45	0.44
4:C:253:PRO:HA	4:C:256:TRP:CE2	2.53	0.44
7:F:188:TYR:O	7:F:191:LYS:HG3	2.18	0.44
10:I:205:ARG:HD2	10:I:205:ARG:HA	1.64	0.44
13:L:20:LYS:HE3	13:L:20:LYS:HB2	1.77	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:b:36:LYS:O	29:b:77:CYS:HA	2.18	0.44
1:2:223:C:H3'	1:2:225:G:OP2	2.18	0.43
1:2:659:G:H2'	1:2:663:C:C5	2.53	0.43
1:2:1324:G:O2'	1:2:1325:G:H5'	2.18	0.43
4:C:212:LYS:O	4:C:216:MET:HG3	2.18	0.43
5:D:211:VAL:O	19:R:20:TYR:OH	2.31	0.43
9:H:39:GLN:HG2	9:H:75:ILE:HG21	1.99	0.43
9:H:80:VAL:HG22	9:H:92:VAL:HG12	2.00	0.43
9:H:82:GLU:O	9:H:85:LYS:HB3	2.18	0.43
27:Z:49:LEU:H	27:Z:49:LEU:HD12	1.82	0.43
28:a:75:VAL:O	28:a:78:VAL:HG12	2.17	0.43
30:c:17:VAL:HG13	30:c:30:VAL:HG12	1.99	0.43
32:e:36:MET:HE3	32:e:36:MET:HB3	1.94	0.43
34:g:32:LEU:HD22	34:g:40:ILE:CG2	2.48	0.43
1:2:96:C:H2'	1:2:97:U:H6	1.83	0.43
1:2:113:G:N2	1:2:293:C:H5'	2.32	0.43
1:2:598:G:O2'	1:2:605:A:N1	2.48	0.43
1:2:630:U:O4	36:n:165:TYR:HA	2.18	0.43
1:2:1718:G:H22	1:2:1814:G:P	2.41	0.43
1:2:1797:U:H2'	1:2:1798:C:C6	2.53	0.43
4:C:166:ARG:NH2	4:C:254:ASP:OD2	2.44	0.43
5:D:122:VAL:O	5:D:126:ILE:HG12	2.17	0.43
6:E:107:GLY:HA2	6:E:189:LEU:HG	1.99	0.43
7:F:110:GLN:HG3	7:F:114:ASN:ND2	2.33	0.43
11:J:111:GLN:CD	11:J:127:ARG:HB2	2.43	0.43
11:J:134:HIS:ND1	11:J:163:SER:HB2	2.32	0.43
17:P:123:TYR:HD1	17:P:123:TYR:O	2.02	0.43
18:Q:13:PHE:CG	18:Q:22:VAL:HA	2.53	0.43
21:T:84:ARG:NE	21:T:84:ARG:HA	2.33	0.43
25:X:7:LEU:HD12	25:X:7:LEU:HA	1.83	0.43
34:g:24:THR:HG23	34:g:27:PHE:H	1.82	0.43
34:g:39:THR:HG23	34:g:59:LEU:C	2.42	0.43
34:g:90:TRP:CZ3	34:g:96:THR:HA	2.53	0.43
1:2:1232:U:H2'	1:2:1233:G:H8	1.83	0.43
1:2:1269:G:N7	1:2:1270:G:N2	2.62	0.43
4:C:163:VAL:HG22	4:C:164:PRO:HD2	2.01	0.43
5:D:54:ARG:HH12	5:D:57:ASN:ND2	2.16	0.43
7:F:27:ASP:N	7:F:27:ASP:OD1	2.51	0.43
7:F:165:ASN:ND2	7:F:167:LYS:O	2.37	0.43
10:I:206:LYS:HD2	10:I:206:LYS:HA	1.84	0.43
12:K:21:MET:HE3	12:K:21:MET:HB3	1.86	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:Q:100:VAL:HG13	18:Q:101:ASP:N	2.33	0.43
26:Y:20:ARG:HH21	26:Y:22:GLN:NE2	2.16	0.43
31:d:11:PRO:HB2	31:d:13:LYS:HG2	2.00	0.43
1:2:447:A:H4'	6:E:3:ARG:HG2	1.99	0.43
1:2:849:A:H2'	1:2:850:C:H6	1.83	0.43
1:2:962:A:H1'	16:O:65:ASP:OD1	2.17	0.43
1:2:1615:U:H3'	1:2:1616:U:H4'	1.99	0.43
7:F:100:ILE:O	7:F:104:THR:OG1	2.32	0.43
16:O:64:ALA:HB3	16:O:67:ASP:OD1	2.19	0.43
28:a:36:ILE:HG21	28:a:78:VAL:HG21	1.99	0.43
34:g:189:ILE:HG22	34:g:190:GLY:H	1.82	0.43
1:2:1249:C:H2'	1:2:1250:A:H2	1.81	0.43
10:I:83:TYR:HB3	10:I:101:ILE:HB	1.99	0.43
11:J:22:LYS:HA	11:J:22:LYS:HD2	1.72	0.43
22:U:94:PRO:HD2	22:U:97:ILE:HD13	1.99	0.43
34:g:156:PHE:HA	34:g:164:ILE:O	2.18	0.43
34:g:198:VAL:HG21	34:g:208:ALA:O	2.18	0.43
1:2:582:U:H1'	26:Y:33:ALA:HB2	2.00	0.43
1:2:638:C:C2'	1:2:639:C:H5'	2.48	0.43
1:2:829:C:OP1	6:E:21:ASP:HB2	2.19	0.43
1:2:1517:G:H4'	1:2:1518:C:OP1	2.18	0.43
5:D:54:ARG:HG3	5:D:54:ARG:NH1	2.32	0.43
16:O:17:LEU:HB3	16:O:18:GLY:H	1.65	0.43
18:Q:82:TYR:HA	18:Q:85:ARG:HG2	2.00	0.43
29:b:34:ASP:N	29:b:34:ASP:OD1	2.51	0.43
34:g:17:TRP:HZ2	34:g:305:ASN:HA	1.83	0.43
34:g:121:VAL:HG22	34:g:131:LEU:HD23	2.00	0.43
34:g:305:ASN:O	34:g:306:LEU:HD13	2.18	0.43
1:2:472:C:O2'	1:2:474:G:OP1	2.36	0.43
1:2:918:U:H2'	1:2:919:A:O4'	2.19	0.43
1:2:1586:U:O2'	1:2:1587:G:OP2	2.31	0.43
1:2:1605:G:N1	1:2:1635:C:O2	2.51	0.43
1:2:1679:A:P	7:F:60:ARG:HH11	2.41	0.43
7:F:117:ILE:N	7:F:183:GLY:H	2.06	0.43
20:S:55:ARG:HH11	27:Z:83:LEU:HD13	1.84	0.43
34:g:118:ARG:NH1	34:g:179:LEU:HD13	2.34	0.43
1:2:1022:U:H4'	15:N:128:TYR:HB2	1.99	0.43
1:2:1562:C:H2'	1:2:1563:G:C8	2.54	0.43
5:D:109:LEU:HD22	5:D:115:VAL:HG12	2.00	0.43
12:K:24:LYS:HG3	12:K:66:HIS:CE1	2.54	0.43
17:P:37:TYR:CD2	17:P:112:ILE:HG13	2.53	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:c:50:VAL:HG13	30:c:54:ASP:CG	2.43	0.43
33:f:89:LYS:HD3	33:f:89:LYS:HA	1.86	0.43
34:g:88:ARG:HG3	34:g:97:THR:HG21	2.01	0.43
1:2:1240:A:H1'	1:2:1516:G:O6	2.19	0.43
1:2:1630:A:N7	20:S:31:THR:HA	2.34	0.43
1:2:1851:A:H5''	35:h:1:MET:HG3	2.01	0.43
3:B:110:MET:HA	3:B:113:MET:HE2	2.01	0.43
9:H:76:GLN:HG3	9:H:94:PHE:CD1	2.54	0.43
13:L:88:ILE:HD11	13:L:109:MET:HE2	2.01	0.43
21:T:116:ASP:OD1	21:T:116:ASP:N	2.45	0.43
21:T:133:ARG:HA	21:T:133:ARG:HD2	1.58	0.43
25:X:98:ASP:OD2	25:X:140:ARG:NH2	2.52	0.43
25:X:107:ARG:HB3	25:X:110:HIS:HB3	2.00	0.43
1:2:520:A:H5''	11:J:12:THR:HG23	2.00	0.43
1:2:864:A:H4'	1:2:864:A:OP1	2.19	0.43
1:2:910:G:H2'	1:2:911:C:C5	2.54	0.43
1:2:1146:C:O2'	1:2:1150:A:N1	2.47	0.43
1:2:1189:A:H2'	1:2:1190:A:H8	1.84	0.43
1:2:1265:A:H5'	33:f:81:LYS:HE2	2.01	0.43
1:2:1298:G:C8	17:P:79:HIS:HD2	2.37	0.43
6:E:31:PRO:HG2	6:E:38:LEU:HG	2.00	0.43
7:F:147:VAL:CG2	7:F:151:ILE:HD13	2.42	0.43
21:T:66:LEU:C	21:T:67:ARG:HD2	2.44	0.43
22:U:24:LEU:HD23	22:U:112:VAL:HG22	2.01	0.43
25:X:125:VAL:HG12	25:X:130:LEU:HD12	2.00	0.43
1:2:39:A:H2'	1:2:40:A:O4'	2.18	0.42
1:2:162:C:H2'	1:2:163:U:O4'	2.19	0.42
1:2:407:G:C5	25:X:36:LEU:HD21	2.54	0.42
1:2:1105:G:O3'	29:b:69:GLY:HA3	2.18	0.42
4:C:81:ILE:HG23	4:C:86:LEU:HB2	2.01	0.42
9:H:83:LEU:HD13	9:H:83:LEU:HA	1.92	0.42
19:R:80:ARG:C	19:R:80:ARG:HD2	2.43	0.42
1:2:1622:U:H3	17:P:81:ARG:NH1	2.17	0.42
16:O:67:ASP:OD1	16:O:67:ASP:N	2.50	0.42
18:Q:22:VAL:HB	18:Q:71:ARG:HB2	2.01	0.42
19:R:114:LEU:HD12	19:R:117:LEU:HD11	2.00	0.42
1:2:1487:A:H4'	1:2:1488:C:OP1	2.19	0.42
1:2:1589:A:H2'	1:2:1589:A:N3	2.34	0.42
1:2:1609:C:OP1	20:S:130:ARG:HA	2.18	0.42
5:D:23:GLU:O	5:D:27:ARG:HG2	2.18	0.42
6:E:143:ASP:OD1	6:E:143:ASP:O	2.38	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:L:49:GLU:HG3	13:L:118:ARG:HD2	2.01	0.42
21:T:82:ARG:HD2	21:T:82:ARG:C	2.45	0.42
33:f:136:PHE:CD1	33:f:137:ASP:N	2.86	0.42
1:2:145:G:OP1	8:G:139:SER:OG	2.34	0.42
1:2:846:G:H2'	6:E:19:MET:HG2	2.01	0.42
1:2:921:G:C2	29:b:22:LYS:HD3	2.54	0.42
2:A:200:ASP:OD1	19:R:89:SER:OG	2.27	0.42
3:B:180:ASP:OD1	3:B:183:GLU:HG3	2.19	0.42
7:F:44:LYS:HD2	18:Q:117:ARG:HB2	2.01	0.42
7:F:51:HIS:HD2	7:F:90:VAL:HG21	1.84	0.42
7:F:141:VAL:HG11	7:F:145:ARG:CB	2.50	0.42
9:H:64:VAL:O	9:H:96:ALA:HA	2.19	0.42
11:J:130:ILE:HG22	11:J:143:ASN:O	2.19	0.42
12:K:44:HIS:O	12:K:48:ALA:HB2	2.19	0.42
17:P:84:ILE:C	17:P:86:LEU:H	2.26	0.42
27:Z:107:VAL:HB	27:Z:109:TYR:CE2	2.54	0.42
1:2:967:C:H2'	1:2:968:U:H5'	2.01	0.42
1:2:1093:A:H2'	1:2:1094:C:C6	2.54	0.42
1:2:1603:G:O6	20:S:24:ARG:NH2	2.52	0.42
1:2:1603:G:O2'	1:2:1604:G:H5'	2.20	0.42
10:I:194:GLU:HG2	13:L:10:TYR:CD2	2.55	0.42
26:Y:103:SER:HB3	26:Y:106:GLN:HG3	2.01	0.42
1:2:802:A:C2	1:2:863:U:H5	2.38	0.42
1:2:1181:A:H2'	1:2:1182:A:C8	2.55	0.42
1:2:1269:G:H3'	33:f:83:LYS:NZ	2.35	0.42
1:2:1627:C:H1'	21:T:38:LYS:HE3	2.00	0.42
1:2:1681:U:H2'	1:2:1682:C:C6	2.55	0.42
5:D:93:THR:HB	5:D:96:LEU:HD12	2.02	0.42
5:D:130:GLY:O	5:D:191:PRO:HG3	2.18	0.42
7:F:175:ASP:O	7:F:178:ILE:HG22	2.19	0.42
14:M:49:LEU:HD11	14:M:111:VAL:HG13	2.02	0.42
15:N:98:VAL:HG11	15:N:115:LEU:HB2	2.00	0.42
18:Q:47:LEU:HD23	18:Q:47:LEU:HA	1.83	0.42
22:U:95:SER:O	22:U:98:VAL:HG12	2.19	0.42
1:2:129:C:O2	1:2:182:C:O2'	2.35	0.42
1:2:376:A:H2'	1:2:377:G:O4'	2.20	0.42
1:2:560:A:N6	1:2:588:G:H1	2.17	0.42
1:2:650:A:H2'	1:2:651:U:O4'	2.19	0.42
1:2:1208:A:H2'	1:2:1209:A:H8	1.85	0.42
1:2:1293:A:N3	1:2:1293:A:H3'	2.35	0.42
1:2:1349:G:H2'	1:2:1350:U:C6	2.55	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:E:100:ARG:HG2	6:E:102:ILE:HG12	2.01	0.42
16:O:101:GLY:HA3	16:O:134:PRO:HG2	2.02	0.42
17:P:45:LEU:HD22	17:P:85:ILE:O	2.19	0.42
1:2:1348:G:C8	1:2:1349:G:C8	3.07	0.42
1:2:1567:G:N3	1:2:1567:G:H2'	2.34	0.42
1:2:1751:C:N4	1:2:1782:G:OP2	2.53	0.42
2:A:203:PHE:CZ	19:R:91:LEU:HD21	2.54	0.42
6:E:63:LYS:O	6:E:67:GLN:NE2	2.53	0.42
16:O:39:ASP:OD1	16:O:39:ASP:C	2.63	0.42
16:O:61:LYS:HE3	16:O:76:LEU:HB3	2.01	0.42
21:T:101:ARG:NH2	21:T:121:ARG:NH2	2.68	0.42
34:g:62:HIS:CE1	34:g:82:SER:HB2	2.55	0.42
1:2:1868:U:N3	28:a:98:PRO:O	2.27	0.42
5:D:110:LEU:HD23	5:D:110:LEU:HA	1.84	0.42
5:D:215:ASP:N	5:D:215:ASP:OD1	2.52	0.42
7:F:153:LEU:HA	7:F:153:LEU:HD23	1.81	0.42
9:H:36:LEU:O	9:H:39:GLN:HB2	2.20	0.42
18:Q:8:GLN:HB3	18:Q:9:SER:H	1.56	0.42
22:U:19:ARG:HB3	22:U:117:ALA:HB3	2.01	0.42
34:g:24:THR:OG1	34:g:73:SER:O	2.32	0.42
1:2:1016:U:H5'	15:N:15:ALA:O	2.20	0.42
2:A:208:GLU:O	2:A:211:GLU:HG2	2.20	0.42
7:F:61:PHE:CE1	7:F:65:GLN:HG3	2.55	0.42
7:F:145:ARG:HG2	30:c:49:PRO:HD2	2.01	0.42
16:O:78:ALA:HB3	16:O:118:ALA:HB3	2.02	0.42
22:U:18:HIS:O	22:U:92:HIS:HA	2.20	0.42
28:a:41:ILE:HG12	28:a:68:TYR:CD2	2.54	0.42
29:b:48:SER:HG	29:b:49:HIS:CE1	2.38	0.42
34:g:23:THR:OG1	34:g:24:THR:N	2.52	0.42
1:2:113:G:O6	1:2:290:U:H1'	2.20	0.41
1:2:615:C:OP2	25:X:64:SER:OG	2.37	0.41
2:A:128:ARG:NH2	2:A:151:ASP:O	2.34	0.41
2:A:203:PHE:HZ	19:R:91:LEU:HD21	1.85	0.41
4:C:258:GLU:H	4:C:258:GLU:HG2	1.60	0.41
8:G:46:LYS:HB2	8:G:46:LYS:HE2	1.72	0.41
11:J:26:ASP:O	11:J:30:LYS:HG3	2.20	0.41
18:Q:60:LYS:HA	18:Q:63:PHE:CZ	2.55	0.41
29:b:67:THR:HG22	29:b:68:GLY:N	2.34	0.41
1:2:429:C:H2'	1:2:430:C:C6	2.55	0.41
1:2:834:C:O3'	1:2:841:G:N2	2.54	0.41
1:2:1428:G:N7	21:T:7:LYS:NZ	2.67	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:1533:A:N7	1:2:1604:G:H1'	2.35	0.41
1:2:1848:U:H2'	1:2:1850:A:OP2	2.20	0.41
8:G:6:SER:HB3	8:G:13:GLN:HG2	2.00	0.41
8:G:159:ARG:HE	8:G:171:THR:HB	1.85	0.41
9:H:47:ALA:HB3	9:H:63:PHE:HD2	1.85	0.41
11:J:33:GLY:HA3	32:e:38:TYR:CG	2.55	0.41
21:T:90:SER:O	21:T:91:HIS:ND1	2.53	0.41
22:U:19:ARG:HG2	22:U:92:HIS:NE2	2.34	0.41
22:U:56:MET:HG3	22:U:86:LYS:HE3	2.02	0.41
1:2:339:A:H3'	1:2:339:A:N3	2.35	0.41
1:2:507:G:OP2	26:Y:104:ARG:NH1	2.46	0.41
1:2:1344:A:H4'	1:2:1345:G:OP1	2.20	0.41
1:2:1474:A:H2'	1:2:1475:G:C8	2.55	0.41
1:2:1516:G:H5'	1:2:1517:G:OP2	2.20	0.41
1:2:1598:G:H2'	27:Z:80:ARG:HG3	2.02	0.41
1:2:1724:A:H3'	1:2:1724:A:N3	2.35	0.41
2:A:118:GLU:HB2	4:C:65:LYS:HD3	2.02	0.41
7:F:77:MET:HB2	7:F:84:GLY:N	2.34	0.41
10:I:184:ARG:HE	10:I:184:ARG:HB3	1.64	0.41
12:K:11:ILE:O	12:K:15:LEU:HD23	2.20	0.41
20:S:113:ARG:O	20:S:117:ILE:HG23	2.19	0.41
25:X:133:LEU:HD23	25:X:138:LYS:O	2.19	0.41
30:c:59:LEU:HD23	30:c:59:LEU:H	1.84	0.41
1:2:34:U:H4'	1:2:564:A:H4'	2.02	0.41
1:2:171:A:H5'	8:G:177:GLN:HG2	2.03	0.41
1:2:174:C:N4	1:2:175:A:H62	2.17	0.41
1:2:579:C:C2'	1:2:580:U:O5'	2.67	0.41
2:A:25:LEU:O	2:A:164:ASN:HB2	2.21	0.41
9:H:37:LYS:HZ1	9:H:37:LYS:H	1.68	0.41
13:L:80:MET:HE2	13:L:86:ILE:HG21	2.02	0.41
21:T:83:GLN:CB	21:T:90:SER:HB2	2.49	0.41
1:2:681:U:O2'	1:2:1160:U:OP1	2.36	0.41
1:2:1619:A:H2'	1:2:1620:A:H5'	2.03	0.41
1:2:1626:C:H5''	21:T:88:MET:HE3	2.02	0.41
4:C:232:THR:OG1	4:C:235:ASN:HB2	2.20	0.41
5:D:142:LEU:HD13	5:D:142:LEU:HA	1.95	0.41
6:E:37:LYS:HB2	6:E:40:GLU:HG2	2.01	0.41
7:F:70:GLU:O	7:F:74:ASN:HB3	2.21	0.41
1:2:29:G:OP1	25:X:124:LYS:NZ	2.46	0.41
1:2:318:A:C8	1:2:333:G:C2	3.08	0.41
3:B:71:LEU:HD13	3:B:84:PHE:CE1	2.56	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:147:ASN:OD1	3:B:148:ASN:N	2.53	0.41
5:D:95:GLY:HA2	5:D:101:GLN:OE1	2.20	0.41
7:F:104:THR:HG22	7:F:106:GLU:HG3	2.03	0.41
8:G:200:LYS:HE3	8:G:200:LYS:HB3	1.87	0.41
16:O:75:MET:HB2	16:O:118:ALA:HB2	2.02	0.41
18:Q:109:LYS:HG2	18:Q:113:ILE:HG12	2.03	0.41
21:T:33:TRP:H	21:T:102:ARG:HH21	1.68	0.41
31:d:38:MET:HE1	31:d:50:ILE:HD11	2.02	0.41
33:f:91:ASN:OD1	33:f:91:ASN:N	2.51	0.41
1:2:822:U:H2'	1:2:824:C:OP2	2.20	0.41
2:A:57:LYS:HD2	2:A:57:LYS:HA	1.85	0.41
10:I:66:SER:HA	10:I:73:THR:HA	2.03	0.41
18:Q:100:VAL:HG13	18:Q:101:ASP:H	1.86	0.41
19:R:99:ASP:OD1	19:R:102:THR:HG22	2.21	0.41
20:S:106:LYS:HD2	20:S:111:LEU:HD11	2.02	0.41
1:2:952:G:N2	16:O:52:THR:HG21	2.30	0.41
1:2:1138:C:O2'	23:V:61:ARG:HD2	2.21	0.41
1:2:1571:G:HO2'	1:2:1572:C:P	2.41	0.41
1:2:1755:C:O2'	1:2:1756:C:OP1	2.31	0.41
2:A:99:ILE:HD11	2:A:117:ARG:HD2	2.02	0.41
3:B:29:ASP:OD1	3:B:51:ARG:NE	2.48	0.41
7:F:142:SER:HB2	30:c:56:LEU:HD22	2.03	0.41
8:G:41:LEU:HB3	8:G:45:TRP:CD1	2.56	0.41
9:H:25:GLN:HA	9:H:28:LEU:HD12	2.02	0.41
9:H:78:ARG:HA	9:H:78:ARG:NH1	2.36	0.41
10:I:165:GLN:HE22	10:I:172:LEU:HG	1.86	0.41
12:K:24:LYS:HB2	12:K:32:HIS:NE2	2.35	0.41
17:P:43:ARG:O	17:P:43:ARG:HG2	2.20	0.41
34:g:87:LEU:HB3	34:g:101:PHE:HB2	2.03	0.41
1:2:77:A:H1'	8:G:176:ILE:HG12	2.02	0.41
1:2:484:A:OP2	1:2:484:A:H8	2.03	0.41
1:2:952:G:C6	1:2:975:G:C6	3.09	0.41
1:2:1036:A:N3	1:2:1844:U:O2'	2.51	0.41
1:2:1271:C:O2'	1:2:1272:C:O4'	2.36	0.41
1:2:1296:U:H3'	1:2:1297:U:H4'	2.02	0.41
1:2:1745:A:H1'	8:G:31:ARG:NH2	2.33	0.41
1:2:1782:G:H5''	1:2:1783:C:H5''	2.03	0.41
2:A:121:LEU:HD12	2:A:143:PRO:HG2	2.03	0.41
5:D:202:LYS:H	5:D:202:LYS:HG2	1.58	0.41
6:E:19:MET:SD	6:E:108:ARG:HD2	2.60	0.41
7:F:116:ILE:HA	7:F:119:SER:HB2	2.03	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:F:141:VAL:HG11	7:F:145:ARG:HB3	2.02	0.41
7:F:173:LEU:HG	7:F:174:ALA:N	2.36	0.41
13:L:75:GLY:HA3	13:L:88:ILE:HD12	2.03	0.41
14:M:103:VAL:HG13	14:M:104:VAL:HG23	2.02	0.41
16:O:93:LEU:HD11	16:O:124:MET:HE3	2.03	0.41
17:P:45:LEU:HB2	17:P:85:ILE:HA	2.01	0.41
17:P:87:PRO:HD2	17:P:112:ILE:HD11	2.03	0.41
33:f:133:ALA:HB3	33:f:139:HIS:HE2	1.86	0.41
34:g:14:HIS:CE1	34:g:41:ILE:HG21	2.56	0.41
34:g:310:TRP:N	34:g:310:TRP:CD1	2.88	0.41
6:E:23:LEU:HD22	6:E:23:LEU:HA	1.89	0.41
7:F:51:HIS:CD2	7:F:90:VAL:HG21	2.56	0.41
10:I:36:THR:HG21	10:I:60:LEU:HD12	2.03	0.41
10:I:172:LEU:HD23	10:I:172:LEU:HA	1.91	0.41
15:N:86:GLU:OE1	15:N:90:HIS:ND1	2.52	0.41
16:O:142:ARG:HE	16:O:142:ARG:HB3	1.73	0.41
20:S:7:GLU:HG3	20:S:58:GLU:HB3	2.03	0.41
22:U:51:LYS:HE3	22:U:51:LYS:HB3	1.75	0.41
1:2:120:U:H1'	6:E:33:THR:O	2.22	0.40
1:2:875:A:N6	1:2:911:C:H42	2.19	0.40
6:E:95:THR:HG22	26:Y:16:ARG:HB2	2.03	0.40
7:F:109:LEU:O	7:F:113:VAL:HG12	2.21	0.40
9:H:8:ILE:HB	9:H:28:LEU:HD21	2.01	0.40
11:J:143:ASN:OD1	11:J:143:ASN:C	2.64	0.40
12:K:2:LEU:HD21	12:K:28:HIS:CE1	2.56	0.40
15:N:132:LYS:HB3	15:N:134:VAL:HG23	2.03	0.40
17:P:80:LEU:HD23	17:P:80:LEU:HA	1.92	0.40
19:R:78:ARG:HE	19:R:78:ARG:HB2	1.64	0.40
22:U:30:LYS:HD3	22:U:30:LYS:HA	1.82	0.40
28:a:3:LYS:HE3	28:a:8:ASN:OD1	2.22	0.40
28:a:78:VAL:HG22	28:a:83:VAL:HB	2.01	0.40
34:g:38:LYS:HB2	34:g:60:ARG:HH21	1.86	0.40
34:g:41:ILE:HD13	34:g:41:ILE:HA	1.86	0.40
34:g:79:LEU:HB3	34:g:111:VAL:HG11	2.04	0.40
1:2:1339:U:O2'	1:2:1340:U:OP1	2.35	0.40
1:2:1455:A:C2	1:2:1456:G:C8	3.09	0.40
1:2:1475:G:H8	1:2:1475:G:OP2	2.04	0.40
1:2:1544:C:H5''	21:T:78:ILE:O	2.20	0.40
1:2:1627:C:H4'	21:T:41:LYS:HB3	2.02	0.40
3:B:83:LYS:HE3	3:B:83:LYS:HB2	1.85	0.40
4:C:246:LYS:HE2	4:C:246:LYS:HB2	1.94	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:E:87:MET:HE1	6:E:236:ILE:HD13	2.03	0.40
7:F:82:ASN:ND2	7:F:88:MET:HG3	2.36	0.40
8:G:147:LEU:HD11	8:G:156:TYR:CE2	2.57	0.40
18:Q:102:GLU:O	18:Q:106:LYS:HB2	2.22	0.40
21:T:63:HIS:CG	21:T:75:MET:HB2	2.56	0.40
30:c:13:ARG:O	30:c:32:VAL:HA	2.21	0.40
34:g:157:SER:HA	34:g:158:PRO:HD3	1.93	0.40
1:2:34:U:H2'	1:2:35:C:O4'	2.21	0.40
1:2:414:A:H3'	1:2:415:A:C8	2.56	0.40
1:2:618:C:O4'	1:2:632:C:H4'	2.21	0.40
1:2:986:G:C8	16:O:137:SER:O	2.75	0.40
1:2:1253:A:C8	1:2:1666:C:H4'	2.56	0.40
1:2:1413:G:H22	1:2:1427:C:H2'	1.87	0.40
1:2:1535:U:C5	7:F:88:MET:HE3	2.56	0.40
1:2:1535:U:O4	7:F:169:ILE:HB	2.21	0.40
2:A:206:ASP:O	2:A:210:ILE:HG13	2.21	0.40
3:B:187:LYS:HG2	3:B:192:SER:HB3	2.03	0.40
14:M:122:ASP:O	14:M:126:GLU:HG2	2.21	0.40
15:N:4:MET:HE3	15:N:4:MET:HB2	1.91	0.40
17:P:120:SER:N	20:S:120:HIS:HD2	2.20	0.40
29:b:35:VAL:HA	29:b:78:SER:O	2.21	0.40
34:g:33:SER:OG	34:g:41:ILE:O	2.29	0.40
1:2:1182:A:OP1	35:h:18:ARG:NH1	2.54	0.40
2:A:149:ASN:HB2	2:A:165:ASN:OD1	2.21	0.40
5:D:117:ARG:HA	5:D:117:ARG:HD2	1.70	0.40
7:F:71:ARG:HD2	7:F:75:SER:OG	2.21	0.40
11:J:107:GLU:OE1	11:J:107:GLU:HA	2.22	0.40
12:K:3:MET:CE	12:K:48:ALA:HA	2.51	0.40
16:O:119:LEU:O	16:O:124:MET:HB2	2.22	0.40
24:W:83:LEU:HD12	24:W:83:LEU:HA	1.92	0.40
34:g:78:ALA:HB3	34:g:90:TRP:HB3	2.04	0.40
34:g:90:TRP:CH2	34:g:96:THR:HA	2.56	0.40
1:2:1010:G:H2'	1:2:1011:A:H8	1.86	0.40
1:2:1093:A:H2'	1:2:1094:C:H6	1.87	0.40
1:2:1538:C:C4'	21:T:44:GLU:H	2.34	0.40
3:B:67:PHE:CE2	16:O:48:SER:HB3	2.57	0.40
4:C:176:LYS:HD2	4:C:176:LYS:HA	1.70	0.40
6:E:136:ILE:HG23	6:E:149:TYR:CE1	2.57	0.40
8:G:22:ARG:HA	8:G:25:ARG:HG2	2.04	0.40
8:G:162:LEU:O	8:G:169:PRO:HA	2.22	0.40
10:I:133:GLU:O	10:I:136:ILE:HG22	2.22	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:K:5:LYS:O	12:K:9:ILE:HG12	2.21	0.40
17:P:32:GLN:HE21	17:P:32:GLN:HB2	1.64	0.40
18:Q:43:GLU:O	18:Q:45:ARG:NH1	2.54	0.40
19:R:54:VAL:O	19:R:58:MET:HG3	2.21	0.40
26:Y:47:MET:HE3	26:Y:47:MET:HB2	1.78	0.40
27:Z:96:LEU:HD12	27:Z:96:LEU:HA	1.87	0.40
27:Z:107:VAL:HB	27:Z:109:TYR:CZ	2.57	0.40
31:d:17:GLY:O	31:d:27:ARG:NH2	2.54	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	A	211/295 (72%)	205 (97%)	5 (2%)	1 (0%)	24	53
3	B	211/264 (80%)	206 (98%)	5 (2%)	0	100	100
4	C	216/293 (74%)	211 (98%)	5 (2%)	0	100	100
5	D	223/243 (92%)	216 (97%)	7 (3%)	0	100	100
6	E	260/263 (99%)	251 (96%)	9 (4%)	0	100	100
7	F	187/204 (92%)	157 (84%)	29 (16%)	1 (0%)	24	53
8	G	228/249 (92%)	220 (96%)	7 (3%)	1 (0%)	30	58
9	H	184/194 (95%)	169 (92%)	12 (6%)	3 (2%)	7	24
10	I	203/208 (98%)	195 (96%)	8 (4%)	0	100	100
11	J	178/194 (92%)	172 (97%)	6 (3%)	0	100	100
12	K	95/165 (58%)	88 (93%)	7 (7%)	0	100	100
13	L	149/158 (94%)	144 (97%)	5 (3%)	0	100	100
14	M	119/132 (90%)	108 (91%)	9 (8%)	2 (2%)	7	23

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
15	N	147/151 (97%)	145 (99%)	2 (1%)	0	100	100
16	O	133/151 (88%)	128 (96%)	5 (4%)	0	100	100
17	P	124/145 (86%)	109 (88%)	13 (10%)	2 (2%)	7	24
18	Q	136/146 (93%)	122 (90%)	13 (10%)	1 (1%)	18	45
19	R	130/135 (96%)	127 (98%)	3 (2%)	0	100	100
20	S	141/152 (93%)	124 (88%)	16 (11%)	1 (1%)	18	45
21	T	142/145 (98%)	130 (92%)	12 (8%)	0	100	100
22	U	99/119 (83%)	91 (92%)	8 (8%)	0	100	100
23	V	80/83 (96%)	80 (100%)	0	0	100	100
24	W	127/130 (98%)	124 (98%)	3 (2%)	0	100	100
25	X	139/143 (97%)	138 (99%)	1 (1%)	0	100	100
26	Y	122/130 (94%)	120 (98%)	2 (2%)	0	100	100
27	Z	70/125 (56%)	63 (90%)	7 (10%)	0	100	100
28	a	97/101 (96%)	95 (98%)	1 (1%)	1 (1%)	12	36
29	b	80/82 (98%)	77 (96%)	3 (4%)	0	100	100
30	c	60/62 (97%)	53 (88%)	7 (12%)	0	100	100
31	d	53/55 (96%)	51 (96%)	2 (4%)	0	100	100
32	e	54/56 (96%)	54 (100%)	0	0	100	100
33	f	72/74 (97%)	64 (89%)	8 (11%)	0	100	100
34	g	312/315 (99%)	283 (91%)	28 (9%)	1 (0%)	36	64
35	h	20/25 (80%)	20 (100%)	0	0	100	100
36	n	33/180 (18%)	33 (100%)	0	0	100	100
All	All	4835/5567 (87%)	4573 (95%)	248 (5%)	14 (0%)	37	64

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	189	ILE
7	F	79	HIS
9	H	43	LEU
14	M	109	VAL
14	M	110	VAL
18	Q	17	LYS
28	a	63	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
34	g	314	ILE
9	H	30	LEU
9	H	29	GLU
17	P	92	SER
20	S	100	ALA
8	G	151	ASP
17	P	87	PRO

5.3.2 Protein sidechains [i](#)

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	A	179/243 (74%)	171 (96%)	8 (4%)	24	57
3	B	194/231 (84%)	183 (94%)	11 (6%)	18	47
4	C	184/225 (82%)	181 (98%)	3 (2%)	55	82
5	D	189/202 (94%)	177 (94%)	12 (6%)	16	43
6	E	224/225 (100%)	220 (98%)	4 (2%)	51	80
7	F	159/170 (94%)	149 (94%)	10 (6%)	16	43
8	G	200/218 (92%)	191 (96%)	9 (4%)	24	57
9	H	167/174 (96%)	155 (93%)	12 (7%)	13	37
10	I	178/180 (99%)	172 (97%)	6 (3%)	32	66
11	J	160/168 (95%)	157 (98%)	3 (2%)	50	79
12	K	88/136 (65%)	85 (97%)	3 (3%)	32	66
13	L	135/142 (95%)	131 (97%)	4 (3%)	36	69
14	M	102/108 (94%)	102 (100%)	0	100	100
15	N	130/131 (99%)	126 (97%)	4 (3%)	35	69
16	O	105/119 (88%)	100 (95%)	5 (5%)	23	54
17	P	112/130 (86%)	97 (87%)	15 (13%)	4	13
18	Q	114/121 (94%)	109 (96%)	5 (4%)	25	58
19	R	119/122 (98%)	111 (93%)	8 (7%)	15	40

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
20	S	124/132 (94%)	113 (91%)	11 (9%)	9	27
21	T	114/115 (99%)	108 (95%)	6 (5%)	20	50
22	U	93/107 (87%)	89 (96%)	4 (4%)	26	58
23	V	66/67 (98%)	63 (96%)	3 (4%)	24	57
24	W	112/113 (99%)	106 (95%)	6 (5%)	20	50
25	X	113/115 (98%)	111 (98%)	2 (2%)	51	80
26	Y	108/112 (96%)	104 (96%)	4 (4%)	30	63
27	Z	64/103 (62%)	60 (94%)	4 (6%)	16	43
28	a	87/89 (98%)	87 (100%)	0	100	100
29	b	74/74 (100%)	74 (100%)	0	100	100
30	c	55/55 (100%)	52 (94%)	3 (6%)	19	49
31	d	48/48 (100%)	45 (94%)	3 (6%)	16	43
32	e	45/45 (100%)	44 (98%)	1 (2%)	45	77
33	f	67/67 (100%)	65 (97%)	2 (3%)	36	69
34	g	272/273 (100%)	254 (93%)	18 (7%)	15	41
35	h	21/24 (88%)	21 (100%)	0	100	100
36	n	30/147 (20%)	29 (97%)	1 (3%)	33	67
All	All	4232/4731 (90%)	4042 (96%)	190 (4%)	26	57

All (190) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	28	THR
2	A	75	SER
2	A	87	VAL
2	A	104	THR
2	A	111	GLN
2	A	136	GLU
2	A	152	SER
2	A	154	LEU
3	B	50	THR
3	B	52	THR
3	B	62	LEU
3	B	69	VAL
3	B	76	ASN
3	B	77	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
3	B	108	ASP
3	B	125	VAL
3	B	131	ASP
3	B	143	THR
3	B	179	ASN
4	C	163	VAL
4	C	259	THR
4	C	267	GLN
5	D	9	ARG
5	D	38	GLU
5	D	55	THR
5	D	139	SER
5	D	143	ARG
5	D	149	SER
5	D	156	LEU
5	D	157	MET
5	D	184	ILE
5	D	190	LEU
5	D	193	ASP
5	D	209	SER
6	E	12	VAL
6	E	23	LEU
6	E	173	ILE
6	E	204	SER
7	F	33	ILE
7	F	47	LYS
7	F	49	LEU
7	F	72	LEU
7	F	89	THR
7	F	101	HIS
7	F	102	LEU
7	F	116	ILE
7	F	126	THR
7	F	169	ILE
8	G	20	ASP
8	G	26	THR
8	G	34	THR
8	G	125	THR
8	G	143	LYS
8	G	144	LEU
8	G	153	VAL
8	G	163	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
8	G	204	GLU
9	H	9	VAL
9	H	28	LEU
9	H	30	LEU
9	H	32	MET
9	H	33	ASN
9	H	36	LEU
9	H	37	LYS
9	H	41	ARG
9	H	42	GLU
9	H	60	ILE
9	H	61	ILE
9	H	172	THR
10	I	81	VAL
10	I	88	ASN
10	I	95	THR
10	I	130	THR
10	I	159	SER
10	I	160	SER
11	J	128	VAL
11	J	153	SER
11	J	159	PHE
12	K	22	VAL
12	K	26	ASP
12	K	72	THR
13	L	12	LYS
13	L	15	THR
13	L	97	ARG
13	L	146	THR
15	N	9	LYS
15	N	53	ILE
15	N	57	SER
15	N	143	SER
16	O	46	ASP
16	O	57	THR
16	O	75	MET
16	O	81	VAL
16	O	133	THR
17	P	32	GLN
17	P	38	SER
17	P	40	ARG
17	P	45	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
17	P	49	LEU
17	P	52	LYS
17	P	79	HIS
17	P	83	MET
17	P	85	ILE
17	P	86	LEU
17	P	88	GLU
17	P	89	MET
17	P	90	VAL
17	P	102	PHE
17	P	103	ASN
18	Q	24	HIS
18	Q	42	ILE
18	Q	93	VAL
18	Q	116	ASP
18	Q	119	LEU
19	R	6	THR
19	R	9	VAL
19	R	74	GLN
19	R	91	LEU
19	R	98	VAL
19	R	115	SER
19	R	124	VAL
19	R	127	ASN
20	S	12	ILE
20	S	16	LEU
20	S	46	ARG
20	S	84	LEU
20	S	86	ARG
20	S	87	GLN
20	S	111	LEU
20	S	121	ARG
20	S	131	VAL
20	S	138	THR
20	S	145	THR
21	T	9	VAL
21	T	34	VAL
21	T	51	ASN
21	T	82	ARG
21	T	90	SER
21	T	107	LEU
22	U	24	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
22	U	29	VAL
22	U	63	ILE
22	U	79	ARG
23	V	39	VAL
23	V	52	THR
23	V	59	ILE
24	W	30	CYS
24	W	47	ILE
24	W	74	VAL
24	W	76	SER
24	W	80	ASP
24	W	121	THR
25	X	53	GLU
25	X	72	VAL
26	Y	24	VAL
26	Y	35	VAL
26	Y	39	GLU
26	Y	104	ARG
27	Z	48	VAL
27	Z	72	VAL
27	Z	92	LEU
27	Z	97	ILE
30	c	18	LEU
30	c	45	ASN
30	c	46	VAL
31	d	10	HIS
31	d	14	PHE
31	d	30	LEU
32	e	29	THR
33	f	115	SER
33	f	136	PHE
34	g	10	THR
34	g	31	ILE
34	g	94	THR
34	g	102	VAL
34	g	109	LEU
34	g	111	VAL
34	g	137	VAL
34	g	194	TYR
34	g	199	THR
34	g	200	VAL
34	g	217	MET

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
34	g	221	LEU
34	g	241	PHE
34	g	260	ASP
34	g	270	LEU
34	g	285	GLN
34	g	291	TRP
34	g	309	VAL
36	n	147	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	24	HIS
2	A	111	GLN
3	B	75	GLN
3	B	186	ASN
4	C	113	GLN
4	C	172	ASN
5	D	57	ASN
5	D	101	GLN
6	E	17	HIS
6	E	67	GLN
6	E	98	ASN
6	E	142	HIS
6	E	161	GLN
6	E	188	ASN
7	F	82	ASN
8	G	13	GLN
9	H	33	ASN
10	I	9	HIS
10	I	111	GLN
13	L	94	HIS
15	N	13	GLN
17	P	32	GLN
17	P	35	GLN
17	P	104	GLN
21	T	126	GLN
22	U	100	GLN
24	W	82	GLN
25	X	26	GLN
26	Y	22	GLN
27	Z	64	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
31	d	4	GLN
33	f	135	HIS
34	g	14	HIS
34	g	51	ASN
34	g	119	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	2	1661/1869 (88%)	650 (39%)	27 (1%)

All (650) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	2	3	C
1	2	4	C
1	2	17	C
1	2	30	C
1	2	33	G
1	2	36	U
1	2	41	G
1	2	42	A
1	2	46	A
1	2	56	G
1	2	62	G
1	2	66	G
1	2	67	C
1	2	68	A
1	2	70	G
1	2	72	C
1	2	73	C
1	2	74	G
1	2	75	G
1	2	76	U
1	2	77	A
1	2	79	A
1	2	102	A
1	2	103	A
1	2	110	U
1	2	113	G
1	2	115	U

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	126	G
1	2	127	C
1	2	128	U
1	2	129	C
1	2	130	G
1	2	142	C
1	2	143	U
1	2	145	G
1	2	146	G
1	2	147	A
1	2	155	G
1	2	160	U
1	2	163	U
1	2	168	C
1	2	170	A
1	2	173	A
1	2	176	U
1	2	177	G
1	2	178	C
1	2	179	C
1	2	181	A
1	2	182	C
1	2	183	G
1	2	185	G
1	2	186	C
1	2	187	G
1	2	188	C
1	2	190	G
1	2	191	A
1	2	192	C
1	2	193	C
1	2	197	U
1	2	198	U
1	2	199	C
1	2	200	G
1	2	204	G
1	2	205	G
1	2	206	G
1	2	208	G
1	2	209	A
1	2	210	U
1	2	211	G

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	215	G
1	2	216	C
1	2	217	A
1	2	219	U
1	2	220	U
1	2	221	A
1	2	222	U
1	2	224	A
1	2	225	G
1	2	291	G
1	2	293	C
1	2	296	U
1	2	297	A
1	2	299	A
1	2	301	A
1	2	302	A
1	2	304	C
1	2	306	C
1	2	308	G
1	2	309	G
1	2	311	C
1	2	317	C
1	2	318	A
1	2	320	G
1	2	321	C
1	2	322	C
1	2	330	G
1	2	332	G
1	2	335	G
1	2	338	G
1	2	340	C
1	2	341	C
1	2	344	U
1	2	345	U
1	2	346	C
1	2	347	G
1	2	351	G
1	2	352	U
1	2	353	C
1	2	362	C
1	2	363	A
1	2	364	A

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	368	U
1	2	370	G
1	2	377	G
1	2	381	C
1	2	385	G
1	2	386	C
1	2	387	C
1	2	398	A
1	2	400	C
1	2	407	G
1	2	408	A
1	2	409	C
1	2	414	A
1	2	415	A
1	2	417	C
1	2	418	A
1	2	421	G
1	2	448	A
1	2	449	A
1	2	450	C
1	2	464	A
1	2	470	G
1	2	471	G
1	2	472	C
1	2	473	A
1	2	474	G
1	2	476	A
1	2	482	G
1	2	484	A
1	2	485	A
1	2	487	U
1	2	488	U
1	2	492	C
1	2	493	A
1	2	500	A
1	2	501	C
1	2	502	C
1	2	513	G
1	2	514	U
1	2	516	A
1	2	523	A
1	2	528	A

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	529	A
1	2	530	U
1	2	533	A
1	2	534	G
1	2	539	C
1	2	540	U
1	2	541	U
1	2	542	U
1	2	544	G
1	2	545	A
1	2	546	G
1	2	547	G
1	2	548	C
1	2	549	C
1	2	550	C
1	2	553	U
1	2	554	A
1	2	555	A
1	2	556	U
1	2	559	G
1	2	564	A
1	2	566	U
1	2	567	C
1	2	570	C
1	2	571	U
1	2	576	A
1	2	577	U
1	2	578	C
1	2	579	C
1	2	580	U
1	2	588	G
1	2	589	G
1	2	590	A
1	2	591	U
1	2	603	C
1	2	604	A
1	2	605	A
1	2	606	G
1	2	607	U
1	2	608	C
1	2	614	C
1	2	621	C

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	626	G
1	2	627	U
1	2	629	A
1	2	631	U
1	2	639	C
1	2	640	A
1	2	643	A
1	2	644	G
1	2	655	A
1	2	660	C
1	2	666	U
1	2	668	A
1	2	669	A
1	2	671	A
1	2	672	A
1	2	673	G
1	2	685	A
1	2	688	U
1	2	747	U
1	2	748	C
1	2	749	U
1	2	752	G
1	2	793	G
1	2	794	A
1	2	797	C
1	2	799	U
1	2	800	U
1	2	801	U
1	2	802	A
1	2	804	U
1	2	805	U
1	2	808	A
1	2	809	A
1	2	821	G
1	2	822	U
1	2	823	U
1	2	824	C
1	2	832	G
1	2	833	C
1	2	834	C
1	2	847	A
1	2	853	C

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	854	A
1	2	859	G
1	2	860	G
1	2	861	A
1	2	862	A
1	2	863	U
1	2	864	A
1	2	869	A
1	2	870	A
1	2	871	U
1	2	872	A
1	2	873	G
1	2	874	G
1	2	875	A
1	2	878	G
1	2	879	C
1	2	882	U
1	2	883	U
1	2	886	A
1	2	887	U
1	2	888	U
1	2	889	U
1	2	890	U
1	2	893	U
1	2	894	G
1	2	895	G
1	2	898	U
1	2	899	U
1	2	901	G
1	2	902	G
1	2	903	A
1	2	906	U
1	2	907	G
1	2	908	A
1	2	912	C
1	2	913	A
1	2	914	U
1	2	916	A
1	2	920	A
1	2	922	A
1	2	933	G
1	2	940	U

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	941	C
1	2	943	U
1	2	955	A
1	2	961	G
1	2	962	A
1	2	963	A
1	2	964	A
1	2	967	C
1	2	968	U
1	2	969	U
1	2	970	G
1	2	981	A
1	2	988	C
1	2	990	A
1	2	992	A
1	2	999	G
1	2	1002	U
1	2	1008	A
1	2	1017	U
1	2	1022	U
1	2	1023	A
1	2	1026	C
1	2	1027	A
1	2	1028	A
1	2	1045	U
1	2	1049	A
1	2	1060	A
1	2	1061	U
1	2	1062	A
1	2	1078	C
1	2	1083	A
1	2	1085	C
1	2	1096	G
1	2	1100	A
1	2	1109	C
1	2	1114	U
1	2	1115	U
1	2	1117	C
1	2	1118	C
1	2	1119	A
1	2	1120	U
1	2	1123	C

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	1126	G
1	2	1130	G
1	2	1133	A
1	2	1146	C
1	2	1154	U
1	2	1155	U
1	2	1157	G
1	2	1168	G
1	2	1171	G
1	2	1181	A
1	2	1182	A
1	2	1183	A
1	2	1195	A
1	2	1207	G
1	2	1208	A
1	2	1212	G
1	2	1215	C
1	2	1216	C
1	2	1220	A
1	2	1221	G
1	2	1222	G
1	2	1227	G
1	2	1228	A
1	2	1229	G
1	2	1231	C
1	2	1235	G
1	2	1236	G
1	2	1238	U
1	2	1239	U
1	2	1240	A
1	2	1241	A
1	2	1242	U
1	2	1243	U
1	2	1244	U
1	2	1245	G
1	2	1250	A
1	2	1251	A
1	2	1253	A
1	2	1254	C
1	2	1256	G
1	2	1257	G
1	2	1258	A

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	1259	A
1	2	1260	A
1	2	1263	U
1	2	1264	C
1	2	1267	C
1	2	1269	G
1	2	1272	C
1	2	1274	G
1	2	1276	A
1	2	1277	C
1	2	1278	A
1	2	1279	C
1	2	1280	G
1	2	1281	G
1	2	1282	A
1	2	1283	C
1	2	1284	A
1	2	1286	G
1	2	1287	A
1	2	1289	U
1	2	1290	G
1	2	1291	A
1	2	1292	C
1	2	1294	G
1	2	1295	A
1	2	1297	U
1	2	1298	G
1	2	1299	A
1	2	1300	U
1	2	1301	A
1	2	1303	C
1	2	1304	U
1	2	1305	C
1	2	1308	U
1	2	1309	C
1	2	1310	U
1	2	1311	C
1	2	1312	G
1	2	1313	A
1	2	1314	U
1	2	1315	U
1	2	1316	C

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	1317	C
1	2	1319	U
1	2	1320	G
1	2	1322	G
1	2	1323	U
1	2	1324	G
1	2	1325	G
1	2	1331	C
1	2	1340	U
1	2	1341	C
1	2	1343	U
1	2	1348	G
1	2	1352	G
1	2	1353	A
1	2	1371	U
1	2	1372	U
1	2	1375	G
1	2	1378	A
1	2	1382	A
1	2	1393	G
1	2	1394	G
1	2	1396	A
1	2	1397	U
1	2	1398	G
1	2	1403	C
1	2	1405	A
1	2	1408	U
1	2	1412	C
1	2	1413	G
1	2	1414	A
1	2	1416	C
1	2	1425	G
1	2	1426	U
1	2	1427	C
1	2	1428	G
1	2	1429	G
1	2	1430	C
1	2	1431	G
1	2	1432	U
1	2	1439	A
1	2	1440	C
1	2	1441	U

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	1447	G
1	2	1448	A
1	2	1449	G
1	2	1453	C
1	2	1454	A
1	2	1462	U
1	2	1463	U
1	2	1471	C
1	2	1472	C
1	2	1474	A
1	2	1475	G
1	2	1477	U
1	2	1478	U
1	2	1480	A
1	2	1488	C
1	2	1489	A
1	2	1490	G
1	2	1493	C
1	2	1494	U
1	2	1495	G
1	2	1497	G
1	2	1504	U
1	2	1505	U
1	2	1506	A
1	2	1507	G
1	2	1508	A
1	2	1513	C
1	2	1514	G
1	2	1515	G
1	2	1516	G
1	2	1517	G
1	2	1518	C
1	2	1520	G
1	2	1521	C
1	2	1522	A
1	2	1523	C
1	2	1524	G
1	2	1530	U
1	2	1531	A
1	2	1533	A
1	2	1534	C
1	2	1535	U

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	1539	U
1	2	1540	G
1	2	1541	G
1	2	1542	C
1	2	1543	U
1	2	1544	C
1	2	1545	A
1	2	1546	G
1	2	1547	C
1	2	1553	C
1	2	1555	U
1	2	1557	C
1	2	1560	U
1	2	1564	C
1	2	1565	C
1	2	1566	G
1	2	1567	G
1	2	1568	C
1	2	1569	A
1	2	1570	G
1	2	1571	G
1	2	1572	C
1	2	1579	A
1	2	1580	A
1	2	1585	U
1	2	1586	U
1	2	1587	G
1	2	1588	A
1	2	1589	A
1	2	1590	C
1	2	1593	C
1	2	1594	A
1	2	1595	U
1	2	1596	U
1	2	1598	G
1	2	1599	U
1	2	1600	G
1	2	1601	A
1	2	1603	G
1	2	1604	G
1	2	1605	G
1	2	1607	A

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	1608	U
1	2	1609	C
1	2	1611	G
1	2	1612	G
1	2	1613	G
1	2	1614	A
1	2	1615	U
1	2	1616	U
1	2	1618	C
1	2	1619	A
1	2	1620	A
1	2	1621	U
1	2	1622	U
1	2	1623	A
1	2	1624	U
1	2	1627	C
1	2	1630	A
1	2	1631	U
1	2	1633	A
1	2	1637	A
1	2	1638	G
1	2	1639	G
1	2	1641	A
1	2	1642	U
1	2	1643	U
1	2	1647	A
1	2	1654	G
1	2	1661	A
1	2	1662	U
1	2	1663	A
1	2	1665	G
1	2	1669	G
1	2	1671	G
1	2	1677	U
1	2	1695	A
1	2	1698	C
1	2	1699	A
1	2	1707	U
1	2	1708	C
1	2	1709	G
1	2	1710	C
1	2	1713	C

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	1715	A
1	2	1717	C
1	2	1718	G
1	2	1719	A
1	2	1720	U
1	2	1721	U
1	2	1722	G
1	2	1723	G
1	2	1724	A
1	2	1725	U
1	2	1728	U
1	2	1729	U
1	2	1730	U
1	2	1737	G
1	2	1738	C
1	2	1739	C
1	2	1740	C
1	2	1741	U
1	2	1742	C
1	2	1743	G
1	2	1746	U
1	2	1750	C
1	2	1751	C
1	2	1752	C
1	2	1753	C
1	2	1755	C
1	2	1756	C
1	2	1757	G
1	2	1758	G
1	2	1760	G
1	2	1761	U
1	2	1772	C
1	2	1773	C
1	2	1774	C
1	2	1775	U
1	2	1777	G
1	2	1778	C
1	2	1779	G
1	2	1780	G
1	2	1781	A
1	2	1783	C
1	2	1784	G

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	1785	C
1	2	1788	A
1	2	1792	G
1	2	1793	A
1	2	1794	C
1	2	1796	G
1	2	1797	U
1	2	1800	A
1	2	1801	A
1	2	1802	C
1	2	1807	C
1	2	1808	U
1	2	1809	A
1	2	1811	C
1	2	1813	A
1	2	1815	A
1	2	1816	G
1	2	1822	A
1	2	1823	A
1	2	1825	A
1	2	1826	G
1	2	1827	U
1	2	1828	C
1	2	1829	G
1	2	1830	U
1	2	1831	A
1	2	1835	A
1	2	1838	U
1	2	1849	G
1	2	1852	C
1	2	1861	G
1	2	1862	G
1	2	1863	A
1	2	1864	U
1	2	1865	C

All (27) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	2	301	A
1	2	345	U
1	2	414	A

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	2	500	A
1	2	501	C
1	2	549	C
1	2	604	A
1	2	628	A
1	2	792	C
1	2	808	A
1	2	963	A
1	2	980	A
1	2	1022	U
1	2	1339	U
1	2	1342	U
1	2	1471	C
1	2	1487	A
1	2	1517	G
1	2	1566	G
1	2	1571	G
1	2	1588	A
1	2	1589	A
1	2	1592	C
1	2	1619	A
1	2	1642	U
1	2	1712	A
1	2	1755	C

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

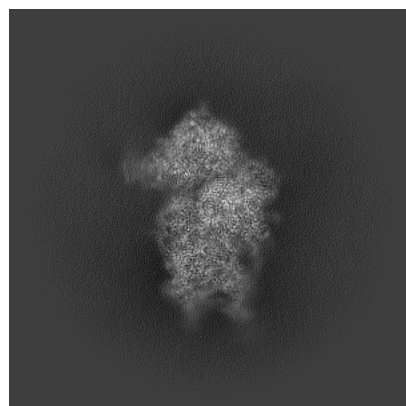
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-62444. These allow visual inspection of the internal detail of the map and identification of artifacts.

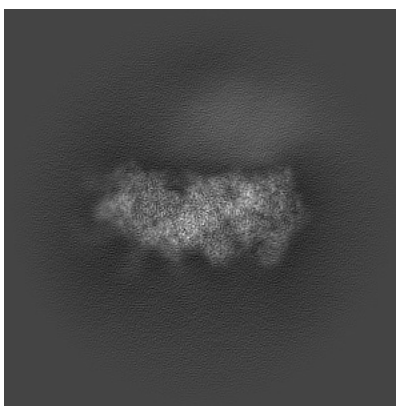
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

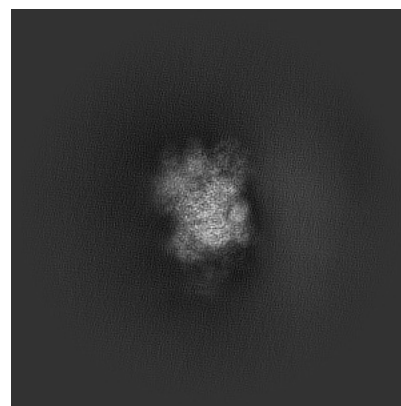
6.1.1 Primary map



X

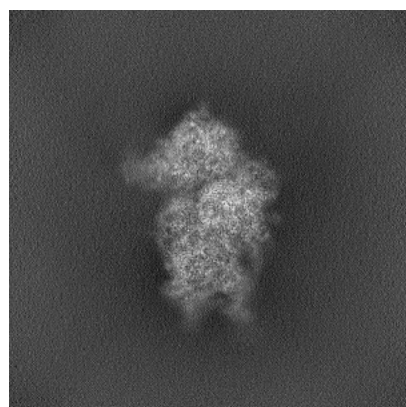


Y

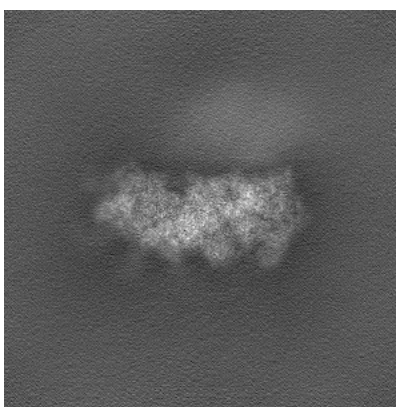


Z

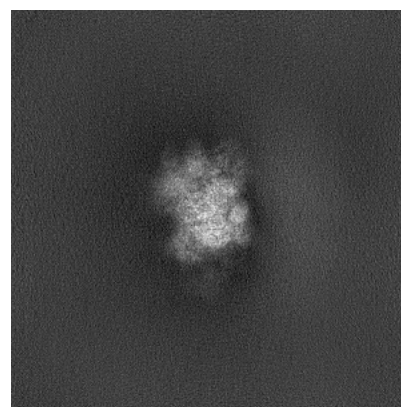
6.1.2 Raw map



X



Y

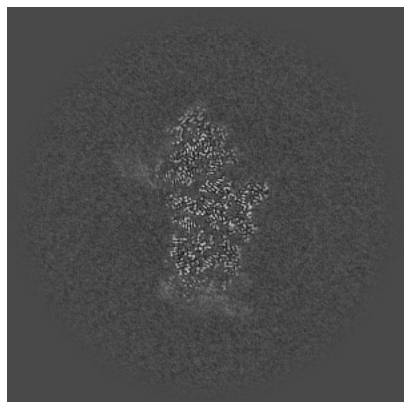


Z

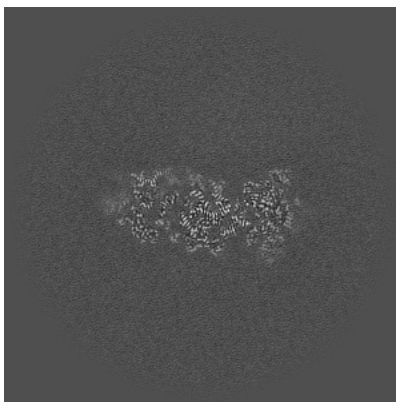
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

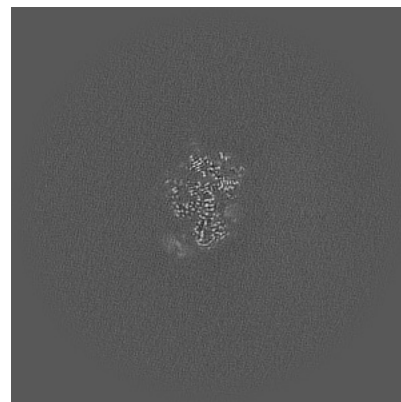
6.2.1 Primary map



X Index: 256

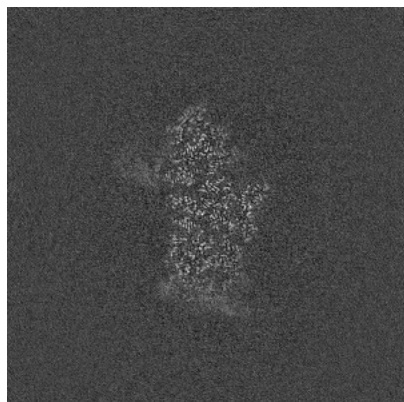


Y Index: 256

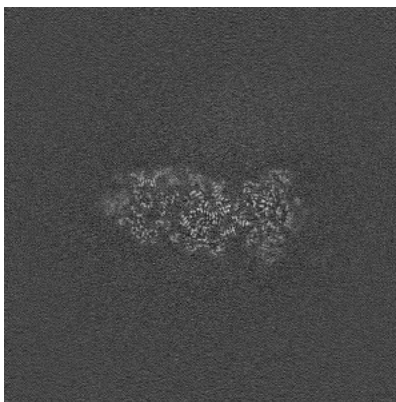


Z Index: 256

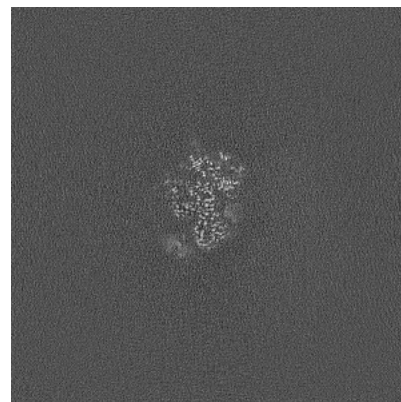
6.2.2 Raw map



X Index: 256



Y Index: 256

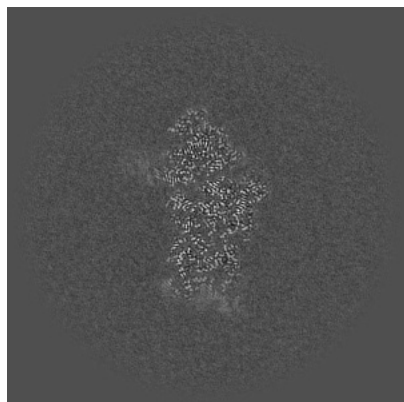


Z Index: 256

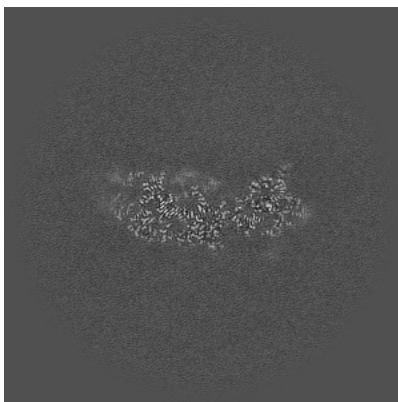
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

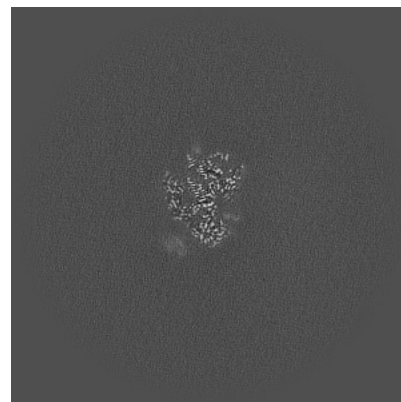
6.3.1 Primary map



X Index: 258

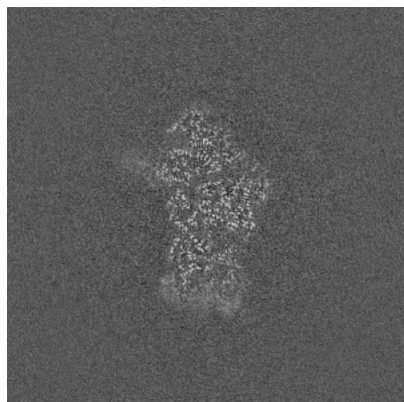


Y Index: 247

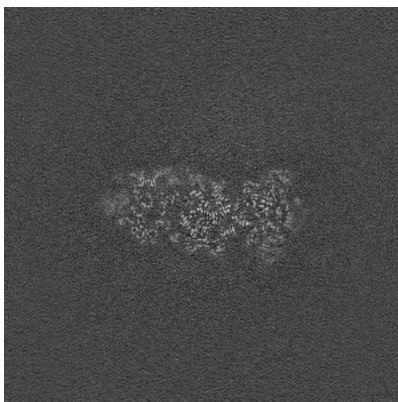


Z Index: 260

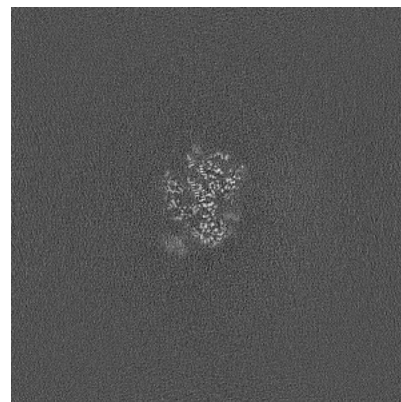
6.3.2 Raw map



X Index: 260



Y Index: 256

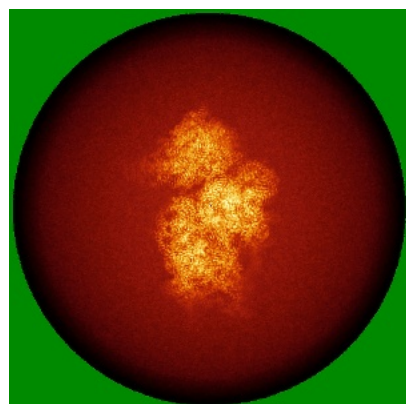


Z Index: 260

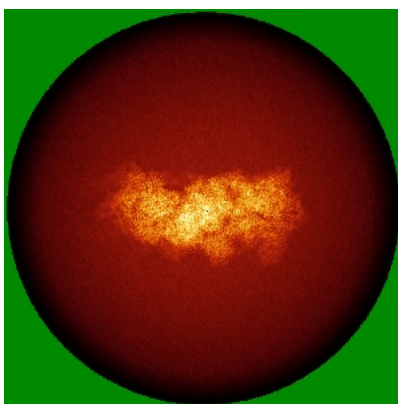
The images above show the largest variance slices of the map in three orthogonal directions.

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

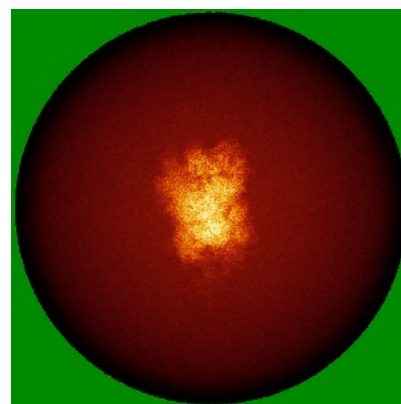
6.4.1 Primary map



X

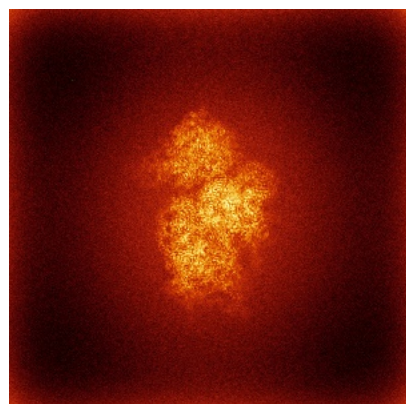


Y

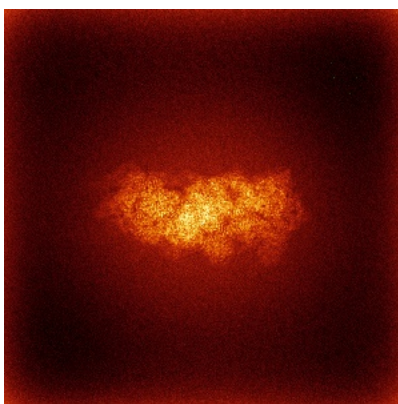


Z

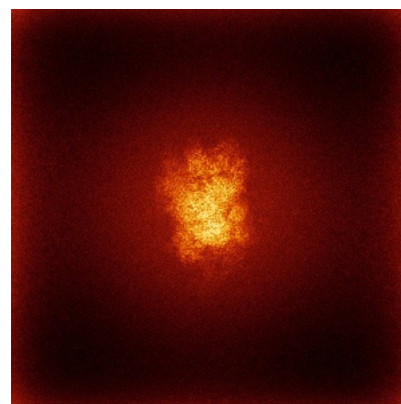
6.4.2 Raw map



X



Y

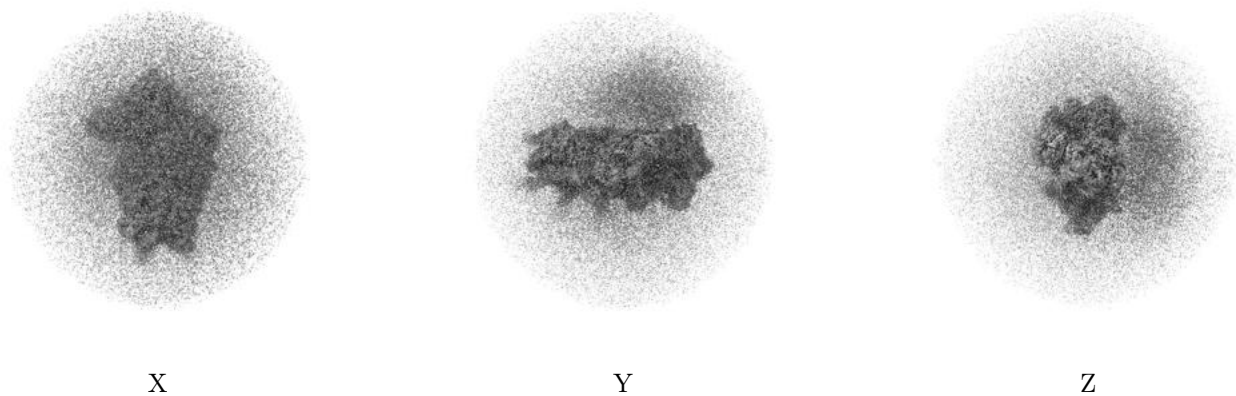


Z

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.

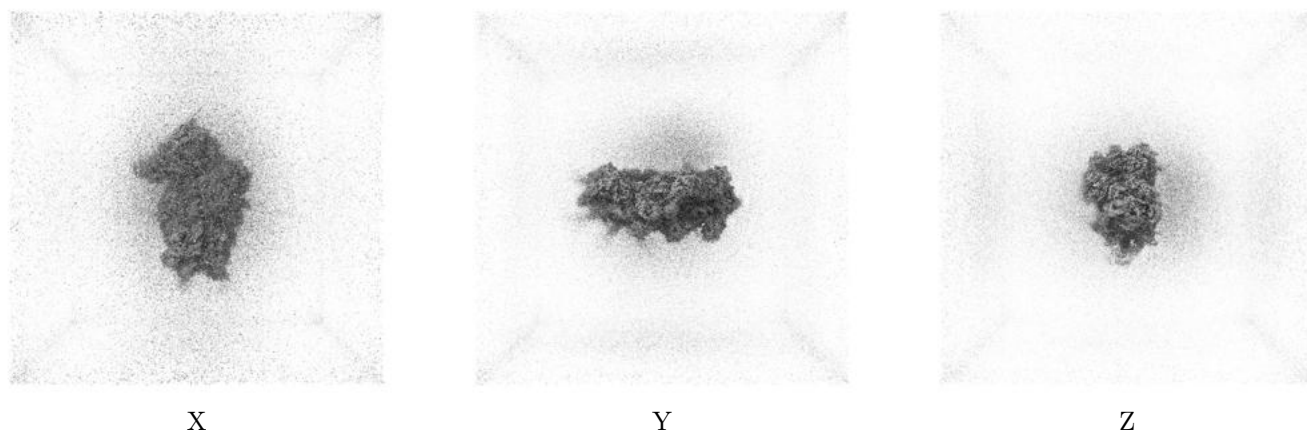
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.177. 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.

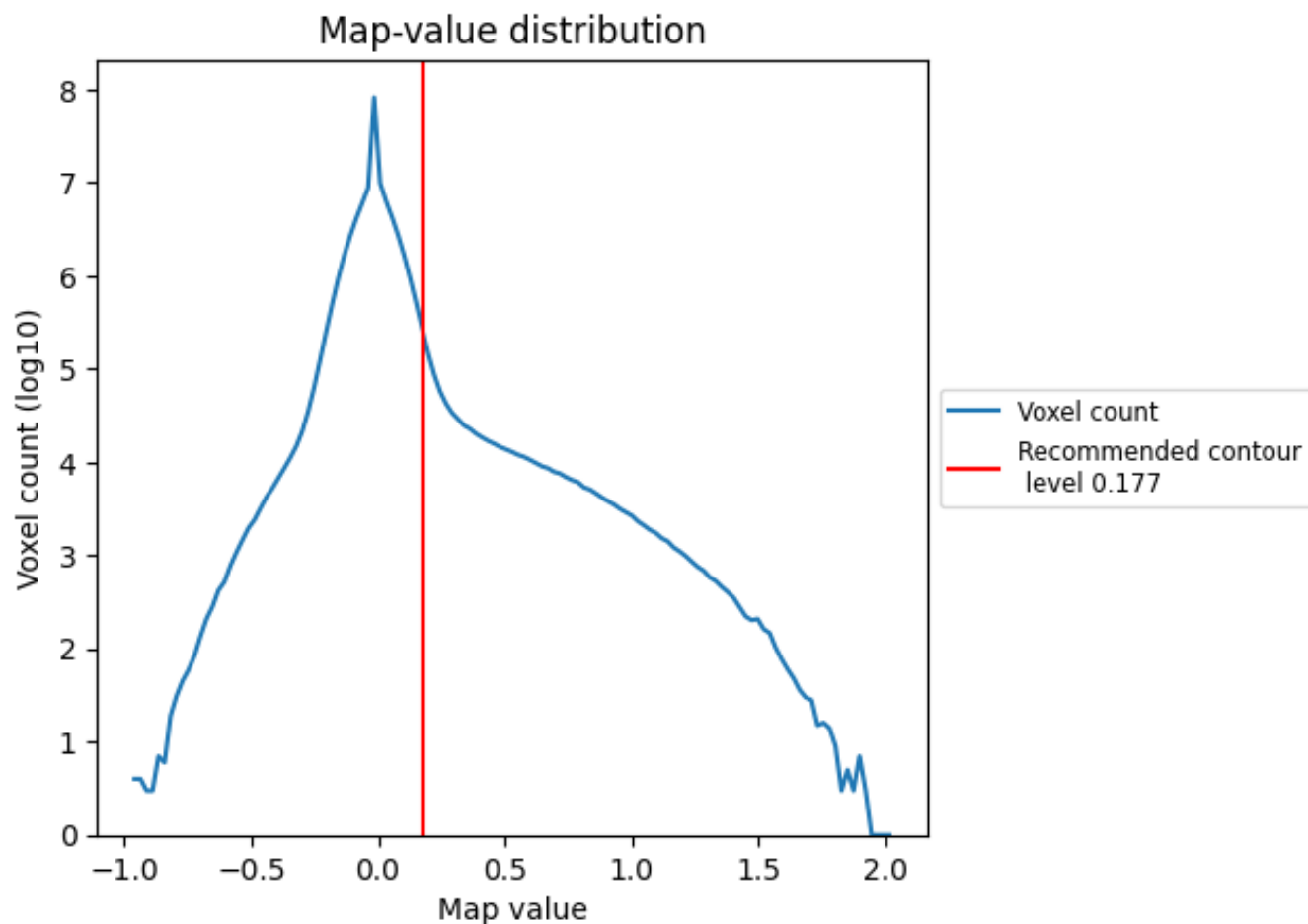
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

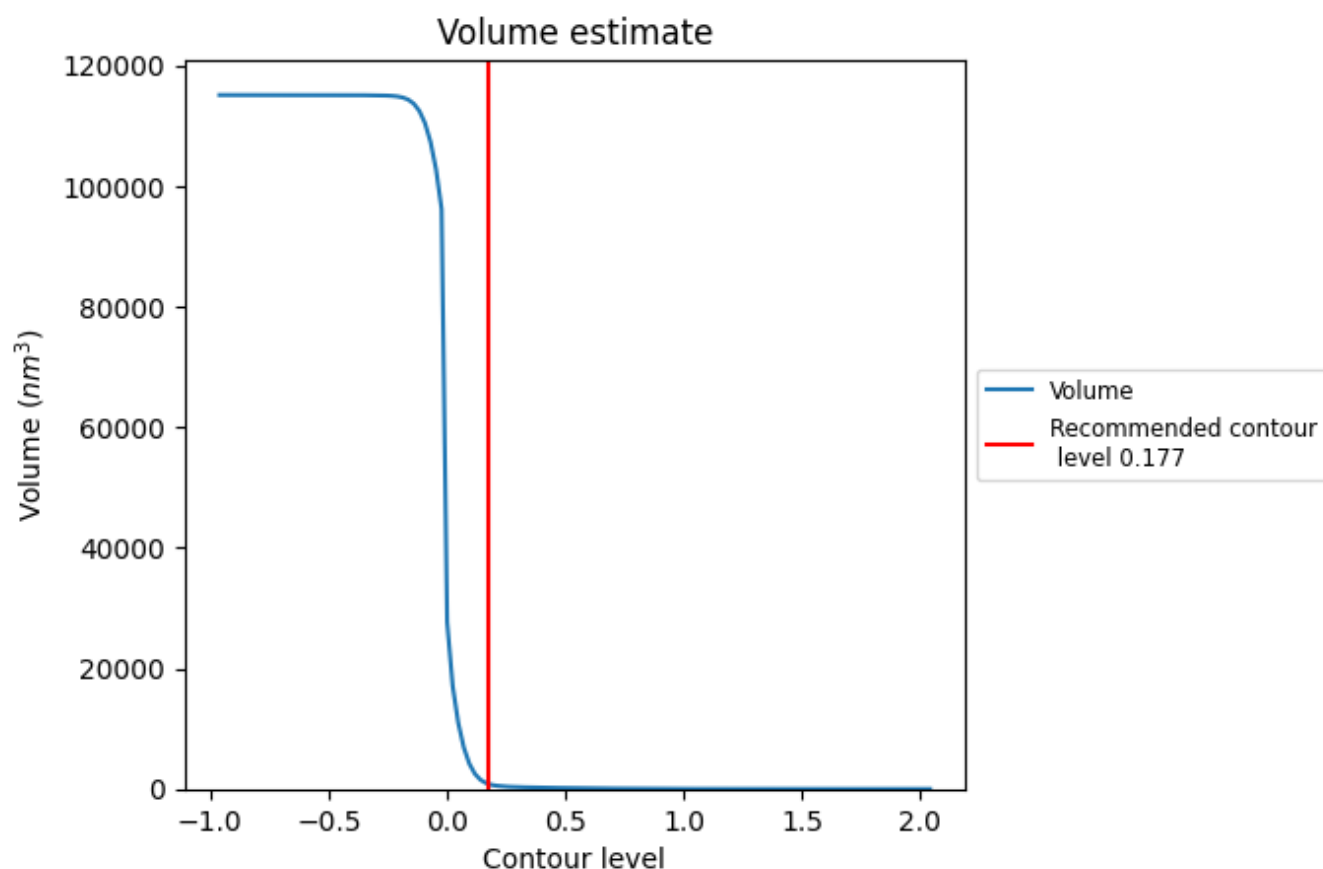
This section contains the results of statistical analysis of the map.

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.

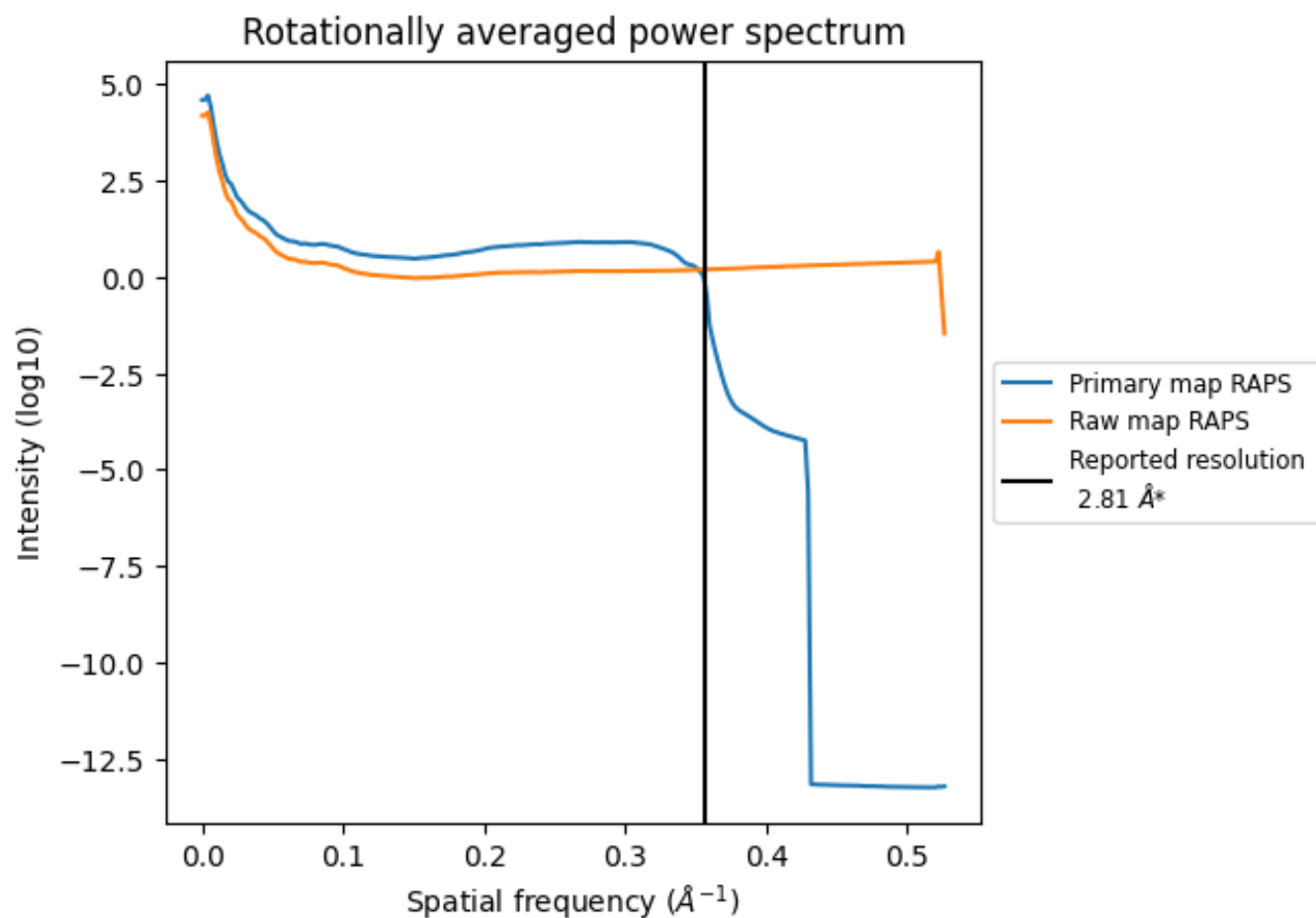
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 826 nm^3 ; this corresponds to an approximate mass of 746 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 ⓘ

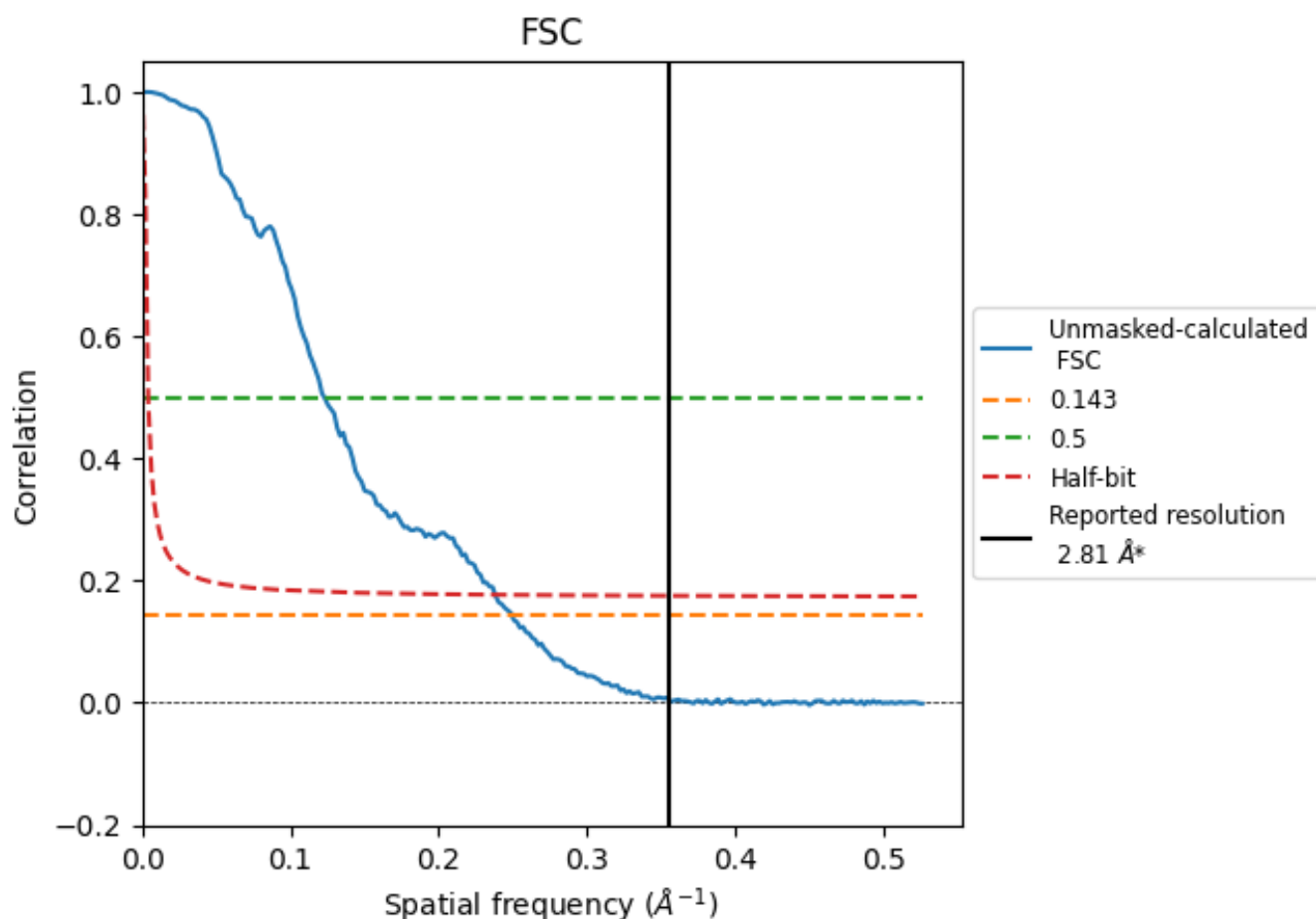


*Reported resolution corresponds to spatial frequency of 0.356 Å⁻¹

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 $0.356 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

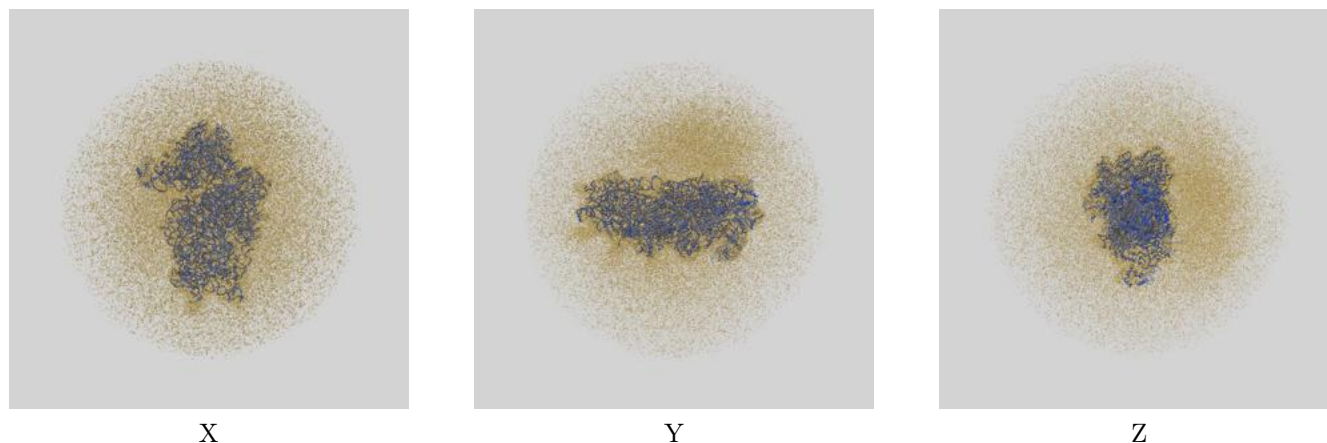
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.81	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.01	8.14	4.20

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.01 differs from the reported value 2.81 by more than 10 %

9 Map-model fit [i](#)

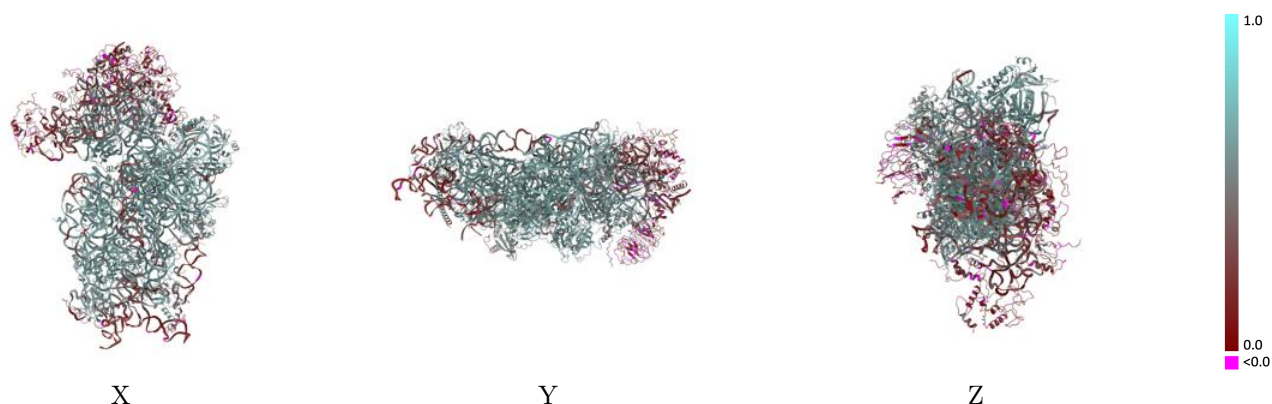
This section contains information regarding the fit between EMDB map EMD-62444 and PDB model 9KMT. Per-residue inclusion information can be found in section [3](#) on page [11](#).

9.1 Map-model overlay [i](#)



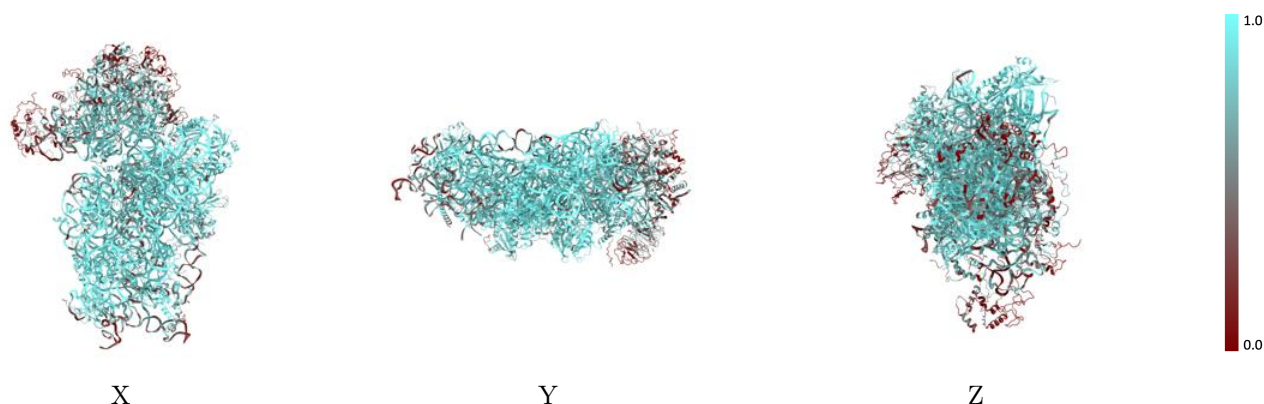
The images above show the 3D surface view of the map at the recommended contour level 0.177 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



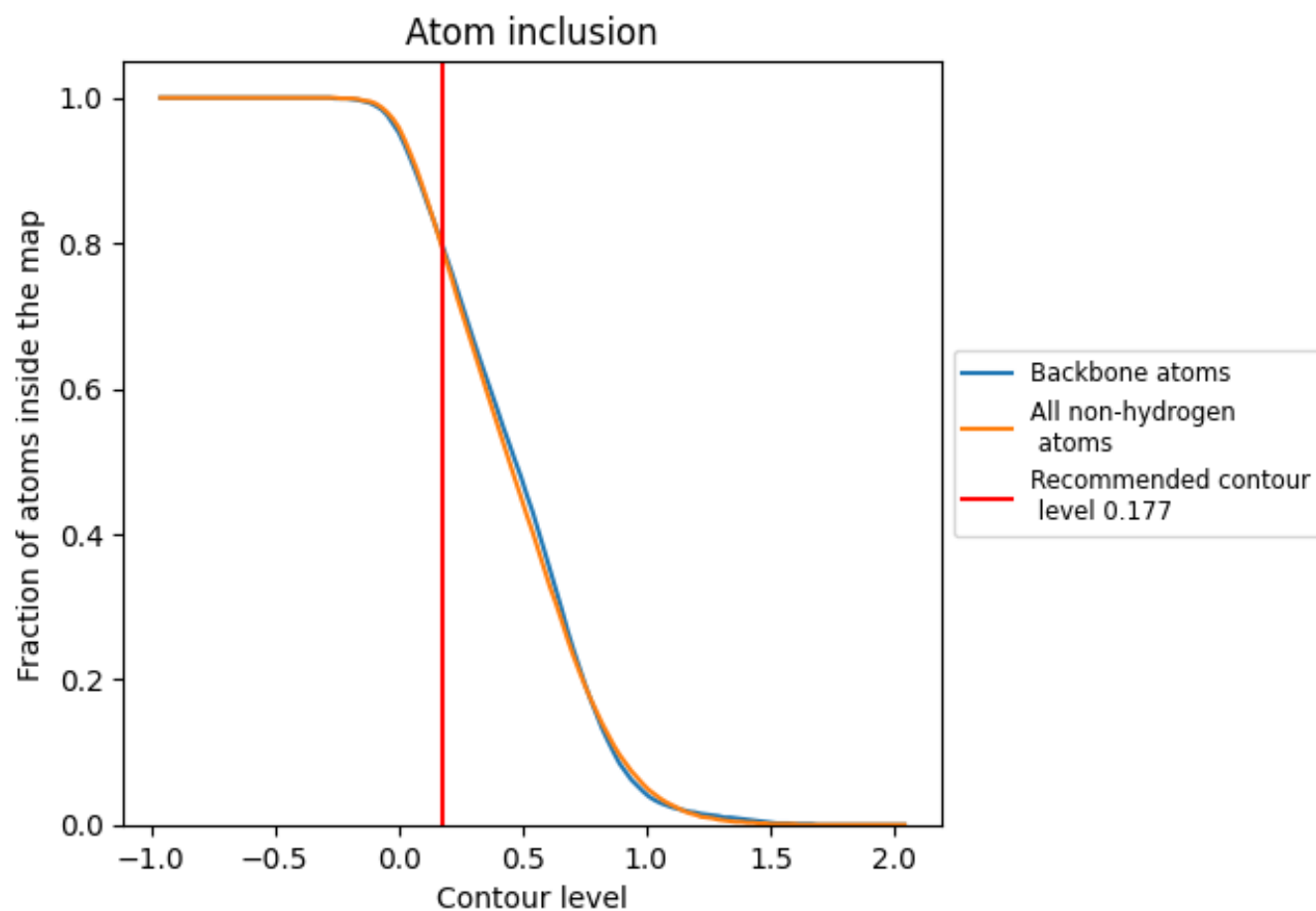
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.177).

9.4 Atom inclusion [i](#)



At the recommended contour level, 80% of all backbone atoms, 79% 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 (0.177) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7920	 0.4870
2	 0.8340	 0.4950
A	 0.9490	 0.6060
B	 0.9370	 0.5910
C	 0.9350	 0.6090
D	 0.8300	 0.5430
E	 0.9480	 0.6070
F	 0.4480	 0.2680
G	 0.8500	 0.5210
H	 0.7890	 0.5140
I	 0.8340	 0.5230
J	 0.9510	 0.6090
K	 0.5630	 0.3090
L	 0.9010	 0.5880
M	 0.2770	 0.2110
N	 0.9370	 0.5960
O	 0.9180	 0.5830
P	 0.3710	 0.2740
Q	 0.6040	 0.3520
R	 0.8630	 0.5570
S	 0.4490	 0.2570
T	 0.5550	 0.2730
U	 0.8070	 0.5130
V	 0.9490	 0.6090
W	 0.9730	 0.6240
X	 0.9510	 0.6060
Y	 0.9390	 0.5990
Z	 0.2100	 0.1950
a	 0.9270	 0.5990
b	 0.8700	 0.5520
c	 0.3690	 0.2530
d	 0.7750	 0.5240
e	 0.8480	 0.5550
f	 0.0990	 0.2010
g	 0.3870	 0.2210



Continued on next page...

Continued from previous page...

Chain	Atom inclusion	Q-score
h	 0.8710	 0.5800
n	 0.9180	 0.6020